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Bio-Aquatic Testing 2501 Mayes Rd

Suite 100 Carrollton, TX 75006 (972) 242-7750

Bioremediation Agent Effectiveness Test

SpillRemed (Marine)

Sarva Bio Remed, LLC

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Prepared by: 9/11/2006 Vice Pres Date

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EXECUTIVE SUMMARY

Bio-Aquatic Testing, Inc. located at 2501 Mayes Rd. Suite 100 Carrollton, Texas 75006 was contracted by Sarva Bio Remed, LLC to test effectiveness of their bioremediation product, SpillRemed (Marine), using Environmental Protection Agency (EPA) protocol listed in 40 CFR Chapter 1 (7-1-99) Pt. 300 Appendix C, Item 4.0. The test protocol calls for application of products onto ANS 521 oil. The product was applied to test flasks according to manufacturer's specifications. Samples were sacrificed on Day 0, Day 7, and Day 28 of the test period. Day 0 and Day 7 samples were sampled for microbiological analysis and then frozen at -10° C until GC/MS results were known for the Day 28 samples. Each replicate of product and control were tested for continued microbiological viability over time, reduction in weight via gravimetric analysis, and reduction in alkane and/or aromatic constituents via Gas Chromatography/Mass Spectroscopy (GC/MS). The product was deemed effective if the data showed the GC/MS product results for Day 28 treatments to be statistically less than Day 0 treatments.

GC/MS data for Days 0, 7, and 28, were consolidated and analyzed with Minitab Statistical program 13.3. Data was analyzed for a significant difference between controls and treatments (products) using a General Linear ANOVA Model with Dunnett's and/or Tukey's means comparison test. GC/MS analysis showed significant reduction of both alkane and aromatic constituents of the test oil as indicated by the statistically significant difference between the Day 28 controls and Day 28 treatments as well as between Day 0 and Day 28 treatments. Day 7 results also showed a statistically significant reduction of treatments as compared to controls.

The surrogate compounds, d-10 phenanthrene and $5-\alpha$ and rostane showed excellent recovery indicating the test to be valid.

Microbiological results showed continued viability of the oil-eating microorganisms over time. Day 7 and Day 28 gravimetric analysis showed a statistically significant reduction from the controls to the treatments.

Based on the parameters of this test, the product should be included on the NCP list of approved bio-remediation products.

BIOREMEDIATION AGENT EFFECTIVENESS TEST USING SARVA BIO REMED, LLC PRODUCT "SPILLREMED (MARINE)"

Introduction

The bioremediation agent effectiveness testing protocol is designed to determine a product's ability to biodegrade oil by quantifying changes in the oil composition resulting from biodegradation. The protocol quantifies the disappearance of saturated hydrocarbons and polynuclear aromatic hydrocarbons (PAHs) as well as weight loss. The protocol also tests for microbial activity over time to ascertain continued viability of oil degrading microorganisms.

Summary of Method

The protocol calls for gas chromatography/mass spectrophotometry and gravimetric analyses to quantify saturated hydrocarbons and PAHs, and determine weight loss respectively. The sample preparation procedure extracts the oil phase into dicloromethane (DCM), with a subsequent distillation to 1-3-mL using a K-D apparatus and Snyder column. To effectively accomplish the goals of the testing protocol, it is necessary to normalize the concentration of the various analytes in oil to a non-biodegradable marker, either C_2 - or C_3 – phenanthrene, C_2 –chrysene, or hopane. The test method targets the relatively easy to degrade normal alkanes and the more resistant and toxic PAHs. It normalizes their concentrations to C2 or C3 phenanthrene, C2 -chrysene, or $C_{30}17\alpha(H)$, 21 β (H)-hopane on an oil weight basis (mg marker/kg oil, mg target analyte/kg). The analytical technique uses a high-resolution gas chromatography/mass spectrophotometer (GC/MS) because of its high degree of chemical separation and spectral resolution. GC/MS has long been used to study the weathering and fate of oil spilled into the environment. For quantitative analyses, the instrument is operated in the selective ion detection mode (SIM) at a scan rate of greater than 1.5 scans per second to maximize the linear quantitative range and precision of the instrument. The sample preparation method does not exclude analysis of selected samples by GC/MS in the full scanning mode of operation to qualitatively assess changes in the oil not accounted for by the SIM approach. Gravimetric analysis is used to support the GC/MS analysis by measuring weight loss of samples over time as compared to controls by drying the extracted samples using nitrogen a blowdown technique.

Performed concurrently with the chemical analysis described above is a microbiological analysis. The microbiological analysis is performed to determine and monitor the viability of relative concentrations of the microbial cultures being studied. Using this method, continued viability is measured over time by comparing serial dilutions of microorganisms, to determine statistical significance between treatments and controls.

MATERIALS AND METHODS

The following methods* were obtained from 40 CFR Chapter 1 (7-1-99) Pt. 300 Appendix C, item 4.0 Bioremediation Effectiveness test, as submitted by the Environmental Protection Agency. Some modifications were made to these methods as discussed below.

The procedure consists of an experimental orbital shaker flask setup using 250-mL Erlenmeyer flasks labeled with unique identifiers using the following treatment design:

*Details from these methods can be found in the aforementioned 40 CFR Chapter 1 (7-1-99) Pt. 300 Appendix C, item 4.0. A copy is available upon request.

Treatment	Number of	samples at san	npling times	Total number of analytical determinations		
	Day 0 Day 7 Day 28		ANALYSES			
				Microbial counts	Gravimetric	GC/MS
Control	3	3	3	9	9	9
Nutrient	3	3	3	9	9	9
SpillRemed	3	3	3	9	9	9

Number of replicates per treatment or control per sampling event - 3 Number of replicates per treatment or control - 9 Total replicates - 27 Control - Oil + Seawater Nutrient - Oil + Seawater + EPA Nutrient SpilRemed - Oil + Seawater + Product + EPA Nutrient

Using sterile technique, each appropriately labeled replicate flask has 100-mL of seawater added. The seawater obtained was from the Gulf of Mexico by faculty at LSU. Each flask is placed on a balance and the weight recorded. Approximately one half-gram (0.5 g) of artificially weathered oil (Alaska North Slope 521)* is then added to each flask while still on the balance and the weight recorded again.

*The ANS 521 oil was obtained from John Haines of the Environmental Protection Agency's Environmental Monitoring and Support Laboratory, Cincinnati, Ohio, 45268

The control flasks were prepared by adding oil to the natural seawater.

The nutrient flasks were prepared as instructed in 40 CFR Chapter 1 (7-1-99) Pt. 300 Appendix C.

The product mix was prepared according to the manufacturer's instructions. The product was applied to each oil + product flask at a ratio of 10:1 (V/V).

After preparing all treatments and controls, three replicates of each treatment and control were shaken on an orbital shaker at 190 - 200 rpm and incubated at 20° C until sacrificed for the Day 0, 7, and 28 analyses. At each sampling (sacrifice) day, a 0.5-mL aliquot was set aside for microbiological analysis and the remaining solution is prepared for chemical analysis.

A phosphate buffer solution was made from a recipe obtained from Jan Kurtz of the Microbial Ecology Branch of the Environmental Protection Agency's Gulf Breeze Ecology Division. This recipe is provided in Appendix 1. A 0.5-mL aliquot from each replicate was added to a test tube containing 4.5-mL of a sterile phosphate buffer for the microbiological analysis. Aseptic technique was then used to make serial dilutions down to a 10^{-8} dilution. Microtiter plates were prepared by adding 1.75-mL of Bushnell-Haas broth into to each well. Six replicates per dilution are used per treatment or control giving a total of forty-eight wells, (48) per treatment or control. Each of the wells was inoculated with 0.1-mL of solution from each of the serial dilutions made from the original aliquot of 0.5-mL of sample. 20 µl of sterile No. 2 fuel oil was then carefully placed on top of the solution in each well. Each microtiter plate was then incubated for fourteen (14) days at 20° C. At the conclusion of the fourteen-day incubation period, 100 µl of p-iodotetrazolium violet dye was added to each well and the results were recorded after at least 45 minutes to 2 hours of reaction time. Appearance of a pink to purple color constituted a positive test (continued microbial viability).

Each replicate sacrificed was extracted with an initial volume of 50-mL dichloromethane (DCM) for the chemical analysis. The sample was first extracted three times with 10-mL aliquots of the DCM. The remaining 20-mL was used to rinse the separatory funnel and added to the first 30mL of extract. Just prior to the initial extraction, each replicate is spiked with 100 µl of a surrogate-recovery standards stock solution. This stock solution was made up of 500 mg/L 5α androstane and d_{10} -phenanthrene. The separatory funnel was then capped and shaken vigorously for approximately thirty seconds to insure good mixing between phases. After mixing, the separatory funnel was allowed to sit for up to three hours to insure the greatest amount of separation between phases. This was done because of the presence of thick emulsions caused by microbiological activity. After a period of up to three hours, a 10-mL aliquot of the extract is poured into a 40-mL amber vial with a Teflon[™] lined cap, and taped with Teflon[™] tape. The samples were then stored in a 4° C walk-in refrigerator until retrieval for gravimetric analysis. The extraction was completed by filtering the remaining 40ml of DCM through a glass filter containing 20 grams of anhydrous sodium sulfate (Na₂SO₄) and into a 250-ml flat-bottom distillation flask. The Na₂SO₄ was rinsed with DCM until all traces of oil were removed from the funnel. The 250-ml flat-bottom distillation flask was placed on a Rotovap distillation unit until a volume of 10-ml was attained. Approximately 50-ml of hexane was added to the DCM extract and distilled to a volume of 10-ml. Another 50-ml of hexane was added to the hexane extract and distilled down to a final volume of 10-ml. A 1-ml aliquot of the final extract was removed and prepared for analysis on the GC/MS.

The gravimetric analysis was accomplished by first weighing an empty 40-mL vial and recording the weight. The 10-mL aliquot of extract was then placed in the vial, weighed and concentrated to dryness using a nitrogen gas blowdown technique. The remaining sample was then weighed and subjected to nitrogen blowdown for another ten to fifteen minutes. This was repeated once more to insure that the weight had changed no more than 5% weight difference between the second and third blowdown. If there was greater than a 5% difference, the sample was subjected a final blowdown to insure complete dryness. Weights were recorded after each blowdown, and then subjected to statistical analysis discussed below.

*The GC/MS analysis was subcontracted to Louisiana State University-IES, 42 Atkinson Hall, Baton Rouge, Louisiana, 70803.

STATISTICAL METHODS

GC/MS Data

Surrogate-adjusted data or rank-transformed surrogate adjusted data were analyzed using the MinitabTM 13.3 program. The computer program, unlike many others, is powerful enough to analyze unbalanced sets (uneven replication) of data using a general linear multiple factor ANOVA model. The probability of a type I error (α) was set apriori to 0.05.

Data sets were first analyzed for normality using the Anderson-Darling Goodness of Fit test. This test compares plot points with the normal theoretical distribution. Minitab calculates the statistic, above which there is a danger of non-normality. This is then compared to the chosen (preset by program), alpha (α) level of 0.01. For least-squares estimation, Minitab calculates a Pearson correlation coefficient. If the distribution fits the data well, then the plot points on a probability plot will fall on a straight line. The correlation measures the strength of the linear relationship between the X and Y variables on a probability plot. The correlation will range between 0 and 1, with higher values indicating a better fitting distribution.

Data passing a formal test for normality may not, strictly speaking, come from a normal distribution. Data that has sufficient linearity as shown by the passing results of a formal test for normality, may have attributes that weaken the ANOVA and Dunnett's test's ability to detect statistically significant differences between treatments (Zar, 1984).

Routine transformations were not amenable to non-normal data so an acceptable procedure for multiple-comparison ANOVA was found by using the rank-transformation test (Helsel, 1993). This technique first rank transforms the data and subjects it to the same multiple factor ANOVA test. This allows for an acceptable multiple comparison non-parametric test. After the program calculated the "F" and "P" statistics, the data were automatically subjected to Dunnett's means comparison test for comparison between treatments and controls.

Tables below give the final adjusted P-Values. Values of less than 0.05 (chosen α) indicate statistical significance. The T-Value is a ratio of the Difference of Means and Standard Error of Difference and indicates the degree and direction of the difference.

Microbiological Data

Microbiological data was analyzed with the Environmental Protection Agency's Most Probable Number Calculator, designed by the Risk Reduction Engineering Laboratory, Cincinnati, Ohio. This program calculates the most probable number (mpn) per mL with Salama correction for bias, and a Spearman-Karber Estimate. The program is based on the number of positive reactions in each of six replicates per serial dilution made. Confidence limits are included in the output of the program.

Gravimetric Data

Gravimetric data were analyzed with a simple two sample t-test available on the MinitabTM 13.3 program which compares the Day 0, 7, or 28 control means with their respective treatment means for statistical significance. The calculated p-Value is then compared to the chosen alpha (α) level of 0.05, as in the ANOVA analysis above. If the calculated value exceeds the 0.05, there is no statistical significance.

RESULTS AND DISCUSSION

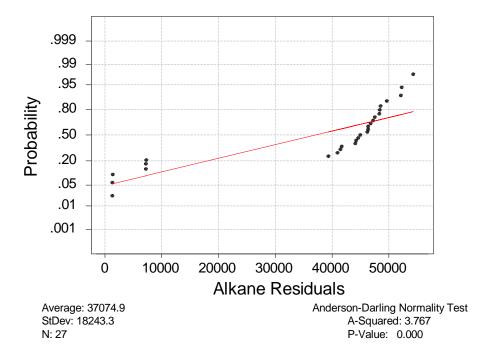
GC/MS Data

Results of the statistical analysis for the surrogate-adjusted data are reported and discussed below. Results for transformed data, if transformations were necessary, are discussed first, followed by the non-transformed data. Actual data (raw followed by surrogate-adjusted) are presented in the Appendices. GC/MS spectra appear in APPENDIX I along with computer printouts of the MinitabTM ANOVA analysis discussed below, which appear in APPENDIX II.

Sarva Bio Remed, LLC "SpillRemed" Marine Product & EPA Nutrient Solution

Surrogate-Adjusted Alkane Data

Preliminary analysis of surrogate-adjusted alkane data for normality (fig.1) showed the raw data to be non-normal with an Anderson-Darling P-statistic of 0.000. This is below the selected α -level of 0.01 and indicates the data are not normally distributed. Further visual evidence of the data's non-linearity can be seen in the probability plot for residuals of the data (fig.2). The data were rank-transformed and reanalyzed for normality (fig.3) giving an Anderson-Darling statistic of 0.585, well above the chosen α -level of 0.01. The probability plot for the residuals (fig. 4) of the data still show a small degree of non-linearity which can slightly lower the ANOVA and Dunnett's test ability to detect a statistical difference between treatments and controls. More on this subject is discussed in the conclusions.



Normality Test of Non-Transformed Alkane Data

Figure 1. – Anderson-Darling test for normality showing non-linearity of surrogate adjusted alkane data.

Normal Probability Plot for Non-Transformed Alkane Data

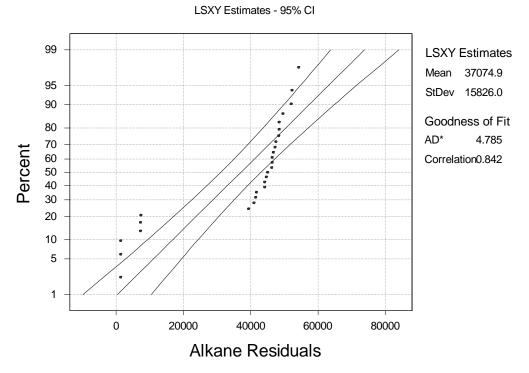


Figure 2. – Probability plot of the surrogate-adjusted alkane residuals showing further evidence of non-linearity.

Normality Test of Rank-Transformed Alkane Data

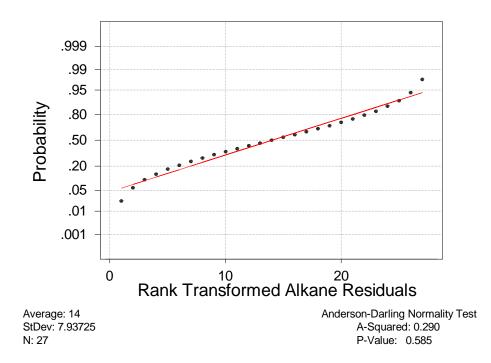


Figure 3. - Anderson-Darling test for normality showing improved linearity of the rank transformed surrogate-adjusted alkane data.

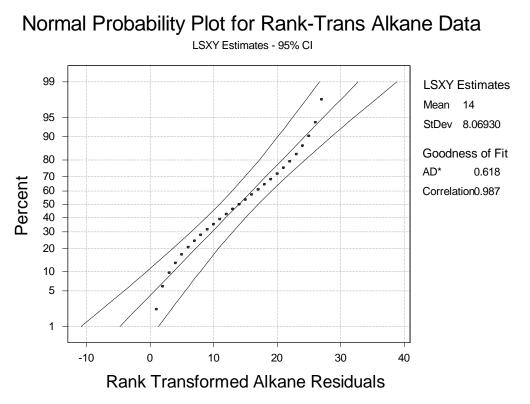


Figure 4. – Probability plot of the rank-transformed surrogate-adjusted alkane residuals showing improved linearity.

Non-transformed and rank-transformed surrogate-adjusted alkane data were analyzed with the General Linear ANOVA Model and Dunnett's multiple comparison tests between treatments and controls. P-statistics calculated by the ANOVA table for non-transformed and transformed treatment main effects, and treatment/day interactions are all under the chosen alpha (α) level of 0.05 indicating a significant difference between one or more treatments over one or more days.

Adjusted P-values for non-transformed and transformed data SPILLREMED Days 7 and 28 are shown to be significantly less than the Day 0 controls (Table 3). Adjusted P-values for non-transformed data SPILLREMED Days 7 and 28 are shown to be significantly less than the Day 7 controls (Table 4). The adjusted P-value for transformed data SPILLREMED Days 7 and 28 is also shown to be significantly less than the Day 7 controls (Table 4). Both transformed and non-transformed data demonstrated statistically significant differences between the Day 28 controls and the SPILLREMED Day 28 results (Table 5).

The Nutrient control behaved in the same manner as the product, showing the same significant differences between the Days 7 and 28 results from both the Day 0, Day 7, and Day 28 controls. However, using Tukey's pairwise means comparison method on non-transformed data, the Day 28 SPILLREMED product is also significantly less than the Nutrient alone (Table 6).

Table 2. ANOVA on non-transformed

ANOVA non-transformed data							
Source	DF	Seq SS	Adj SS	Adj MS	F	Р	
Day	2	5046904846	5046904846	2523452423	945.67	0.000	
Treatment	2	1425384113	1425384113	712692056	267.08	0.000	
Treatment*Day	4	2132924217	2132924217	533231054	199.83	0.000	
Error	18	48031825	48031825	2668435			
Total	26	8653245001					

ANOVA on rank transformed data

Source	DF	Seq SS	Adj SS	Adj MS	F	Р
Day	2	904.22	904.22	452.11	68.97	0.000
Treatment	2	349.56	349.56	174.79	26.66	0.000
Treatment*Day	4	266.22	266.22	66.56	10.15	0.000
Error	18	118.00	118.00	6.56		
Total	26	1638.00				

Table 3. Dunnett's test results using the Day 0 control as the control level vs. all other treatments
and controls (all interactions). Note - non = non-transformed data, trans = transformed data

Truckment		Difference of Means		T-Value		Adjusted P-Value	
Treatment	Day	NON	TRANS	NON	TRANS	NON	TRANS
Nutrient	0	-2000	-4.00	-150	-1.913	0.5604	0.3168
SPILLREMED	0	3100	4.33	2.32	2.073	0.1598	0.2457
Control	7	-1357	-1.00	-1.02	478	0.8693	0.9979
Nutrient	7	-8110	-12.67	-6.08	-6.059	0.0001	0.0001
SPILLREMED	7	-5727	-10.67	-4.29	-5.102	0.0029	0.0005
Control	28	-3496	-6.33	-2.62	-3.030	0.0924	0.0413
Nutrient	28	-41753	-16.33	-31.30	-7.813	0.0000	0.0000
SPILLREMED	28	-47653	-19.33	-35.73	-9.248	0.0000	0.0000

Table 4. Dunnett's test results using the Day 7 control as the control level vs. all other treatments and controls (all interactions). Note - non = non-transformed data, trans = transformed data

		Difference	of Means	T-V	alue	Adjusted	P-Value
Treatment	Day	NON	TRANS	NON	TRANS	NON	TRANS
Control	0	1357	1.00	1.02	0.478	0.8693	0.9979
Nutrient	0	-643	-3.00	048	-1.435	0.9978	0.6039
SPILLREMED	0	4457	5.33	3.34	2.551	0.0218	0.1055
Nutrient	7	-6753	-11.67	-5.06	-5.581	0.0006	0.0002
SPILLREMED	7	-4370	-9.67	-3.28	-4.624	0.0249	0.0014
Control	28	-2139	-5.33	-1.60	-2.551	0.4921	0.1055
Nutrient	28	-40397	-15.33	-30.29	-7.335	0.0000	0.0000
SPILLREMED	28	-46297	-18.33	-34.71	-8.770	0.0000	0.0000

Table 5. Dunnett's test results using the Day 28 control as the control level vs. all other treatments and controls (all interactions). Note - non = non-transformed data, trans = transformed data

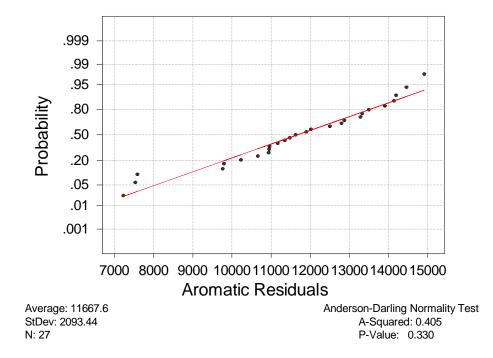
		Difference of Means		T-Value		Adjusted P-Value	
Treatment	Day	NON	TRANS	NON	TRANS	NON	TRANS
Control	0	3496	6.33	2.62	3.030	0.0924	0.0413
Nutrient	0	1496	2.33	1.12	1.116	0.8110	0.8143
SPILLREMED	0	6596	10.67	4.95	5.102	0.0007	0.0005
Control	7	2139	5.33	1.60	2.551	0.4921	0.1055
Nutrient	7	-4614	-6.33	-3.46	-3.030	0.0170	0.0413
SPILLREMED	7	-2231	-4.33	-1.67	-2.073	0.4494	0.2457
Nutrient	28	-38257	-10.00	-28.68	-4.783	0.0000	0.0010
SPILLREMED	28	-44157	-13.00	-33.11	-6.218	0.0000	0.0001

Table 6. Tukey's pairwise means comparison results between the Day 28 Nutrient and the Day 28 SPILLREMED non-transformed alkane data.

Treatment	Day	Difference of Means	T-Value	Adjusted P-Value
SPILLREMED	28	-5900	-4.424	0.0078

Surrogate-adjusted Aromatic Data

Preliminary statistical analysis for normality showed the raw data to be normal with an Anderson-Darling statistic of 0.033 (fig.5). This is above the selected alpha (α) level of 0.01 and indicates the data are marginally normal. Further visual inspection of data residuals reveals that the data residuals have fair linearity (fig.6). The data were then analyzed with the General Linear multiple-comparison ANOVA model and Dunnett's means comparison test using Day 0, 7, and 28 controls as control levels.



Normal Probability Plot

Figure 5. - Anderson-Darling test for normality showing linearity of the surrogate adjusted aromatic data.

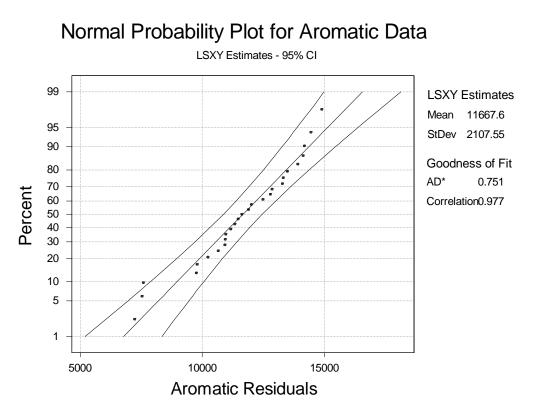


Figure 6. – Probability plot of the surrogate-adjusted aromatic residuals showing further evidence of linearity.

Surrogate-adjusted aromatic weight data were analyzed with the General Linear multiplecomparison ANOVA model and Dunnett's means comparison test using Day 0, 7, and 28 control for control levels (Table's 8, 9, and 10). P-statistics calculated by the ANOVA model on nontransformed data treatment main effects are under the chosen alpha (α) level of 0.05 indicating a significant difference between one or more treatments over one or more days.

ANOVA						
Source	DF	Seq SS	Adj SS	Adj MS	F	Р
Day	2	68510796	68510796	34255398	114.47	0.000
Treatment	2	10785102	10785102	5392551	18.02	0.000
Treatment*Day	4	29262580	29262580	7315645	24.45	0.000
Error	18	5386510	5386510	299251		
Total	26	113944989				

Table 7. ANOVA on Surrogate-adjusted Aromatic Data

Analysis of the non-transformed data shows SPILLREMED Day 7 and Day 28 adjusted P-values to be significantly less than the Day 0 controls (Table 8). The data shows adjusted P-values for SPILLREMED Day 28 to be significantly less than Day 7 controls, but the Day 7 SPILLREMED p-value is not significantly different than the Day 7 control (Table 9). Adjusted P-values for SPILLREMED Day 28 are also significantly less than the Day 28 controls (Table 10).

The Nutrient alone behaved in the same manner as the product, showing the same significant differences between the Days 7 and 28 results from the Day 0 controls. The Nutrient at Day 28 was also significantly less than the Day 28 control, but not the Day 7 control. Also, using Tukey's pairwise means comparison method on the non-transformed data, the Day 28 SPILLREMED product is also significantly less than the Nutrient alone (Table 11).

Treatment	Day	Difference of Means	T-Value	Adjusted P-Value
Nutrient	0	295	0.66	0.9843
SPILLREMED	0	659	1.48	0.5766
Control	7	-2221	-4.97	0.0007
Nutrient	7	-2155	-4.82	0.0009
SPILLREMED	7	-2407	-5.39	0.0003
Control	28	-924	-2.07	0.2474
Nutrient	28	-3556	-7.96	0.0000
SPILLREMED	28	-6041	-13.53	0.0000

Table 8. Dunnett's test using Day 0 control as the control level vs. all other treatments and controls (all interactions). No transformation.

Table 9. Dunnett's test using Day 7 Control as the control level vs. all other treatments and controls (all interactions). No transformation.

Treatment	Day	Difference of Means	T-Value	Adjusted P-Value
Control	0	2221	4.973	0.0007
Nutrient	0	2516	5.634	0.0002
SPILLREMED	0	2880	6.449	0.0000
Nutrient	7	66	0.149	1.0000
SPILLREMED	7	-186	-0.416	0.9992
Control	28	1297	2.905	0.0532
Nutrient	28	-1335	-2.989	0.0449
SPILLREMED	28	-3820	-8.552	0.0000

Table 10. Dunnett's test using Day 28 control as the control level vs. all other treatments and controls (all interactions). No transformation.

Treatment	Day	Difference of Means	T-Value	Adjusted P-Value
Control	0	924	2.07	0.2474
Nutrient	0	1219	2.73	0.0751
SPILLREMED	0	1583	3.54	0.0142
Control	7	-1297	-2.90	0.0532
Nutrient	7	-1231	-2.76	0.0713
SPILLREMED	7	-1483	-3.32	0.0227
Nutrient	28	-2632	-5.89	0.0001
SPILLREMED	28	-5117	-11.46	0.0000

Table 11. Tukey's pairwise means comparison results between the Day 28 Nutrient and the Day 28 SPILLREMED non-transformed aromatic data.

Treatment	Day	Difference of Means	T-Value	Adjusted P-Value
SPILLREMED	28	-2485	-5.564	0.0008

Microbiological Analysis Data

The following tables show the most probable number calculated by EPA's most probable number calculator Version 4.04. The data show the continued viability of the organisms in the product and to a lesser extent, the controls. Print outs appear in APPENDIX III

Day 0	Day 7	Day 28
6704	7120	9549
6195	6956	8724
6330	7647	9279
6014		7,068,457
	<i>i</i>	7,235,388
	,	7,250,977
		1.38E+08
		1.55E+08 1.57E+08
	6794 6195 6330	6794 7129 6195 6956 6330 7647 6014 28,566 6330 27,514 15,565 1,535,074 15,685 1,399,693

Table 12. Micro Results, MPN (per mL)

<u>Gravimetric Data</u>

The following tables show the P-Values calculated by the two-sample t-test of the MinitabTM program. Table 13 shows the calculated values for Day 28 controls the Day 28 product, and the p-value of the comparison is lower than the chosen alpha (α) level of 0.05 and therefore indicate statistical significance. A computer printout of the analyses can be seen in APPENDIX IV. Table 14 shows that the calculated values for the Day 28 controls and both the Day 7 and Day 28 nutrient are both statistically significant.

Table 13.	P-Values calculated by the two-sample t-test for product (SPILLREMED) and the
controls		

Treatments	Day	Treatment Weight Means (mg)	T-test Scores	p-value
Controls	0	0.101	-2.79	0.068
SPILLREMED	0	0.108	-2.19	0.000
Controls	7	0.098	1.04	0.373
SPILLREMED	7	0.096	1.04	0.575
Controls	28	0.095	42.25	0.001
SPILLREMED	28	0.016	72.23	0.001

Table 14. P-Values calculated by the two-sample t-test for the nutrient and the controls

Treatments	Day	Treatment Weight Means (mg)	T-test Scores	p-value
Controls	0	0.101	1.36	0.308
NUTRIENT	0	0.098	1.50	0.508
Controls	7	0.098	10.07	0.001
NUTRIENT	7	0.085	10.07	0.001
Controls	28	0.095	33.84	0.000
NUTRIENT	28	0.024	55.04	0.000

Conclusions

Our conclusions will begin with a discussion of the GC/MS due to its relative importance in judging the tested product effective. A discussion of the microbiological results and gravimetric results will follow.

GC/MS Data

SarvaBio Remed, LLC Product (SpillRemed) & EPA Nutrient mix (NUT)

Surrogate-adjusted Alkane Data

Surrogate-adjusted alkane SpillRemed data was shown to be non-normal and had to be ranktransformed to attain an acceptable degree of linearity. Analysis of the surrogate-adjusted data with ANOVA and Dunnett's test did however show the product treatments at Day 7 and 28 to be significantly less than Day 0, 7, and 28 controls. The extreme non-linearity of the nontransformed data makes the results of the ANOVA and Dunnett's test less reliable. The data, upon rank-transformation, achieved the desired linearity showing Day 7 and 28 product to be significantly less that the respective Day 0 and 28 controls. Based on this parameter the product appears to be effective.

Surrogate-adjusted alkane nutrient data was shown to be non-normal and had to be ranktransformed to attain an acceptable degree of linearity. Analysis of the surrogate-adjusted data with ANOVA and Dunnett's test did however show the nutrient treatments at Day 7 and 28 to be significantly less than their respective controls. The non-linearity of the non-transformed data may make the results of the ANOVA and Dunnett's test less reliable, however. The data, upon rank-transformation, achieved the desired linearity showing Day 28 nutrient to be significantly less than the respective Day 0, 7, and 28 controls. Based on this parameter the nutrient treatment alone appears to be effective.

Tukey's test on untransformed alkane data also showed a significant difference between the Day 28 SpillRemed results and Day 28 Nutrient results, indicating that the product seems to be more effective than nutrient treatment alone.

Surrogate-adjusted Aromatic Data

Surrogate-adjusted aromatic SpillRemed data were normally distributed based on the Anderson Darling normality test and needed no transformation to attain linearity. Residuals of the data were also fairly linear. The data showed the product at Day 28 to be significantly less than Day 0, 7, and 28 controls. Based on this parameter, the product should be deemed effective.

Surrogate-adjusted aromatic nutrient data were normally distributed based on the Anderson Darling normality test and needed no transformation to attain linearity. Residuals of the data were also fairly linear. The data showed the nutrient at Day 28 to be significantly less than Day 0, 7, and 28 controls. Based on this parameter, the nutrient treatment should be deemed effective.

Tukey's test on the aromatic data also showed a significant difference between the Day 28 SpillRemed results and Day 28 Nutrient results, indicating that the product seems to be more effective than nutrient treatment alone.

Microbiological Results

Sarva Bio Remed, LLC Product (SpillRemed)& EPA Nutrient (NUT)

The microbiological results speak for themselves. They show a definite continued microbiological viability over time for the product treatments.

Similar to the product treatment and the nutrient treatments show a definite continued microbiological viability over time.

Gravimetric Results

SarvaBio Remed, LLC Product (SpillRemed)& EPA Nutrient (NUT)

Gravimetric results showed statistical significance between products and controls by Day 28. This tends to support the bulk of the data seen in both GC/MS analysis and microbiological analysis.

Gravimetric results showed a statistical significance between the Nutrient and the control on Day 7 and Day 28. This data tends to support the bulk of the data in both GC/MS analysis and microbiological analysis.

Discussion on Surrogate Recovery – QA/QC

The purpose of incorporating surrogate recovery percentages into the raw data is to check the efficiency of extraction techniques and in most cases is a valid quality control check. The acceptable range of surrogate recovery percentages is given in the cited Federal Register document titled Environmental Protection Agency, (EPA) Pt. 300, Appendix C, page 237, as 70%-120%. Percentage recoveries for product and controls for Day 0, Day 7 and Day 28 are given in Table 15 below.

,	Treatment	Day 0	Day 7	Day 28
Control	5-Alpha Andorstane	0.97	0.90	0.97
Rep#1	Phenanthrene-d10	0.97	0.94	0.88
Control	5-Alpha Andorstane	0.98	0.91	0.97
Rep#2	Phenanthrene-d10	0.98	0.96	0.90
Control	5-Alpha Andorstane	0.99	0.92	0.96
Rep#3	Phenanthrene-d10	0.99	0.93	0.89
NUT	5-Alpha Andorstane	1.00	0.79	0.99
Rep#1	Phenanthrene-d10	1.00	0.82	0.74
NUT	5-Alpha Andorstane	0.99	0.86	1.00
Rep# 2	Phenanthrene-d10	0.99	0.93	0.77
NUT	5-Alpha Andorstane	0.99	0.92	1.00
Rep# 3	Phenanthrene-d10	1.00	0.97	0.75
SpillRemed	5-Alpha Andorstane	0.98	0.81	0.83
Rep #1	Phenanthrene-d10	0.98	0.82	0.75
SpillRemed	5-Alpha Andorstane	0.99	0.91	0.78
Rep# 2	Phenanthrene-d10	0.99	0.97	0.71
SpillRemed	5-Alpha Andorstane	1.00	0.88	0.79
Rep# 3	Phenanthrene-d10	1.00	0.89	0.71

Table 15. Surrogate recovery percentages.

As Table 15 shows, the bulk of surrogate recovery percentages decline over time. However, all surrogate recoveries were within specifications of the protocol.

Statistical Analysis

Lastly, we feel that the nature of the data may reduce the ANOVA and Dunnett's means comparison test to detect a legitimate statistical effect between treatments and controls. Before the data can be subjected to the ANOVA analysis, it must pass a "normality" test where a calculated P-value is compared to a chosen alpha (α) level (usually 0.01). ANOVA has reduced power to detect a significant statistical difference when analyzing non-normal data (Zar, 1984). However, data that passes a formal test for normality is not necessarily from a "normal distribution" strictly speaking. A test for normality looks for linearity, which is only one aspect of the assumptions of normality. The data may also be skewed to the left or right as indicated by measurement of the median, may have 'heavy tails" in the distribution or may contain outliers. Normality after all, is usually a matter of degrees and not just whether the data are, or are not normally distributed. If data are not normal in the strictest sense, we feel the test's ability to detect subtle but significant statistical differences may be compromised to some degree.

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APPENDIX I

8	CONTROL, REPLICATE 1	-	0	CONTROL, REPLICATE 2		8	CONTROL, REPLICATE 3	
	Testing Date: Day 0 Initial Oil Weight: 510 mg Final Extracted Volume: 10 ml	mg 10 ml	Einal	Testing Date: Day 0 Initial Oil Weight: 500 mg Final Extracted Volume: 10 ml	Ξ	Linit Final F	Testing Date: Day 0 Initial Oil Weight: 490 mg Final Extracted Volume: 10 ml	6 m
	(encontration (nq/mq)	Surrogate Corrected (rq/mq)		Concentration (ng/mg)	Surrogate Corrected (aq/mg)		Concentration (ng/mg)	Surrogate Corrected (ng/mg)
Alkane Analyte:			Alkane Analyte:			Alkane Analyte:		
nC-10 Decare	1400	1400	nc-10 Decane	1400	1400	nC-10 Decane	1400	1400
nC-11 Undecare nC-12 Dodecare	2600	2700	nc-11 undecane nc-12 Dodecane		2700	nC-11 Undecane nC-12 Dodecane	2300	2300
nC 13 Tridecare	2400	2500	nC 13 Tridecane		2700	nC 13 Tridecane	2400	2400
nC-14 Tetradecane	2400	2500	nC-14 Tetradecane	2600	2700	nC-14 Tetradecane nC-15 Pentadecane	2300	2300
nC-16 Hexadecane	2200	2300	nC-16 Hexadecane	2300	2400	nC-16 Hexadecane	2000	2000
nC-17 Heptadecane	2200	2300	nC-17 Heptadecane	2300	2400	nC-17 Heptadecane	2100	2100
Pristane nC-18 Octadecane	2100	1500	Pristane IIC-18 Octadecane	1600	1600 2600	Pristane nC-18 Octadecane	1500	1500
2 5	1400	1400	Phytane	1500	1500	Phyteno	1300	1300
nC-19 Nonadecane	1900	2000	nC-19 Nonadecane	2200	2300	nC-19 Nonadecane	2000	2000
nC-20 Elcosare	2000	2100	nC-20 Ecosane	2300	2400	nu-20 bicosane	2000	2000
nC-21 Henelcosane	1900	2000	nC-21 Heneicosane	2100	2200	nC-21 Heneicosane	1900	1900
nC.23 Trimsare	1500	1500	nC-23 Trimsane	1700	1700	nC-23 Tricnsane	1500	1500
nC-24 Tetracosane	1400	1400	nC-24 Tetracosane	1500	1500	nC-24 Tetracosane	1500	1500
n0-25 Pentacosane	2000	2100	nC-25 Pentacosane		2000	nC-25 Pentacosane	1900	1900
nC-26 Hexacosane	1500	2100	nC-26 Hexacosane		2300	nC-26 Hexacosane	1900	1400
nC-26 Octacosane	1200	1200	nc-28 Octacosane		0021	nC-28 Octacosane	1200	1200
nC-29 Nonacosane	1200	1200	nC-29 Nonacosane	1200	1200	nC-29 Nonacosane	1200	1200
nC-30 Triacontane	1200	1200	nC-30 Triacontane	1300	1300	nC-30 Triacontane	1200	1200
nC-31 Hentriacontane	1100 880	1100	nC-31 Hentriacontane	1200	1200	nC-31 Hentriacontane	1100	1100
nC-33 Tritriacontane	940	026	nC-33 Tritriacontane	1000	1000	nC-33 Tritriacontane	870	88
nC-34 Tetratriacontane	740	760	nC-34 Tetratriacontare	780	800	nC-34 Tetratriacontane	740	750
	8	6	UC-00 Periodiana	0071	0071		8	8
Total Alkancs	47060	48430	Total Alkancs	50760	52100	Total Alkancs	46330	46360
Aromatic Analyte:			Aromatic Analyte:	ŝ		Aromatio Analyte:	į	Į
C1-Naphthalenes	1600	1700	C1-Naphthalenes	1700	1700	C1-Naphthalenes	1500	1500
C2-Naphthalenes	2100	2200	C2-Naphthalenes	2400	2400	C2-Naphthalenes	2000	2000
C3-Naphthalonca	1700	1800	C3-Naphthalones	1900	1900	C3-Naphthalonca	1600	1600
	110	110	Fluorene	120	071	+140rene	100	100
C1-Fluorenes	260	270	C1-Fluorenes	280	290	C1-Fluorenes	260	260
C2-Fluorenes	000	046	C2-Fluorenes	410 260	420 270	C2-Fluorenes	020 239	020
Diberzothiophene	240	250	Dibenzothionhene	270	280	Dibenzothiophene	240	240
C1-Dibenzothiophenes	440	460	C1-Dibenzothiophenes	500	510	C1-Dibenzothicphenes	420	420
C2-Dibenzothiophenes	540	560 360	C2-Dibenzothiophenes	600	610	C2-Dibenzothicphenes	500	500
Fhenanthrene	310	320	Phenanthrene	320	330	Phenanthrene	300	300
C1-Phenanthrenes	530	550	C1-Phenanthrenes	065	600	C1-Phenarthrenes	530	530
C3-Phenanthrenes	420	440	C3-Ptrenanthrenes	490	500	C3-Phenarthrenes	420	420
C4-Phenanthrenes	130	140	C4-Phenandirenes	130	130	C4-Pilenar Uneries	130	130
Fluorentere	2.7	3.1	Anthracene	2.4	25	Anthracene Fluoranthene	3.0	3.0
Pyrene	8.2	8.5	Pyrene	8.7	8.8	Pyrene	6	8.9
C1- Pyrenes	110	110	C1- Pyrenes	110	110	C1- Pyrenes	110	110
C2 Hyrenes	140	150	C2 Pyrenec	160	160	C2 Pyrenec	140	140
C4- Pyrenes	92	8	C4- Pyrenes	100	100	C4- Pyrenes	88	88
Naphthohenzothinphene	54	95	Naphthohen7othinphere		60	Naphthohenzothinphene	50	50
C-2 Naphthobenzothiophenes	180	190	C-2 Naphthobenzothiophenes		210	C-1 reprint over connormer es	180	8
C-3 Naphthobenzothiophenes	150	160	C-3 Naphthobenzothiophenes	150	150	C-3 Naphthobenzothiophenes	140	140
Chrysene	5 05	6	Chrysene		99	Chrysene	- 95	
	69	7.6	C1- Chrysenes		£	C1- Chrysenes	93	84 8
C2- Chrysenes C3- Chrysenes	120	120	C2- Chrysenes	5 8	120	C2- Chrysenes C3- Chrysenes	1 18	110
C4- Chrysenes	64	5	C.4- Chrysenes	47	47	C.4- Chrysenes	47	8
Benzo (b) Fluoranthene	6.C	0.4	Benzo (b) Fluoranthere	3.7	3.8	Benzo (b) Fluoranthene	9 9 9	3.5
Benzo (e) Pyrene	7.9	8.2	Benzo (e) Pyrene	8.8	8.9	Benzo (e) Pyrene	3.7	2.6
Benzo (a) Pyrene	1.2	1.2	Benzo (a) Pyrene	12	12	Benzo (a) Pyrene	1.1	1.1
heryrene Incieno (1,2,3 - cd) Pyrene	0.0	0.00	Perylene Indeno (1,2,3 - cd) Pyrene	0.0	0.00	Perylene Indeno (1,2,5 - cd) Pyrene	*0:0	8.0
Dibenzo (a,h) anthracene	0.00	0.00	Dibenzo (a,h) anthracere	0.00	0.00	Elbenzo (a,h) anthracene	0.00	0.00
I otal Aromatics	12912	13499	I otal Aromatics	14307	14455	I otal Aromatics	1244/	12449
% Surrogata Recovery			% Surrogate Recovery			% Surrogate Recovery	-	
5 Alpha Androetane Phonorthrono d-10	0.97	1.00	5 Alpha Androetane Phonorthrond du/0	0.98 0 0	1.00	5 Alpha Androetane Phonanthrond d_10	0:08 00:00	00 F
0	0.07	1.00	l'honanthrono a-10	0.08	1.00	I'honanthrono a-ru	0.09	1.00

			-					
	Testing Date: Day 0 Initial Oil Weight: 510 mg	6	= <u>e</u>	Testing Date: Day 0 Initial Oil Weight: 510 mg			Testing Date: Day 0 Initial Oil Weight: 510 mg	5
Lina	Final Extracted Volume: 10 mL	10 mL	Final	Final Extracted Volume: 10 mL) mL	Final	Final Extracted Volume: 10 mL	0 mL
	Concentration (ng/mg)	Surregate Correted (rg/mg)		Concentration (ng/mg)	Surrogate Corrected (ng/mg)		Concentration (ng/mg)	Surrogate Corrected (ng/mg)
Alkane Analyte: nC-10 Decare	730	730	Alkane Analyte: nC-10 Decane	200	710	Alkane Analyte: nC-10 Decane	720	730
nC-11 Undecare	1800	1800	nC-11 Undecane	1700	1700	nC-11 Undecane	1800	1800
nC 13 Tridecare		1800	nC 13 Tridecane	1800	1800	nC 13 Tridecane	1900	1900
nC-14 Tetradecane	2200	2200	nC-14 Tetradecane	2000	2000	nC-14 Tetradecane	2200	2200
nc-15 rentarecane nc-16 Hexadecane	1800	1800	nC-15 Hexadecane	1800	1800	nC-16 Hexadecane	1800	1800
nC-17 Heptadecane	1900	1900	nC-17 Heptadecane	1900	1900	nC-17 Heptadecane	2000	2000
Pristane - 18.0 claderate	7100	1600	Pristane	1700	1700	Pristane IIC-18 Orderbersene	1600 2400	1600
Phytano	1500	1500	Phytano	1600	1600	Phytene	1600	1600
nC-19 Nonadecane	1800	1800	nC-19 Nonadecane	1800	1800	nC-19 Nonadecane	1900	1900
nC-20 Elosare nC-21 Heneionsane	2000	2000	nC-20 Ecosane nC-21 Heneionsane	2000	2000	nC-20 bicosane nC-21 Heneicneane	2000	2000
nc-22 Docosare	1600	1600	nC-22 Docosane	1700	1700	nC-22 Docosane	1800	1800
nC-23 Tribnsare	1600	1600	nC-23 Trionsane	1600	1600	nC-23 Trinnsane	1800	1800
nC-24 Tetracosane	1700	1600	nC-24 Tetracosane	1800	1800	nC-24 Tetracosane	1800	1800
n-20 Haveneere	0012	0012	nu-zu remacosane	0022	0077	PU-20 Fentacosane	0002	0002
nC-27 Heptacosane	1900	1900	nc-27 Heptacosane	2000	2000	nC-27 Heptacosane	2100	2100
nC-26 Octacosane	1700	1700	hC-28 Octacosane	0061	1900	nC-28 Octacosane	1700	1700
nC-29 Nonacosane	1800	1800	nC-29 Nonacosane	1800	1800	nC-29 Nonacosane	1700	1700
nC-30 Triacontane	1200	1200	nC-30 Triacontane	1200	1200	nC-30 Triacontane	1200	1200
nc-or remanuation nc-32 Dutriacontane	1100	1100	nc-or nerrinacontaire nc-32 Dutriacontaire	1100	1100	nc-32 Dutriacontane	1100	1100
nC-33 Tritriacontane	1100	1100	nC-33 Tritriacontane	1000	1000	nC-33 Tritriacontane	1200	1200
nC-34 Tetratriacontane	980	980	nC-34 Tetratriacontare	970 050	080	nC-34 Tetratriacontane	1100	1100
Total Alkancs	46310	46210	Total Alkancs	46320	46350	Total Alkanes	48320	48330
Aromatic Analyte: Manhthalana	440	440	Aromatic Analyte: Nanhthalene	480	480	Aromatic Analyte: Nanhthalene	510	510
C1-Naphthalenes	1600	1600	C1-Naphthalenes	1800	1800	C1-Naphthalenes	1700	1700
C2-Naphthalenes	2200	2200	C2-Naphthalenes	2400	2400	C2-Naphthalenes	2300	2300
C3-Naphthalonca	1800 860	1800 860	C3-Naphthalonca	1900	1900	C3-Naphthalonca C4 Naphthalanee	1000	1000
Fluorene	110	110	Fluorene	110	110	Fluorene	110	110
C1-Fluorenes	250	250	C1-Fluorenes	250	250	C1-Fluorenes	250	250
C2-Fluorences	000	000	C2-Fluorenes	020	020	C2-Fluorenes	000 220	800
Diberzothiophene	240	240	Dibenzothiphene	250	250	Dibenzothiophene	250	250
C1-Dibenzothiophenes	450	450	C1-Dibenzothiophenes	470	470	C1-Dibenzothicphenes	450	450
C2-Dibenzothiophenes	560 380	560 380	C2-Dibenzothiophenes	590 260	590 360	C2-Dibenzothicphenes	370	570
Phenanthrere	320	320	Phenanthrene	330	330	Phenanthrene	330	330
C1-Phenanthrenes	065	065	C1-Phenanthrenes	ROO	RDA	C1-Phenarthrenes	600	600
C3-Phenanthrenes	420	420	C3-Prenantrenes	470	470	C3-Phenarthrenes	470	470
C4-Plicitalibiliteties	130	130	C4-Pt enanthrenes	130	130	C4-Pienal Ilirenes	140	140
Anthracene	2.0	2.0	Anthracene	22	22	Auflitacene	2.1	2.1
Pyrene	5.3	23	Pyrene	5.9	5.9	Pyrene	6.1	6.1
	110	110	-	120	120		120	120
C2 Pyrenee	140	140	C2 Pyrenec	150	150	C2 Pyrenec	150 160	150
C4- Pyrenes	64	40	C4- Pyrenes	67	67	C4- Pyrenes	91	5
Naphthohenzothinphene	5	55	Naphthohen7ofthiphere		25	Naphthohenzothinphene	95	95
C-1 Naphthobenzothiophenes C-2 Naphthobenzothiophenes	1/0	1/0 180	C-1 Naphthobenzothiophenes C-2 Naphthobenzothiophenes		180	C-1 Naphtrobenzothiophenes C-2 Naphtrobenzothiophenes	190 200	200
C-3 Naphthobenzothiophenes	150	150	C-3 Naphthobenzothiophenes		150	C-3 Naphthobenzothiophenes	150	150
Bonzo (a) Anthracono Chrysene	± 8	± 68	Bonzo (u) Anthrucono Chrysone	1 8	‡ %	Bonzo (a) Anthracono Christene	13	13
C1- Chrysenes	1/20	120	C1- Chrysenes	130	130	C1- Chrysenes	140	140
C2- Chrysenes	150	150	C2- Chrysenes	160	160	C2- Chrysenes	170	170
C.4- Chrysenes	75	74	C.4- Chrysenes	70	70	C.4- Chrysenes	32	25
Benzo (b) Fluoranthene	2.9	2.9	Benzo (b) Fluoranthere	2.7	2.7	Benzo (b) Fluoranthene	2.7	2.7
Denzo (K) I luoranmene Benzo (e) Pyrene	7.6 6.6	F 6	Denzo (K) Iuorantinene Benzo (e) Pyrene	÷ e	- 0	Benzo (k) I luoranmene Benzo (e) Pyrene	2 6 2 6	63
Benzo (a) Pyrene	13	1.3	Benzo (a) Pyrene	14	1.4	Benzo (a) Pyrene	13	13
Perylene Inceno (1.2.3 - cd) Pvrene	0.00	0.00	Perylene Indeno (1.2.3 - cd) Pvrere	0.00	0.71	Perylene Indeno (1.2.3 - cd) Pyrene	0.72	0.72
Dibenzo (a,h) anthracene Benzo (g,h,i) perylene	0.00 1.6	0.00 1.6	Dibenzo (a,h) anthracene Benzo (g,h,i) perylene	0.00 1.7	0.00 1.7	Elbenzo (a,h) anthracene Benzo (g,h,i) perylene	0.00 1.6	0.00 1.6
Total Aromatics	13291	13287	Total Aromatics	14144	14140	Total Aromatics	13916	13911
% Surrogate Recovery 5 Aloha Androstane	1.00	1.00	% Surrogate Recovery 5 Aloha Androstane	0.99	1.00	% Surrogate Recovery 5 Aloha Androetane	0.99	1.00
Phonanthrono d-10	1.00	1.00	Phonanthrono d-10	1.00	1.00	Phonanthronc d-10	1.00	1.00

SARV	SARVA PRODUCT, REPLICATE 1	ATE 1	SARV	SARVA PRODUCT, REPLICATE 2	TE 2	SARVA	SARVA PRODUCT, REPLICATE 3	TE3
Ini	Initial Oil Weight: 510 mg Final Extracted Volume: 10 mL	mg 10 mL	Final	Initial Oil Weight: 510 mg Final Extracted Volume: 10 mL	0 mL	Init Final E	Initial Oil Weight: 510 mg Final Extracted Volume: 10 mL	0 m. 0
	Concentration (ng/mg)	Surregate Corrected (rg/mg)		Concentration (ng/mg)	Surragate Corrected (ng/mg)		Concentration (ng/mg)	Eurrogata Corrected (ng/mg)
Alkane Analyte: n0_10 Decare	1100		Alkane Analyte: nC_10 Decane	1200	unct	Alkane Analyte: nC_10 Decane	1200	1200
nC-11 Undecare	2100	2100	nC-11 Undecane	2100	2100	nC-11 Undecane	2100	2100
nC-12 Uotecare nC 13 Tridecare		2000	nu-12 Uddecane nC 13 Tridecane		2200	nC-12 Undecane nC 13 Tridecane	2300	2300
nC-14 Tetradecane	2200	2200	nC-14 Tetradecane	2500	2500	nC-14 Tetradecane	2400	2400
nC-16 Hexadecane	2300	2300	nc-16 Hexadecane	2600	2600	nC-16 Hexadecane	2500	2500
nC-17 Heptadecane	2400	2400	nC-17 Heptadecane	2600	2600	nC-17 Heptadecane	2400	2400
nc-18 Ocladecare	2200	2200	nic-18 Octadecare	2500	1000	InC-18 Ocladecane	2500	2500
Phytane P	1600	1600	Phytano	1700	1700	Phytene	1600	1600
nC-19 Nonadecane nC-20 Elcosare	2200	1900	nC-19 Nonapecane nC-20 Ecosane	2400	2400	nC-19 Nonadecane nC-20 Elcosane	2500	2200
nC-21 Heneicosane	2000	2000	nC-21 Heneicosane	2300	2300	nC-21 Heneicosane	2200	2200
nC-22 Docosare nC-23 Trinosare	1900	1900	nC-22 Docosane nC-23 Trimeane	2100	2100	nC-22 Docosane nC-23 Trionsane	2000	2000
nC-24 Tetracosane	2000	2000	nC-24 Tetracosane	2000	2000	nC-24 Tetracosane	1900	1900
nC-25 Pentacosane	2200	2200	nC-25 Pentacosane	2400 2600	2400	nC-25 Pentacosane	2300	2000
nc-zo nexacusarie nC-27 Heptacosane	1900	1900	nc-27 Heptacosane	2100	2100	nC-27 Heptacosane	1900	1900
nC-28 Octacosane	1700	1700	nC-28 Octacosane	0061	0061	nC-28 Octacosane	1700	1700
nC-29 Nonacosane	1100	1800	nC-29 Nonacosane	1900	1900	nC-29 Nonacosane	1700	1700
nC-31 Hentriacontane	1100	1100	nC-31 Hentriacontane	1300	1300	nC-31 Hentriacontane	1100	1100
nC-32 Dutriacontane	1000	1000	IIC-32 Dulriacontane	1100	1100	IIC-32 Dutriacontaire	1100	1100
nC-34 Tetratriacontane	1100	1100	nC-34 Tetratriacontare	1100	1100	nc-34 Tetratriacontane	1000	1000
nC-35 Pentatriacontane	920	940	nC-35 Pentatriacontare	056	960	nC-35 Pertatriacontane	930	940
Total Alkancs	49590	49630	Total Alkancs	54270	54290	Total Alkanes	52250	52270
Aromatic Analyte: Nanhthalere	570	280	Aromatic Analyte: Nanhthalene	270	580	Aromatio Analyte: Namhthalene	570	122
C1-Naphthalenes	1700	1700	C1-Naphthalenes	1800	1800	C1-Naphthalenes	1600	1600
C2-Naphthalenes C3-Naphthalenes	2300	2300	C2-Naphthalenes C3-Naphthalenes	2400	2400	C2-Naphthalenes C3-Naphthalenes	2100	2100
C4-Naphthalenes	800	810	C4-Naphthalenes	850	860	C4-Naphthalenes	190	790
Fluorene C1.Fluorenes	110	11U 260	Pluorene C1_Flinnenes	11U 280	110 280	Fluorene C1_Fluorenes	110	11U 260
C2-Fluorenes	000	040	C2-Fluorenes	410	420	C2-Fluorenes	000	38
C3- Flurrenes	360	370	C.3. Fluorenes	380	390	C3- Flurrenes	340	340
Ulberzothiophenes C1-Dibenzothiophenes	2/0	790	Ulbenzothiphenes C1-Dibenzothiphenes	290	290	Ulbenzothicphenes C1-Dibenzothicphenes	260 490	790
C2-Dibenzothiophenes	600	610	C2-Dibenzothiophenes	660	029	C2-Dibenzothicphenes	280	590
C3- Dibenzothiophenes Phenanthrere	380	390	C3- Dibenzothiophenes Phenanthrene	380	390	C3- Dibenzothicphenes Phenanthrene	330	350 330
C1-Phenanthrenes	640	650	C1-Phenanthrenes	700	710	C1-Phenarthrenes	610	610
C3-Phenanthrenes	480	490	C3-Phenanthrenes C3-Phenanthrenes	510 510	810 520	C3-Phenarthrenes C3-Phenarthrenes	450	450
C4-PlienanBirenes	150	150	C4-Plenandmenes	150	150	C4-Phenal threncs	150	150
Fluoranthere	2.4	2.4	Fluoranthene	2.5	2.5	Fluoranthene	2.5	2.5
Pyrene C1 Dyranes	7.1	7.2	Pyrene C1 Dyranee	7.0	7.1	Pyrene C1 Dyranas	6.5	6.3
C2 Fyrenee	110	140	C2 Pyrenec	150	150	C2 Pyrenec	130	1 <u>6</u>
C3- Pyrenes C4- Pyrenes	140 76	140	C3- Pyrenes C4- Pyrenes	140 81	140 82	C3- Pyrenes C4- Pyrenes	130	130
llaphthohen7othinphene	65		\aphthoben7othinphere		60	Naphthohen7othinphene.	Υ.	22
C-1 Naphthobenzothiophenes C-2 Naphthobenzothiophenes	180 190		C-1 Naphthobenzothiophenes C-2 Naphthobenzothiophenes		200 200	C-1 Naphthobenzothiophenes C-2 Naphthobenzothiophenes	180	180
C-3 Naphthobenzothiophenes	150		C-3 Naphthobenzothiophenes		160	C-3 Naphthobenzothiophenes	150	150
Chrysene	51 76	51	3	2 23	2 23	P.	16	-1
C1- Chrysenes	120	120	CI- Chrysenes	130	130		120	071
CG- Chrysenes	140	140	C2- Chrysenes C3- Chrysenes	110	110	C2- Chrysenes C3- Chrysenes	140	140
C4- Chrysenes	67	68	C4- Chrysenes	74	75	C4- Chrysenes	89 7	69
Benzo (b) Fluorantnene Denzo (k) Fluoranthene	8.5 9.5	9.9 9.4	Benzo (b) rluoranthere Denzo (k) fluoranthere	3.8 4.2	3.8 4.0	Denzo (b) Fluoranthene Denzo (k) Fluoranthene	9.6	3.4 0.0
Benzo (e) Pyrene Benzo (a) Pyrene	12	12	Benzo (e) Pyrene Benzo (a) Pyrene	5 <u>1</u>	13	Benzo (e) Pyrene Benzo (a) Pyrene	6 <u>6</u>	12
Perylene	0.61	0.62	Perylene	0.62	0.63	Perylene	0.60	0.60
Dbenzo (1,2,3 - cu) -yrans Dbenzo (a,h) anthracene Benzo (a h.) neorlene	0.00	0.00 1.00 1.6	Dibenzo (1,2,5 - cu) Fyrere Dibenzo (a,h) anthracene Barzo (a h i) papilana	0.00	0.0	Eliberto (1,2,5 - cu) ryrele Eliberto (a,h) anthracene Barto (c.h.) randere	0.0	0.0
	02077	20777			1014		10005	00007
I OIAI ALOIIIAIICS	14012	14131	I OLAL AL OFFICIES		11241	I OLAL AL OILIALICS	67001	13362
% Surrogate Recovery	0.08	100	% Surrogate Recovery	00 0	100	% Surrogate Recovery	000	100
5 /lpna /ndroetane Phonanthrono d-10	0.08 0.08	1.00	5 / upna / undroetane Phonanthrono d-10	96:0	1.00	 Aupril Ancrostane Phonanthrond d-10 	1.00	1.00

STATISTIC	CONTRO S FOR SURROGA Testing Date	TE CORRECTED DA	ТА			
Alkane Analyte:	Replicate #1 Conc (ng/mg)	Replicate #2 Conc (ng/mg)	Replicate #3 Conc (ng/mg)	Average	Stdev	%RSD
nC-10 Decane	1400	1400	1400	1400	0.00	0.00
nC-11 Undecane	2400	2400	2200	2333	115	4.9
nC-12 Dodecane	2700	2700	2300	2567	231	9.0
nC-13 Tridecane	2500	2700	2400	2533	153	6.0
nC-14 Tetradecane	2500	2700	2300	2500	200	8.0
nC-15 Pentadecane	2400	2600	2300	2433	153	6.3
nC-16 Hexadecane	2300	2400	2000	2433	208	9.3
nC-17 Heptadecane	2300	2400	2100	2267	153	6.7
Pristane	1500	1600	1500	1533	58	3.8
nC-18 Octadecane	2200	2600	2400	2400	200	8.3
Phytane	1400	1500	1300	1400	100	7.1
nC-19 Nonadecane	2000	2300	2000	2100	173	8.2
nC-20 Eicosane	2100	2400	2000	2167	208	9.6
nC-21 Heneicosane	2000	2200	1900	2033	153	7.5
nC-22 Docosane	1700	1900	1700	1767	115	6.5
nC-23 Tricosane	1500	1700	1500	1567	115	7.4
nC-24 Tetracosane	1400	1500	1500	1467	58	3.9
nC-25 Pentacosane	2100	2300	1900	2100	200	9.5
nC-26 Hexacosane	2100	2300	1900	2100	200	9.5
nC-27 Heptacosane	1500	1600	1400	1500	100	6.7
nC-28 Octacosane	1200	1200	1200	1200	0.00	0.00
		1200		1200		
nC-29 Nonacosane	1200		1200		0.00	0.00
nC-30 Triacontane	1200	1300	1200	1233	58	4.7
nC-31 Hentriacontane	1100	1200	1 100	1133	58	5.1
nC-32 Dotriacontane	900	1000	930	943	51	5.4
nC-33 Tritriacontane	970	1000	880	950	62	6.6
nC-34 Tetratriacontane	760	800	750	770	26	3.4
nC-35 Pentatriacontane	1100	1200	1100	1133	58	5.1
Total Alkanes	48430	52100	46360	48963	2907	5.9
Aromatic Analyte:	Conc (ng/mg)	Conc (ng/mg)	Conc (ng/mg)	Average	Stdev	%RSD
Naphthalene	610	640	550	600	46	7.6
C1-Naphthalenes	1700	1700	1500	1633	115	7.1
C2-Naphthalenes	2200	2400	2000	2200	200	9.1
C3-Naphthalenes	1800	1900	1600	1767	153	8.6
C4-Naphthalenes	810	900	820	843	49	5.8
Fluorene	110	120	100	110	10	9.1
C1-Fluorenes	270	290	260	273	15	5.6
C2-Fluorenes	370	420	350	380	36	9.5
C3- Fluorenes	340	370	330	347	21	6.0
Dibenzothiophene	250	280	240	257	21	8.1
C1-Dibenzothiophenes	460	510	420	463	45	9.7
C2-Dibenzothiophenes	560	610	500	557	55	9,9
C3- Dibenzothiophenes	360	410	340	370	36	9.7
Phenanthrene	320	330	300	317	15	4.8
C1-Phenanthrenes	550	600	530	560	36	6.5
C2-Phenanthrenes	630	710	590	643	61	9.5
C2-Phenanthrenes	440	500	420	453	61 42	9.5 9.2
C4-Phenanthrenes	140	130	130	133	5.8	4.3
Anthracene	2.8	2.5	2.4	2.6	0.19	0.00
Fluoranthene	3.1	3.0	3.0	3.0	0.04	1.4
Pyrene 01 Durana	8.5	8.9	8.9	9	0.23	2.6
C1- Pyrenes	110	110	110	110	0.00	0.00
C2- Pyrenes	150	160	140	150	10	6.7
C3- Pyrenes	160	170	150	160	10	6.3
C4- Pyrenes	96	100	88	95	6 .1	6.5
Naphthobenzothiophene	56	60	50	56	5.0	9.0
C-1 Naphthobenzcthiophenes	200	210	180	197	15	7.8
C-2 Naphthobenzcthiophenes	190	210	180	193	15	7.9
C-3 Naphthobenzothlophenes	160	150	140	150	10	6.7
Benzo (a) Anthracene	8.7	9.0	8.1	8.6	0.42	4.9
Chrysene	61	66	57	61	4.7	7.6
C1- Chrysenes	92	95	94	94	1.5	1.6
C2- Chrysenes	120	120	110	117	5.8	4.9
C3- Chrysenes	90	92	83	88	4.9	5.6
	51	47	48	49	1.8	3.8
C4- Chrysenes						3.8 7.1
Benzo (b) Fluoranthene	4.0	3.8	3.5	3.8	0.27	
Benzo (k) Fluoranthene	3.3	3.6	3.4	3.4	0.18	5.3
Benzo (e) Pyrene	8.2	8.9	7.6	8	0.67	8.2
Benzo (a) Pyrene	1.2	1.2	1.1	1.2	0.05	4.7
Perylene	0.69	0.67	0.65	0.7	0.02	3.1
	0.00	0.00	0.00	0.00	0.00	0.00
Indeno (1,2,3 - cd) Pyrene						
	0.00	0.0	0.00	0.00	0.00	0.00
Indeno (1,2,3 - cd) Pyrene				0.00 1.8		

STATISTIC	NUTRIE S FOR SURROG Testing Date	ATE CORRECTED DA	ATA			
Alkane Analyte:	Replicate #1 Conc (ng/mg)	Replicate #2 Conc (ng/mg)	Replicate #3 Conc (ng/mg)	Average	Stdev	%RSD
nC-10 Decane	730	710	730	723	12	1.6
nC-11 Undecane	1800	1700	1800	1767	58	3.3
nC-12 Dodecane	2100	1900	2000	2000	100	5.0
nC-13 Tridecane	1800	1800	1900	1833	58	3.1
nC-14 Tetradecane	2200	2000	2200	2133	115	5.4
nC-15 Pentadecane	2000	1900	1900	1933	58	3.0
nC-16 Hexadecane	1800	1800	1800	1800	0.00	0.00
nC-17 Heptadecane	1900	1900	2000	1933	58	3.0
Pristane	1600	1700	1600	1633	58	3.5
nC-18 Octadecane	2100	2100	2400	2200	173	7.9
Phytane	1500	1600	1600	1567	58	3.7
nC-19 Nonadecane	1800	1800	1900	1833	58	3.1
nC-20 Eicosane	2000	2000	2000	2000	0.00	0.00
nC-21 Heneicosane	1800	1800	2000	1867	115	6.2
nC-22 Docosane	1600	1700	1800	1700	100	5.9
nC-23 Tricosane	1600	1600	1800	1667	115	6.9
nC-24 Tetracosane	1600	1800	1800	1733	115	6.7
nC-25 Pentacosane	2100	2200	2300	2200	100	4.5
nC-26 Hexacosane	2200	2200	2300	2233	58	2.6
nC-27 Heptacosane	1900	2000	2100	2000	100	5.0
nC-28 Octacosane	1700	1900	1700	1767	115	6.5
nC-29 Nonacosane	1800	1800	1700	1767	58	3.3
nC-30 Triacontane	1200	1200	1200	1200	0.00	0.00
nC-31 Hentriacontane	1200	1200	1300	1200	58	4.7
					0.00	0.00
nC-32 Dotriacontane	1100	1100	1100	1100		
nC-33 Tritriacontane	1100	1000	1200	1100	100	9.1
nC-34 Tetratriacontane	980	980	1100	1020	69	6.8
nC-35 Pentatriacontane Total Alkanes	1000 46210	960 46350	1100 48330	1020 46963	72 1186	7.1 2.5
Aromatic Analyte:	Conc (ng/mg)	Conc (ng/mg)	Conc (ng/mg)	Average	Stdev	%RSD
	440	480	510	477	35	7.4
Naphthalene	1600	1800		1700	100	5.9
C1-Naphthalenes			1700			
C2-Naphthalenes	2200	2400	2300	2300	100	4.3
C3-Naphthalenes	1800	1900	1900	1867	58	3.1
C4-Naphthalenes	860	940	840	880	53	6.0
Fluorene	110	110	110	110	0.42	0.38
C1-Fluorenes	250	250	250	250	0.16	0.07
C2-Fluorenes	360	370	360	363	5.5	1.5
C3- Fluorenes	330	330	330	330	0.22	0.07
Dibenzothiophene	240	250	250	247	5.7	2.3
C1-Dibenzothiophenes	450	470	450	457	12	2.5
C2-Dibenzothiophenes	560	590	570	573	15	2.7
C3- Dibenzothiophenes	380	360	370	370	10.2	2.8
Phenanthrene	320	330	330	327	5.7	1.7
C1-Phenanthrenes	590	600	600	597	5.8	1.0
C2-Phenanthrenes	620	680	670	657	32	4.9
C3-Phenanthrenes	420	470	470	453	29	6.4
C4-Phenanthrenes	130	130	140	133	5.7	4.2
Anthracene	2.0	2.2	2.1	2.1	0.10	4.2
Fluoranthene	2.5	2.2	2.1	2.1	0.10	8.2
Pyrene C1 Byranas	5.3	5.9	6.1	5.8	0.44	7.7
C1- Pyrenes	110	120	120	116	5.9	5.1
C2- Pyrenes	140	150	150	146	5.6	3.8
C3- Pyrenes	140	150	160	150	10	6.5
C4- Pyrenes	84	87	91	87	3.8	4.3
Naphthobenzothiophene	55	57	56	56	1.3	2.3
C-1 Naphthobenzothiophenes	170	180	190	180	10	5.6
C-2 Naphthobenzothiophenes	180	190	200	190	10	5.1
C-3 Naphthobenzothlophenes	150	150	150	150	0.10	0.07
Benzo (a) Anthracene	14	14	13	14	0.50	3.7
Chrysene	83	85	89	86	3.3	3.9
C1- Chrysenes	120	130	140	130	10	7.8
C2- Chrysenes	150	160	170	160	10	6.1
C3- Chrysenes	130	110	130	123	12	9.4
C4- Chrysenes	74	70	75	73	2.5	3.5
						5.3
Benzo (b) Fluoranthene	2.9	2.7	2.7	2.8	0.15	
Benzo (k) Fluoranthene	3.1	3.4	2.9	3.2	0.25	8.0
Benzo (e) Pyrene	10	10	9	10	0.56	5.7
Benzo (a) Pyrene	1.3	1.4	1.3	1.3	0.05	3.6
Perylene	0.66	0.71	0.72	0.70	0.03	4.5
	0.00	0.00	0.00	0.00	0.00	0.00
Indeno (1,2,3 - cd) Pyrene						
Indeno (1,2,3 - cd) Pyrene Dibenzo (a,h) anthracene	0.00	0.00	0.00	0.00	0.00	0.00
		0.00 1.7	0.00 1.6	0.00 1.7	0.00 0.07	0.00 4.0

Alkane Analyte: nC-10 Decane nC-11 Undecane nC-12 Dodecane nC-13 Tridecane nC-14 Tetradecane nC-15 Pentadecane nC-16 Hexadecane nC-17 Heptadecane	Replicate #1 Conc (ng/mg) 1100 2100	Replicate #2 Conc (ng/mg)	Replicate #3	-		
nC-11 Undecane nC-12 Dodecane nC-13 Tridecane nC-14 Tetradecane nC-15 Pentadecane nC-16 Hexadecane nC-17 Heptadecane	2100	1	Conc (ng/mg)	Average	Stdev	%RSD
nC-12 Dodecane nC-13 Tridecane nC-14 Tetradecane nC-15 Pentadecane nC-16 Hexadecane nC-17 Heptadecane		1200	1200	1167	58	4.9
nC-13 Tridecane nC-14 Tetradecane nC-15 Pentadecane nC-16 Hexadecane nC-17 Heptadecane	2100	2100	2100	2100	0.00	0.00
nC-14 Tetradecane nC-15 Pentadecane nC-16 Hexadecane nC-17 Heptadecane	2100	2400	2400	2300	173	7.5
nC-14 Tetradecane nC-15 Pentadecane nC-16 Hexadecane nC-17 Heptadecane	2000	2200	2300	2167	153	7.1
nC-15 Pentadecane nC-16 Hexadecane nC-17 Heptadecane	2200	2500	2400	2367	153	6.5
nC-16 Hexadecane nC-17 Heptadecane	2100	2400	2400	2300	173	7.5
nC-17 Heptadecane	2300	2600	2500	2467	153	6.2
	2400	2600	2400	2467	115	4.7
Pristane	1700	1800	1700	1733	58	3.3
nC-18 Octadecane	2200	2500	2500	2400	173	7.2
Phytane	1600	1700	1600	1633	58	3.5
nC-19 Nonadecane	1900	2200	2200	2100	173	8.2
nC-20 Eicosane	2200	2400	2500	2367	153	6.5
nC-21 Heneicosane	2000	2300	2200	2167	153	7.1
nC-22 Docosane	1900	2100	2000	2000	100	5.0
nC-23 Tricosane	1800	2000	2000	1933	115	6.0
nC-24 Tetracosane	2000	2000	1900	1933	58	2.9
nC-25 Pentacosane	2000	2400	2300	2300	100	4.3
nC-25 Pentacosane nC-26 Hexacosane	2200	2400	2300	2300	100	4.3
nC-26 Hexacosane	1900	2500	1900	2300	1/3	7.5 5.9
nC-27 Heptacosane	1900	1900	1900	1967	115	5.9 6.5
nC-28 Octacosane nC-29 Nonacosane	1800	1900	1700	1/6/	115	5.6
nC-30 Triacontane	1100	1100	1100 1100	1100	0.00 115	0.00 9.9
nC-31 Hentriacontane	1100	1300		1167		
nC-32 Dotriacontane	1000	1100	1100	1067	58	5.4
nC-33 Tritriacontane	990	930	930	950	35	3.6
nC-34 Tetratriacontane nC-35 Pentatriacontane	1100 940	1100 960	1000 940	1067 947	58 11	5.4 1.2
Total Alkanes	49630	54290	52270	52063	2337	4.5
Aromatic Analyte:	Conc (ng/mg)	Conc (ng/mg)	Conc (ng/mg)	Average	Stdev	%RSD
Naphthalene	580	580	570	577	5.8	1.0
C1-Naphthalenes	1700	1800	1600	1700	100	5.9
C2-Naphthalenes	2300	2400	2100	2267	153	6.7
C3-Naphthalenes	1900	1900	1700	1833	115	6.3
C4-Naphthalenes	810	860	790	820	36	4.4
Fluorene	110	110	110	110	0.06	0.05
C1-Fluorenes	260	280	260	267	12	4.4
C2-Fluorenes	370	420	360	383	32	8.4
C3- Fluorenes	370	390	340	367	25	6.9
Dibenzothiophene	280	290	260	277	15	5.6
C1-Dibenzothiophenes	490	520	490	500	17	3.5
C2-Dibenzothiophenes	610	670	590	623	42	6.7
C3- Dibenzothiophenes	390	410	350	383	31	8.0
Phenanthrene	360	390	330	360	30	8.4
C1-Phenanthrenes	650	710	610	657	50	7.7
C2-Phenanthrenes	760	810	700	757	55	7.3
C3-Phenanthrenes	490	520	450	487	35	7.3
C4-Phenanthrenes	150	150	150	150	0.15	0.10
Anthracene	2.1	1.9	2.0	2.0	0.12	5.8
Fluoranthene	2.4	2.5	2.5	2.5	0.07	2.8
Pyrene	7.2	7.1	6.3	6.9	0.48	7.0
C1- Pyrenes	110	110	120	113	5.7	5.0
C2- Pyrenes	140	150	130	140	10	7.2
C3- Pyrenes	140	140	130	137	5.8	4.3
C4- Pyrenes	77	82	70	76	<u>6.0</u>	7.9
Naphthobenzothiophene	53	60	53	55	3.7	6.6
C-1 Naphthobenzcthiophenes	180	200	180	187	12	6.2
C-2 Naphthobenzothiophenes	190	200	170	187	15	8.2
C-3 Naphthobenzothlophenes	150	160	150	153	5.8	3.8
Benzo (a) Anthracene	15	15	14	15	1.0	6.9
Chrysene	78	83	76	79	3.6	4.5
C1- Chrysenes	120	130	120	123	5.8	4.7
C2- Chrysenes	140	150	140	143	5.8	4.1
C3- Chrysenes	120	110	110	113	5.8	5.1
C4- Chrysenes	68	75	69	70	3.5	4.9
Benzo (b) Fluoranthene	3.9	3.9	3.4	3.7	0.25	6.6
Benzc (k) Fluoranthene	4.0	4.3	3.6	4.0	0.32	8.1
Benzo (e) Pyrene	12	13	12	12	0.72	5.7
Benzo <mark>(</mark> a) Pyrene	1.5	1.5	1.3	1.4	0.06	4.5
Perylene	0.62	0.63	0.60	0.62	0.02	2.9
Indeno (1,2,3 - cd) Pyrene	0.00	0.00	0.00	0.00	0.00	0.00
Dibenzo (a,h) anthracene Banzo (a h i) port/and	0.00	0.00	0.00	0.00	0.00	0.00
Benzo (q.h.i) pervlene Total Aromatics	1.6 14197	1.6 14911	1.5 13322	1.6 14143	0.07 796	4.4 5.6

CON Te Initia Fxi	CONTROL, REPLICATE 1 Testing Date: Dsy 7 Initial Oil Weight: 500 mg Final Extracted Volume: 10 ml	Dia la constante de la constante de La constante de la constante de	Einal C	CONTROL, REFLICATE 2 Testing Date: Day 7 Initial Oil Weight: 540 mg Final Extracted Volume: 10 ml	0 0	Einal F	CONTROL, REPLICATE 3 Testing Date: Day 7 Initial Oil Weight: 500 mg Final Extracted Volume: 10 ml	E G
Alkane Analyte:	Concentration (ng/mg)	Surrogate Corrected (ng/mg)	Alkane Analyte:	Concentration (ng/mg)	Surmgate Corrected (ng/mg)	Alkane Analyte:	Concentration (ng/mg)	Surrogate Corrected (ng/mg)
nC-10 Decare	440	490	nC-10 Decane	440	480	nC-10 Decane	4/0	510
nC-11 Undecare	1600	1800	nC-11 Undecane	1800	2000	nC-11 Undecane	1600	1700
nC-12 Uotecare	2000	2200	nC-12 DodeCane nC-13 Tridecane	0002	2300	nC-12 Uodecane nC-13 Tridecane	2000	0022
nC-14 Tetradecane	1800	2000	nC-14 Tetradecane	1900	2100	nC-14 Tetradecane	2000	2200
n0-15 Pentadecane	2200	2400	nC-15 Pentadecane	2100	2000	nC-15 Pentadecane	2000	2200
nC-16 Heotadecane	1900	2400	nC-16 Hexadecane	2100	2300	nC-10 Hexadecane nC-17 Heptadecane	2000	2200
Pristane 6 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1	1400	1600	Pristane	1400	1500	Pristane	1400	1500
nC-18 Octadecane	2200	2400	nC-18 Octadecane	2100	2300	nC-18 Octadecane	2100	2300
€.	1200	1300	Phytane	1300	1400	Phytsne	1200	1300
IIC-19 Nurladecane	1900	2100	IIC-19 Nunadecane	1900	2100	IIC-19 Norradecare	1800	2000
nozo ciudane nº21 Hanaimeana	1900	2100	no-zo Ecoadio	1000	0012	nr. 21 Haneloosne	1800	0000
nC-22 Docosare	1700	1900	nc-22 Docosane	1600	1800	nC-22 Decosane	1600	1700
nC-23 Tricosare	1600	1800	nC-23 Tricosane	1500	1700	nC-23 Tricosane	1400	1500
nC-24 Tetracosane	1500	1700	nC-24 Tetracosane	1500	1700	nC-24 Tetracosane	1400	1500
nC 25 Pentacoeane	1900	2100	nC 25 Pentacosane	2000	2200	nC 25 Pentacocane	2000	2200
nC26 Hexannsane	1800	2000	nC-26 Hexacosane	2000	00000	nC26 Hexacosane	2000	00000
nc-27 Heptacosane	1400	1600	nC-27 Heptacosane	1500	1700	nC-27 Heptacosane	1400	1500
nC-28 Octacosane	1200	1300	nC-28 Octacosane	1300	1400	nC-28 Octacosane	1300	1400
nC-29 Nonacosane	1100	1200	nC-29 Nonacosane	1300	1400	nC-29 Nonacosane	1200	1300
IIC-30 Triacontane	1100	1200	nC-30 Triacontane	1100	1200	nc-30 Triacontane	1100	1200
nu-31 Hentriacontane	750	0011	nC-31 Hentriacontane	0011	0071	nc-31 hentriacontane	0011	0071
nC-33 Tritriacontane	068	066	nC-33 Tritriacontane	1000	1100	nC-33 Tritriacontane	1000	1100
nC-34 letratriacontane	066	1100	nC-34 letratriacontane	910	1000	nC-34 Lefratriacontane	066	1100
nC-35 Pentatriacontane	1200	1300	nC-35 Pentatriacontare	1300	1400	nC-35 Pertatriacontane	1200	1300
Total Albanee	02041	17540	Total Alkanee	01110	48530	Total Alkanee	12870	16700
Aromatic Analyte:	0/674	010.04	Aromatic Analyte:		07004	Aromatic Analyte:	0/074	06/04
Naphthalere	300	320	Naprthalene	087	067	Naphthalene	087	300
C1-Naphthalenes	950	1000	C1-Naphthalenes	1000	1000	C1-Naphthalenes	1100	1200
C2-Naphthalenes	1800	1900	C2-Naphthalenes	1700	1800	C2-Naphthalenes	1700	1800
C.SNaphthalenes	740	700	C.3-Raphthalenes	0071	0051 0020	C 3-Naphinalenes	0071	1300
Flucience	97	100	Fluorere		100	Flurieres	110	120
C1-Fluorenes	250	270	C1-Fluorenes		250	C1-Fluorenes	250	270
C2-Fluorenes	310	330	C2-Fluorenes	300	310	C2-Fluorenes	270	290
C3- Fluorenes	290	310	C3- Fluorenes	310	320	C3- Fluorenes	290	310
C1 Diherrothinshenen	410	740	C1 Dihenzothinhenen	360	240	C1 Dihanzothichhanae	360	240
C2-Dibenzothiophenes	550	280	C2-Dibenzothiophenes	490	510	C2-Dibenzothicphenes	460	490
C3- Dibenzothiophenes	360	360	C3- Dibenzothiophenes	310	320	C3- Dibenzothicphenes	300	320
Phenanthrere	330	350	Phenanthrene	290	300	Phenanthrene	280	300
C2-Phenanthrenes	660 660	200	C2-Prenanthrenes	200 630	980	C2-Phenarthrenes	550	065
C3-Phenanthrenes	380	410	C3-Phenanthrenes	370	390	C3-Phenarthrenes	340	370
C4-Phonanthronca	130	140	C4-Phonanthronca	130	140	C4-Phonarthronca	120	130
Anthracere	1.5	1.6	Anthracene	1.6	1.6	Anthracene	1.6	17
riuoranthere	2. ¢	9: 2	ruoranmene		4 C 0	Linoranthene	2.0	6.0
C1- Pyrenes	120	130	C1- Pvrenes	110	120	C1- Pvrenes	100	110
C.2- Eyrenes	160	170	C.2- Pyrenes	130	140	C2- Pyrenes	130	140
C3- Fyrenes	130 1	140	C3- Pyrenes	110	120	C3- Pyrenes	120	130
Nachthohanzothiochana	2 6	2 2	Vanhthohanzothipohara	3 4	00	Nanhthohanzothinnhara Manhthohanzothinnhara	10	27
C-1 Naphthobenzothiophenes	170	180	C-1 Naphthobenzothiophenes	150	160	C-1 Naphthobenzothiophenes	140	150
C-2 Naphthobenzothiophenes	190	200	C-2 Naphthobenzothiophenes	170	180	C-2 Naphthobenzothiophenes	160	170
C-3 Naphthobenzothiophenes	130	140	C-3 Naphthobenzothiophenes	110	120	C-3 Naphthobenzothiophenes	110	120
Chrysene	2 09	- 75	Chrysene	5°0	25	Chrysene	51	22
C1- Chrysonos	8	8	CI- Chrysenes	8	8	C1- Chrysones	62	: 3
C2- Chrysenes	110	120	C2- Chrysenes	100	100	C2- Chrysenes	94	100
C4- Chrysenes	90	\$ \$	CJ- Chrysenes C4- Chrysenes	8 8	40	C4- Chrysenes	47	8 5
Benzo (b) Fluoranthene	1.9	20	Benzo (b) Fluorenthere	3.6	1	Benzo (b) Fluoranthene	5	21
Benzo (k) Fluoranthene	4.2	4.5	Benzo (k) Fluorsnthene	3.9	4.1	Benzo (k) Fluoranthene	3.6	4.2
Benzo (e) Pyrene	13	14	Benzo (e) Pyrene	12	13	Benzo (e) Pyrene	2 ;	13
Perviene	2	12	Perviene	01	1 🗆	Perviene	91	1 1
Inceno (1,2,3 - cd) Pyrene	00.0	0.00	Indeno (1,2,3 - cd) Pyrere	0.00	0.00	Indeno (1,2,3 - cd) Pyrene	00:0	0.00
Dibenzo (a,h) anthracene	0:00	0:00	Dibenzo (a.h) anthracere	0.00	0:00	Eibenzo (a,h) anthracene	0:00	0:00
				-				A.1
Total Aromatics	11171	11901	Total Aromatics	10420	10931	Total Aromatics	10192	10957
% Surrogata Recovery			% Surrogate Recovery			% Surrogate Recovery		
5 Alpha Androstane	0:00	1.00	5 Alpha Androstane	0.91	1.00	5 Alpha Androstane	0.92	1.00
Phenanthrene c-10	0.94	1.00	Phenanthrene d-10		1.00	Phenanthrene d-10	0.93	1.00

	NUTRIENT, REFLICATE 1 Testing Date: Dsy 7 Initial Oil Weight: 510 mg	5		NUTRIENT, REPLICATE 2 Testing Date: Day 7 Initial Oil Weight: 510 mg	6		NUTRIENT, REPLICATE 3 Testing Date: Day 7 Initial Oil Weight: 530 mg	
Final L	Final Extracted Volume: 10 mL	0 mL	Final	Final Extracted Volume: 10 mL	0 mL	Final E	Final Extracted Volume: 10 mL) mL
Altere Andrew	Cancentration (ng/mg)	Surrogate Corrected (rg/mg)	Allower Sectored	Concentration (ng/mg)	Surmgate Corrected (ag/mg)	Allower Assertation	Concentration (ng/mg)	Surrogate Corrected (ng/mg)
Alkane Analyte: nC-10 Decare	540	980	Alkane Analyte: nC-10 Decane	540	630	Alkane Analyte: nC-10 Decane	540	280
nC-11 Undecare	710	900	nC-11 Undecane	730	850	nC-11 Undecane	830	910 1200
n0-13 Tritecare	1100	1400	n0-13 Tritecane	1300	1500	n0-13 Tridecane	1200	1300
nC-14 Tetradecane	1200	1500	nC-14 Tetradecane	1200	1400	nC-14 Tetradecane	1200	1300
nC-16 Hexadecane	1400	1800	nC-16 Hexadecane	1500	1800	nC-16 Hexadecane	1500	1600
nC-17 Heptadecane	1300	1700	nC-17 Heptadecane	1400	1600 1800	nC-17 Heptadecane	1400 1500	1500
nC-18 Octadecane	1300	1700	nC-18 Octadecane	1300	1500	nC-18 Octadecane	1400	1500
Phytane Control of Con	1200	1500	Phytane P	1200	1400	Phytsne	1300	1400
IIC-19 NUIBIDECAILE	1300	1200	IIC-19 NUIIAUECAIRE	1300	1500	IIC-15 NUIBUECAILE	1500	1600
nC-21 Heneicosane	1200	1500	nC-21 Heneicosane	1200	1400	nC-21 Heneicosane	1300	1400
nC-22 Docosare	1200	1500	nC-22 Docosane	1400	1600	nC-22 Decosane	1300	1400
nC-23 Incosare	1200	1500	nC-23 Incosane	1200	1400	nC-23 Incosane	1200	1300
nC-25 Pentacosane	2000	2500	nC-25 Pentacosane	2100	2500	nC 25 Pentacosane	2200	2400
nC.26 Hexachsane	1900	2400	nC-26 Hexacosane	2200	2600	nC-26 Hexacnsane	2000	2200
nc-27 Heptacosane	1600	2000	nC-27 Heptacosane	1800	Z100	nC-27 Heptacosane	1300	2000
nC-26 Uctacosane	1100	1400	nC-28 Octacosane	1400	1500	nC-28 Octacosane	1300	1400
IIC-30 Triacontaine	800	1000	IIC-30 Triacontane	870	1000	InC-30 Triacontane	000	1100
nC-31 Hentriacontane	066	1300	nC-31 Hentriacontane	1000	1200	nC-31 Hentriacontane	1100	1200
nC-32 Botriacontano	000	1100	nC-32 Dotriacontanc	1000	1200	nC-32 Dotriacontano	000	1100
nC-34 letratriacontane	810	1001	nC-34 Tetratmacontare	080 N80	0001	nC-34 Tetratriacontane	810	060
nC-35 Pentatriacontane	890	1100	nC-35 Pentatriacontare	930	1100	nC-35 Pertatriacontane	1100	1200
Total Alkanes	32880	41680	Total Alkanes	35440	41500	Total Alkanae	36279	39380
Aromatic Analyte:			Aromatic Analyte:			Aromatic Analyte:		
Naphthaere	290	350	Naprthalene	350	3/0	Naphthalene	320	330
C1-Naphthalenes C2-Naphthalenes	1500	1300	C1-Naphthalenes	001	0002	C1-Naphmalenes C2-Manhthalenes	900 1800	1900
C3-Naphthalenes	1100	1300	C3-Naphthalenes	1400	1500	C3-Naphthalenes	1400	1500
C4-Naphthalenes	650	790	C4-Naphthalenes	760	810	C4-Naphthalenes	760	790
C1-Fluorenes	180	220	C1-Fluorenes	210	230	C1-Fluorenes	210	220
C2-Fluorenes	220	270	C2-Fluorenes	250	270	C2-Fluorenes	260	270
C3- Fluorenes	230	280	C3- Fluorenes	260	280	C3- Fluorenes	290	300
Diborzothiaphono C1 Dibenzothiashenes	180 320	300	Dibonzothiphono C1 Dibenzothiophenes	340	210 360	Dibonzothiophono C1 Dibenzothiophenee	210 380	380
C2-Dibenzothiophenes	400	490	C2-Dibenzothiophenes	430	460	C2-Dibenzothicphenes	460	480
C3- Dibenzothiophenes	260	320	C3- Dibenzothiophenes	270	290	C3- Dibenzothicphenes	310	320
C1-Phenanthrenes	480	290	C1-Phenanthrenes	480	510	C1-Phenanthrenes	280 520	540
C2-Phenanthrenes	500	610	C2-Phenan:hrenes		630	C2-Phenarthrenes	600	620
C3-Phenanthrenes	290	350	C3-Phenanthrenes		370	C3-Phenarthrenes	370	38
Anthracere	0.93	1	Anthracene	1.0	11	Anthracene	1.0	1.0
-luoranthere	2.1	2.5	Fluoranthene	2.0	21	Fluoranthene	2.5	2.4
C1- Pyrenes	100	120	C1- Pvrenes	110	120	C1-Pvrenes	11	110
C.2- Pyrenes	110	130	C2- Pyrenes	130	140	C.2- Pyrenes	130	130
C3- Fyrenes	60	120	C3- Pyrenes	£ 5	120	C3- Pyrenes	1 5	110
Naphthobenzothiophene	40	49	Naphthobenzothipphere	45	48	Naphthobenzothiophene	48	50
C-1 Naphthobenzothiophenes	130	160	C-1 Naphthobenzothiophenes	140	150	C-1 Naphthobenzothiophenes	140	150
C-2 Naphthoben Zothiophenes C-3 Naphthoben Zothiophenes	98	1/0	C-2 Naphthobenzothiophenes C-3 Naphthobenzothiophenes	110	1/0	C-2 Naphtrobenzothiophenes C-3 Naphtrobenzothiophenes	110	19 10
Benzo (a) Anthracene	5	ţ	Benzo (a) Anthracene	÷	12	Benzo (a) Anthracene	5	13
Chrysene	65	62	Chrysene	02	75	Chrysene	73	75
C2- Chrysenes	130	160	C2- Chrysenes	130	140	CI- Chrysenes	140	140
C3- Chrysenes	87	110	C3- Chrysenes	88	95	C3- Chrysenes	100	100
C4- Chrysenes	54 10	99 ;	C4- Chrysenes	99 (49 Ç	C4- Chrysenes	57	59
Benzo (k) Fluoranthene	5.4	5.9	Benzo (k) Fluoranthere	4 6i	3.0	Benzo (k) Fluoranthene	28	- ci
Benzo (e) Pyrene	10	12	Benzo (e) Pyrene	; 12	£ ;	Benzo (e) Pyrene	12	¢;
Perviene Perviene	0.31	0.38	Perviene Perviene	0.40	0.43	Perviene Perviene	0.39	0.40
Incieno (1,2,3 - cd) Pyrene	0.00	0.00	Indeno (1,2,3 - cd) Pyrene	0.00	0.00	Indeno (1,2,3 - cd) Pyrene	0.00	0.00
Dibenze (a,h) anthracene Denze (a,h.i) perviene	0.00	0.00	Dibenzo (a.h) anthracere Denzo (a.h.i) perviene	0.00	0.00	Elbenzo (a,h) anthracene Denzo (c,h.i) perviene	0.00	0:00
	000			06207			07207	
I Utal Ar Utilatics		44011	I OLAL AL OFFICIES		0/#11	I Utal AL UIII atics	0101	
% Surrogata Recovery			% Surrogate Recovery		4.00	% Surrogate Recovery	500	0
5 Alpha Androstane Phenanthrene c-10	0.79 0.82	1.00	5 Alpha Androstane Phenanthrene d-10	0.86 0.93	1.00	5 Alpha Ancrostane Phenanthrene d-10	0.92	1.00
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sarva r Te: Initial Final Ext	SARVA PRODUCT, REFLICATE 1 Testing Date: Day 7 Initial UII Weight: 510 mg Final Extracted Volume: 10 mL	≣ 1 9 0mL	SARV In Final	SARVA PRODUCT, REFLICATE 2 Testing Date: Day 7 Initial UII weight: 520 mg Final Extracted Volume: 10 mL	re 2 9 0 mL	SARVA	SARVA PRODUCT, REPLICATE 3 Testing Date: Day 7 Initial UI Weight: 520 mg Final Extracted Volume: 10 mL	ге з 9 0 м.Г
				- - -				
Alkane Analyte:	(gm/gn) notretineano.	Surrogate Corrected (ng/mg)	Alkane Analyte:	tioncentration (ng/mg)	Surmgate Lorrected (ng/mg)	Alkane Analyte:	oncentration (ng/mg)	Surrogate Corrected (ng/mg)
n0-10 Decare	006	1100	nc-10 Decane	09P0	1000	nC-10 Decane	006	1000
nC-11 Undecare nC-12 Dodecare	1300	1600	nC-12 Dodecane	1400	1500	nC-11 Undecane nC-12 Dodecane	1200	1600
nC-13 Trifecare	1500	1800	nC-13 Triflecane	1600	1800	nC-13 Tridecane	1400	1600
nU-15 Pentadecane	1400	1700	nC-15 Pentadecane	1700	1900	nC-15 Pentadecane	1/00	1500
nC-16 Hexadecane	1500	1800	nC-16 Hexadecane	1800	2000	nC-16 Hexadecane	1600	1800
nu-17 neptagecane Pristane	1400	1700	nu-17 neptaueuaile Pristane	1500	1700	Pristane	1400	1600
nC-18 Octadecane	1500	1800	nC-18 Octadecane	1800	2000	nC-18 Octadecane	1500	1700
Inc-19 Nunadecare	1300	1600	Proytane IIC-19 Nunadecane	1600	0051	IIC-19 Nunadecane	1300	1500
IIC-20 Elcusare	1500	1800	IIC-20 Ecusalie	1700	1900	IIC-20 Elcusare	1500	1700
nC-21 Heneicosane nC-22 Docosare	1400	1700	nC-21 Heneicosane nC-22 Docosane	1600 1600	1800	nC-21 Heneicosane nC-22 Docosane	1400	1500
nC-23 Tricosare	1500	1800	nC-23 Tricosane	1700	1900	nC-23 Tricosane	1400	1600
nC-24 Tetracosane	1200	1500	nC-24 Tetracosane	1500	1700	nC-24 Tetracosane	1300	1500
nC-26 Hexansane	1500	1800	nC-26 Hexacosane	1600	1800	nC-26 Hexacosane	1600	1800
nc-27 Heptacosane	1400	1700	nc-27 Heptacosane	1600	1800	nC-27 Heptacosane	1400	1600
nC-28 Uctacosane nC-29 Nonacosane	1500	1/00	nC-28 Octacosane nC-29 Nonacosane	1600	1800	nC-28 Nonacosane	1400	1600
IIC-30 Triacontane	006	1100	IIC-30 Triacontane	850	940	IIC-30 Triacontane	850	026
nC-31 Hentriacontane	870	1100	nC-31 Hentriacontane	0980 140	950	nC-31 Hentriacontane	880	1000
nC-33 Tritriacontane	720	880	nC-33 Tritriacontane	790	870	nC-33 Tritriacontane	190	006
nC-34 Letratriacontane nC-35 Pentatriacontane	040 070	930 1200	nC-35 Pentatriacontare	86U 1100	950 1200	nC-34 Tetratriacontane nC-35 Pertatriacontane	/40 1000	840 1100
Total Alkanes	36240	44110	Total Alkanes	40150	44610	Total Alkanse Armetia Archite	36020	40990
Naphthaiere	270	330	Naphthalene	082	067	Naphthalene	067	330
C1-Naphthalenes	1100	1300	C1-Naphthalenes	1200	1200	C1-Naphthalenes	1100	1200
C2-Naphthalenes	1100	1300	C2-Naphthalenes C3-Naphthalenes	1400	1400	C2-Naphfhalenes	1200	1300
C4-Naphthalenes	610	740	C4-Naphthalenes	680	200	C4-Naphthalenes	590	660
C1-Fluorenes	500	240	C1-Fluorenes	210	2 22	C1-Fluorenes	8 <u>8</u>	200
C2-Fluorenes	240	290	C2-Fluorenes	260	270	C2-Fluorenes	220	250
C3- Fluorenes	220	270	C3- Fluorenes	250	260	C3- Fluorenes	240	270
UIDCFIZOTNIOPHONE C1 Dibenzothiophenes	200	350	Ulbonzothiophenee C1 Dibenzothiophenee	340	350	C1 Dibenzothicphenee	310	350
C2-Dibenzothiophenes	390	480	C2-Dibenzothiophenes	460	480	C2-Dibenzothicphenes	390	440
CJ- UIDENZOTNIOPINENES Phenanthrere	250	300	CJ- VIDEnZOTNOynenes Phenanthrene	270	280 280	C3- Ulbenzotnicpnenes Phenanthrene	240	270
C1-Phenanthrenes	460	560	C1-Phenanthrenes	530	550	C1-Phenarthrenes	440	490
C2-Phenanthrenes C3-Phenanthrenes	520 320	630 390	C2-Phenanthrenes C3-Phenanthrenes	360	590 370	C2-Phenarthrenes C3-Phenarthrenes	500 320	360
C4-Phonanthronca	110	130	C4-Phonanthronea	130	130	C4-Phonarthronoa	120	140
- Iuoranthere	1.F	2.3	Fluoranthene	2 2	22	Fluoranthene	25	2.4
Pyrene C4 Euroneu	7.2	0.0	Pyrene C1 Burnow	0.6	0.0	Pyrene C1 Bureauer	5.5	0.9
C.1- Pyrenes C.2- Pyrenes	110	130	C.2- Pyrenes	3/ 170	120	CI-Pyrenes C2-Pyrenes	110	120
C3- Fyrenes	88	110 65	C3- Pyrenes	110	110	C3- Pyrenes	8	00 2
Naphthobenzothiophene	98	47	Naphthobenzothipphere	48	49	Naphthobenzothiophene	41	46
C-1 Naphthobenzothiophenes C-2 Naphthobenzothiophenes	150	180	C-1 Naphthobenzothiophenes	150	160	C-1 Naphtrobenzothiophenes C-2 Naphtrobenzothiophenes	140	160
C-3 Naphthobenzothiophenes	100	120	C-3 Naphthobenzothiophenes	100	100	C-3 Naphthobenzothiophenes	56	110
Chrysene	2 99	80	Chrysene Chrysene	67	<u>*</u> 69	Chrysene	51	- 89
C1- Chrysonus	100	120	CI- Chrysenes	100	100	Ct- Chrysonus	100	110
C3- Chrysenes	84	145	C3- Chrysenes	021 03	130 96	C3- Chrysenes	78	140
C4- Chrysenes	50	61	C4- Chrysenes	56 1 E	57	C4- Chrysenes	57	64 1 c
Benzo (k) Fluoranthene	4 0; 4	23	Benzo (k) Fluoranthere	23	23	Benzo (k) Fluoranthene	50	23
Benzo (e) Pyrene Benzo (a) Pyrene	10	12	Benzo (e) Pyrene	11	11	Benzo (e) Pyrene	÷ -	12
Perylene	0.25	0.31	Perylene	0.27	0.28	Perylene	0.28	0.31
Inceno (1,2,3 - cd) Pyrene	0.0	0.00	Indeno (1,2,3 - cd) Pyrere	00:0	00:0	Indeno (1,2,3 - cd) Pyrene	0.00	0.0
Ubenzo (a.n) antin'acene Denzo (g.h.i) perytene	1.5	0.00 1.0	Ulbenzo (a.n) antimacere Denzo (g,h,i) perylene	1.7	1.0	Libenzo (a.n.) antinacene Denzo (g,h,i) perylene	1.7	1.9
Total Aromatics	9628	11626	Total Aromatics	10642	10949	Total Aromatics	9565	10657
% Surrocate Recovery			% Surrogate Recovery			Surrocate Recovery		
5 Alpha Androstane	0.81	1.00	5 Alpha And ostane	0.91	1.00	5 Alpha Ancrostane	0.88	1.00
Phenanthrene d-10	0.82	1.00	Phenanthrene d. 10		1.00	Phenanthrene d-10	0.89	1.00

nc-10 Decen 490 100 493 16 2 nc-11 Decen 1800 2000 2200 2233 66 2 nc-13 Tridecan 2200 2200 2200 2200 2200 2200 2200 2200 2200 2200 2200 2200 2200 2207 115 5 nc-14 Fertadecan 2400 2200 2200 2207 115 5 5 nc-16 Neadecan 2400 2200 2200 2233 58 2 115 5 5 nc-19 Nendecan 2100 2100 2000 2067 68 2 116 11 110 56 2 115 5 11	STATISTIC	CONTRO CS FOR SURROG Testing Date	ATE CORRECTED D	ТА			
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nC-24 Tetracosane 1700 1500 1630 165 2.7 56 2.2 nC-24 Featoceane 2000 2200 2200 2200 2167 56 2. nC-24 Heatoceane 1600 1700 1600 <td< td=""><td>nC-22 Docosane</td><td>1900</td><td>1800</td><td>1700</td><td>1800</td><td>100</td><td>5.6</td></td<>	nC-22 Docosane	1900	1800	1700	1800	100	5.6
nC-25 Pentacosane 2100 2200 2200 2133 115 5. nC-27 Heptacosane 1600 1700 1600 1600 100 160 nC-28 Nonacosane 1200 1400 1300 1300 100 17 nC-30 Triscontane 1200 1200 1200 1200 0.00 0.00 nC-31 Teristriacontane 1300 1400 1300 1667 56 4. nC-33 Tristriacontane 990 1100 1100 1667 56 4. 6.	nC-23 Tricosane	1800	1700	1500	1667	153	9.2
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C3- Fluorenes 310 320 310 313 5.8 11 Dibenzothiophenes 280 240 240 223 23 9 C1-Dibenzothiophenes 590 510 490 530 53 10 C2-Dibenzothiophenes 380 320 320 340 35 11 Phenenthrene 350 300 300 300 317 29 9, C1-Phenanthrene 660 590 550 600 56 80 C3-Phenanthrenes 700 660 590 650 56 80 C3-Phenanthrenes 140 140 130 137 5.8 42 Anthracene 1.6 1.6 1.7 1.6 0.09 0.0 Fluoranthene 3.6 3.4 3.5 3.5 0.7 2 Pyrene 11 9 9 10 0.82 8. C1-Pyrenes 170 140 140<							4.3
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C1-Dibenzothiophenes 440 380 380 400 35 8, C2-Dibenzothiophenes 590 510 490 530 53 11 C3-Dibenzothiophenes 380 320 320 340 35 11 Phenanthrenes 660 590 550 600 56 58 C2-Phenanthrenes 700 660 590 650 56 68 C3-Phenanthrenes 140 140 130 137 5.8 44 Anthracene 1.6 1.6 1.7 1.6 0.09 0.0 Fluoranthene 3.6 3.4 3.5 3.5 0.07 2. Pyrene 11 9 9 10 0.82 8. C1- Pyrenes 130 120 110 120 10 8. C2- Pyrenes 170 140 140 150 17 11 C3- Pyrenes 180 160 150 163	C3- Fluorenes	310	320	310	313		1.8
C2-Dibenzothiophenes 590 510 490 530 53 10 C3-Dibenzothiophenes 380 320 320 340 35 11 Phenanthrene 350 300 300 300 317 29 9, C1-Phenanthrenes 660 590 550 600 56 80 C2-Phenanthrenes 700 660 590 650 56 80 C4-Phenanthrenes 140 130 137 5.8 42 Anthracene 1.6 1.6 1.7 1.6 0.09 0.0 Fluoranthene 3.6 3.4 3.5 3.5 0.07 2. Pyrene 11 9 9 10 0.82 8. C1-Pyrenes 130 120 110 120 10 8. C2-Pyrenes 70 140 140 150 17 11 G3-Pyrenes 78 68 72 73	Dibenzothiophene	280	240	240	253	23	9.1
C3- Dibenzothiophenes 380 320 320 340 35 11 Phenanthrene 350 300 300 317 29 9 C1-Phenanthrenes 660 590 550 600 56 99 C2-Phenanthrenes 700 660 590 650 56 80 C3-Phenanthrenes 140 140 130 137 5.8 44 Anthracene 1.6 1.7 1.6 0.09 0.00 Fluoranthene 3.6 3.4 3.5 3.5 0.07 2.0 C1-Pyrenes 130 120 110 120 10 8.2 8. C2-Pyrenes 170 140 140 150 17 11 C3-Pyrenes 140 120 130 130 10 7. C4-Pyrenes 78 68 72 73 4.9 6. Naphthobenzothiophenes 160 150 163 15	C1-Dibenzothiophenes	440	380	380	400	35	8.7
Phenanthrene 350 300 300 317 29 9. C1-Phenanthrenes 660 590 550 600 56 9. C2-Phenanthrenes 700 660 590 650 56 8. C3-Phenanthrenes 140 130 137 5.8 4. Anthracene 1.6 1.6 1.7 1.6 0.09 0.7 Fluoranthene 3.6 3.4 3.5 3.5 0.07 2. Pyrene 11 9 9 10 0.82 8. C1- Pyrenes 130 120 110 120 10 8. C2- Pyrenes 170 140 140 150 17 17. C3- Pyrenes 78 68 72 73 4.9 6. C-1 Naphthobenzothiophenes 160 150 163 15 9. C-1 Asphthobenzothiophenes 100 100 100 100 10	C2-Dibenzothiophenes	590	510	490	530	53	10
Phenenthrene 350 300 300 317 29 9, C1-Phenanthrenes 660 590 550 600 56 9, C2-Phenanthrenes 700 660 590 650 56 8, C3-Phenanthrenes 140 130 137 5,8 4, Anthracene 1,6 1,6 1,7 1,6 0,09 0,2 Fluoranthene 3,6 3,4 3,5 3,5 0,07 2, Pyrene 11 9 9 10 0,82 8, C1- Pyrenes 130 100 110 120 10 8, C2- Pyrenes 170 140 140 150 17 17, C3- Pyrenes 78 68 72 73 4,9 6, C-1 Naphthobenzothiophenes 180 160 150 163 15 9, C-1 Naphthobenzothiophenes 100 100 100 100 <t< td=""><td>C3- Dibenzothiophenes</td><td>380</td><td>320</td><td>320</td><td>340</td><td>35</td><td>10</td></t<>	C3- Dibenzothiophenes	380	320	320	340	35	10
C2-Phenanthrenes 700 660 590 650 56 84 C3-Phenanthrenes 140 140 130 137 5.8 44 Anthracene 1.6 1.6 1.7 1.6 0.09 0.0 Fluoranthene 3.6 3.4 3.5 3.5 0.07 2. Pyrene 11 9 9 10 0.82 8. C1- Pyrenes 130 120 110 120 10 8. C2- Pyrenes 170 140 140 150 17 11 C3- Pyrenes 170 140 130 130 10 7. C3- Pyrenes 78 68 72 73 4.9 6. C-1 Naphthobenzothiophene 56 48 52 52 3.8 7. C-1 Naphthobenzothiophenes 100 150 163 15 9. C-1 Naphthobenzothiophenes 10.8 8.9 10.2 9.9		350	300	300	317	29	9.1
C2-Phenanthrenes 700 660 590 650 56 81 C3-Phenanthrenes 410 390 370 390 20 5. C4-Phenanthrenes 140 140 130 137 5.8 4.4 Anthracene 1.6 1.6 1.7 1.6 0.09 0.0 Fluoranthene 3.6 3.4 3.5 3.5 0.07 2. Pyrene 11 9 9 10 0.82 8. C1- Pyrenes 130 120 110 120 10 8. C2- Pyrenes 170 140 140 150 17 11 G3- Pyrenes 140 120 130 130 10 7. C3- Pyrenes 78 68 72 73 4.9 6. C-1 Naphthobenzothiophenes 160 150 163 15 9. C-1 Naphthobenzothiophenes 10.8 8.9 10.2 9.9 <	C1-Phenanthrenes	660	590	550	600	56	9.3
C3-Phenanthrenes 410 390 370 390 20 5. C4-Phenanthrenes 140 140 130 137 5.8 4. Anthracene 1.6 1.6 1.7 1.6 0.09 0.0 Fluoranthene 3.6 3.4 3.5 3.5 0.07 2. Pyrene 11 9 9 10 0.82 8. C1- Pyrenes 130 120 110 120 10 8. C2- Pyrenes 140 120 130 130 10 7. C3- Pyrenes 78 68 72 73 4.9 6. C-1 Naphthobenzothiophene 56 48 52 52 3.8 7. C-1 Naphthobenzothiophenes 100 150 163 15 9. C-1 Naphthobenzothiophenes 100 100 102 127 12 9. Benzo (a) Anthracene 10.8 8.9 10.2 9.9<							8.6
C4-Phenanthrenes 140 140 130 137 5.8 4.4 Anthracene 1.6 1.6 1.7 1.6 0.09 0.0 Fluoranthene 3.6 3.4 3.5 3.5 0.07 2. Pyrene 11 9 9 10 0.82 8. C1- Pyrenes 130 120 110 120 10 8. C2- Pyrenes 170 140 140 150 17 17. C3- Pyrenes 78 68 72 73 4.9 6. Naphthobenzothiophene 56 48 52 52 3.8 7. C-1 Naphthobenzothiophenes 180 160 150 163 15 9. C-2 Naphthobenzothiophenes 140 120 120 127 12 9. Berzo (a) Anthracene 10.8 8.9 10.2 9.9 0.98 9. C1- Chrysenes 96 90 84							5.1
Anthracene 1.6 1.6 1.7 1.6 0.09 0.0 Fluoranthene 3.6 3.4 3.5 3.5 0.07 2. Pyrene 11 9 9 10 0.82 8. C1-Pyrenes 130 120 110 120 10 8. C2-Pyrenes 170 140 140 150 17 12 C3-Pyrenes 140 120 130 130 10 7. C4-Pyrenes 78 68 72 73 4.9 6. Naphthobenzothiophene 56 48 52 52 3.8 7. C-1 Naphthobenzothiophenes 180 160 150 163 15 9. C-2 Naphthobenzothiophenes 10.8 8.9 10.2 9.9 0.98 9. C-3 Naphthobenzothiophenes 120 120 127 12 9. Benzo (a) Anthracene 0.8 90 84 90							4.2
Fluoranthene 3.6 3.4 3.5 3.5 0.07 2. Pyrene 11 9 9 10 0.82 8. C1- Pyrenes 130 120 110 120 10 8. C2- Pyrenes 170 140 140 150 17 17 C3- Pyrenes 78 68 72 73 4.9 6. C-1 Naphthobenzothiophene 56 48 52 52 3.8 7. C-1 Naphthobenzothiophenes 180 160 150 163 15 9. C-2 Naphthobenzothiophenes 200 180 170 183 15 8. C-3 Naphthobenzothiophenes 10.8 8.9 10.2 9.9 0.88 9. Benzo (a) Anthracene 10.8 8.9 10.2 9.9 0.58 6. C2- Chrysenes 96 90 84 90 5.8 6. C2- Chrysenes 120 100							0.00
Pyrene 11 9 9 10 0.82 8. C1- Pyrenes 130 120 110 120 10 8. C2- Pyrenes 170 140 140 150 17 17 C3- Pyrenes 140 120 130 130 10 7. C4- Pyrenes 78 68 72 73 4.9 6. C-1 Naphthobenzothiophene 56 48 52 52 3.8 7. C-2 Naphthobenzothiophenes 200 180 170 183 15 8. C-3 Naphthobenzothiophenes 200 180 170 183 15 8. C-3 Naphthobenzothiophenes 10.8 8.9 10.2 9.9 0.98 9. Benzo (a) Anthracene 10.8 8.9 10.2 9.9 0.98 9. C1- Chrysenes 96 90 84 90 5.8 6. C2- Chrysenes 120 100 <t< td=""><td></td><td></td><td></td><td></td><td></td><td></td><td>2.1</td></t<>							2.1
C1- Pyrenes 130 120 110 120 10 8. C2- Pyrenes 170 140 140 150 17 17 C3- Pyrenes 140 120 130 130 10 7. C4- Pyrenes 78 68 72 73 4.9 6. Naphthobenzothiophene 56 48 52 52 3.8 7. C-1 Naphthobenzothiophenes 200 180 170 183 15 9. C-3 Naphthobenzothiophenes 100 120 120 127 12 9. Benzo (a) Anthracene 10.8 8.9 10.2 9.9 0.98 9. C1- Chrysenes 64 57 55 59 4.7 8. C2- Chrysenes 120 100 100 107 12 11 C3- Chrysenes 64 62 66 64 2.1 3. C4- Chrysenes 43 40 50							
C2- Pyrenes 170 140 140 150 17 130 C3- Pyrenes 140 120 130 130 10 7 130 C4- Pyrenes 78 68 72 73 4.9 63 Naphthobenzothiophene 56 48 52 52 3.8 7 C-1 Naphthobenzothiophenes 180 160 150 163 15 9 C-2 Naphthobenzothiophenes 200 180 170 183 15 8 C-3 Naphthobenzothiophenes 140 120 120 127 12 9 Benzo (a) Anthracene 10.8 8.9 10.2 9.9 0.98 93 C1- Chrysenes 96 90 84 90 5.8 6. C2- Chrysenes 120 100 100 107 12 11 C3- Chrysenes 64 62 66 64 2.1 3. C4- Chrysenes 43 4							
C3- Pyrenes 140 120 130 130 10 7. C4- Pyrenes 78 68 72 73 4.9 6. Naphthobenzothiophene 56 48 52 52 3.8 7. C-1 Naphthobenzothiophenes 180 160 150 163 15 9. C-2 Naphthobenzothiophenes 200 180 170 183 16 8. C-3 Naphthobenzothiophenes 200 180 170 183 15 8. C-3 Naphthobenzothiophenes 10.0 120 127 12 9. 9.8 9. Garantinearen 10.8 8.9 10.2 9.9 0.88 9. 0.22 6.6 6.4 2.1 100 100 107 12 14. 14. 13 14. 14. 14. 15. 14. 14. 15. 15. 15. 14. 14. 15. 14. 14. 15. 14. 14.							8.3
C4- Pyrenes 78 68 72 73 4.9 64 Naphthobenzothiophene 56 48 52 52 3.8 7. C-1 Naphthobenzothiophenes 180 160 150 163 15 9. C-2 Naphthobenzothiophenes 200 180 170 183 15 8. C-3 Naphthobenzothiophenes 140 120 120 127 12 9. Benzo (a) Anthracene 10.8 8.9 10.2 9.9 0.98 9. C1- Chrysenes 96 90 84 90 5.8 6. C2- Chrysenes 120 100 100 107 12 11 C3- Chrysenes 120 100 100 107 12 11 C3- Chrysenes 43 40 50 44 50 11 Benzo (k) Fluoranthene 2.0 1.7 2.1 1.9 0.20 10 Benzo (c) Pyrene 1.7							
Naphthobenzothiophene 56 48 52 52 3.8 7.7 C-1 Naphthobenzothiophenes 180 160 150 163 15 9. C-2 Naphthobenzothiophenes 200 180 170 183 15 8. C-3 Naphthobenzothiophenes 140 120 120 127 12 9. Benzo (a) Anthracene 10.8 8.9 10.2 9.9 0.98 9. C1- Chrysene 64 57 55 59 4.7 8. C1- Chrysenes 96 90 84 90 5.8 6. C2- Chrysenes 120 100 100 107 12 11 C3- Chrysenes 64 62 66 64 2.1 3. C4- Chrysenes 43 40 50 44 5.0 11 Benzo (b) Fluoranthene 2.0 1.7 2.1 1.9 0.20 100 Benzo (a) Pyrene 1.7 <td< td=""><td></td><td></td><td></td><td></td><td></td><td>1</td><td></td></td<>						1	
C-1 Naphthobenzothiophenes 180 160 150 163 15 9; C-2 Naphthobenzothiophenes 200 180 170 183 15 8; C-3 Naphthobenzothiophenes 140 120 120 127 12 9; Benzo (a) Anthracene 10.8 8.9 10.2 9.9 0.98 9; C1- Chrysene 64 57 55 59 4.7 8; C1- Chrysenes 96 90 84 90 5.8 6; C2- Chrysenes 120 100 100 107 12 11 C3- Chrysenes 64 62 66 64 2.1 3; C4- Chrysenes 43 40 50 44 5.0 11 Benzo (b) Fluoranthene 4.5 4.1 4.2 4.3 0.18 4.2 Benzo (a) Pyrene 1.7 1.6 1.5 1.6 0.09 5.5 Benzo (a) Pyrene 1.7 <td< td=""><td></td><td></td><td></td><td></td><td></td><td></td><td>6.8</td></td<>							6.8
C-2 Naphthobenzothiophenes 200 180 170 183 15 8; C-3 Naphthobenzothiophenes 140 120 120 127 12 9; Benzo (a) Anthracene 10.8 8.9 10.2 9,9 0.98 9; Chrysene 64 57 55 59 4,7 8; C1- Chrysenes 96 90 84 90 5.8 6; C2- Chrysenes 120 100 100 107 12 1 G-3- Chrysenes 64 62 66 64 2,1 3; C4- Chrysenes 43 40 50 44 5.0 11 Benzo (b) Fluoranthene 2.0 1.7 2,1 1.9 0.20 10 Benzo (c) Privenes 14 13 13 0.82 6; Benzo (a) Pyrene 1.7 1.6 1.5 1.6 0.02 5; Benzo (a) Pyrene 1.7 1.6 1.5							7.3
C-3 Naphthobenzothiophenes 140 120 120 127 12 9. Benzo (a) Anthracene 10.8 8.9 10.2 9.9 0.98 9.9 Chrysene 64 57 55 59 4.7 8.0 C1- Chrysenes 96 90 84 90 5.8 6.0 C2- Chrysenes 120 100 100 107 12 11 C3- Chrysenes 64 62 66 64 2.1 3.0 C4- Chrysenes 43 40 50 44 50 11 Benzo (b) Fluoranthene 2.0 1.7 2.1 1.9 0.20 10 Benzo (c) Pyrene 1.4 13 13 13 0.82 6. Benzo (a) Pyrene 1.7 1.6 1.5 1.6 0.09 5. Benzo (a) Pyrene 1.7 1.8 1.06 1.1 0.08 7.1 Benzo (a) Pyrene 0.00 0.00							9.3
Benzo (a) Anthracene 10.8 8.9 10.2 9.9 0.98 9.1 Chrysene 64 57 55 59 4.7 8.6 C1- Chrysenes 96 90 84 90 5.8 6.6 C2- Chrysenes 120 100 100 107 12 11 C3- Chrysenes 64 62 66 64 2.1 3.3 C4- Chrysenes 43 40 50 44 5.0 11 Benzo (b) Fluoranthene 2.0 1.7 2.1 1.9 0.20 10 Benzo (c) Pyrene 14 13 13 13 0.82 6. Benzo (a) Pyrene 1.7 1.6 1.5 1.6 0.09 5. Benzo (a) Pyrene 1.7 1.6 1.5 1.6 0.02 5. Benzo (a) Pyrene 0.00 0.00 0.00 0.00 0.00 0.00 0.00 Derylene 0.00 0.00 <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td>8.3</td>							8.3
Chrysene 64 57 55 59 4.7 8.1 C1- Chrysenes 96 90 84 90 5.8 6. C2- Chrysenes 120 100 100 107 12 11 C3- Chrysenes 64 62 66 64 2.1 3. C4- Chrysenes 43 40 50 44 5.0 11 Benzo (b) Fluoranthene 2.0 1.7 2.1 1.9 0.20 10 Benzo (b) Fluoranthene 4.5 4.1 4.2 4.3 0.18 4.1 Benzo (a) Pyrene 1.4 13 13 0.82 6. Benzo (a) Pyrene 1.7 1.6 1.5 1.6 0.09 5. Benzo (a) Pyrene 1.20 1.08 1.06 1.1 0.028 7.1 Indeno (1.2.3 - cd) Pyrene 0.00 0.00 0.00 0.00 0.00 0.00 0.00 Dibenzo (a,h) anthracene 0.00 <t< td=""><td></td><td></td><td></td><td></td><td></td><td></td><td>9.1</td></t<>							9.1
C1- Chrysenes 96 90 84 90 5.8 6. C2- Chrysenes 120 100 100 107 12 11 C3- Chrysenes 64 62 66 64 2.1 3. C4- Chrysenes 43 40 50 44 5.0 11 Benzo (b) Fluoranthene 2.0 1.7 2.1 1.9 0.20 11 Benzo (e) Pyrene 14 13 13 13 0.82 6. Benzo (a) Pyrene 1.7 1.6 1.5 1.6 0.09 5. Perylene 1.20 1.08 1.06 1.1 0.02 5. Perylene 1.20 1.08 1.06 1.1 0.08 7. Indeno (1.2.3 - cd) Pyrene 0.00 0.00 0.00 0.00 0.00 0.00 Dibenzo (a,h) anthracene 0.00 0.0 0.00 0.00 0.00 0.00							9.8
C1- Chrysenes 96 90 84 90 5.8 6. C2- Chrysenes 120 100 100 107 12 11 C3- Chrysenes 64 62 66 64 2.1 3. C4- Chrysenes 43 40 50 44 5.0 11 Benzo (b) Fluoranthene 2.0 1.7 2.1 1.9 0.20 11 Benzo (e) Pyrene 14 13 13 13 0.82 6. Benzo (a) Pyrene 1.7 1.6 1.5 1.6 0.09 5. Perylene 1.20 1.08 1.06 1.1 0.02 5. Perylene 1.20 1.08 1.06 1.1 0.08 7.0 Indeno (1.2.3 - cd) Pyrene 0.00 0.00 0.00 0.00 0.00 0.00 0.00	Chrysene	64	57	55	59	4.7	8.0
C2- Chrysenes 120 100 100 107 12 11 C3- Chrysenes 64 62 66 64 2.1 3. C4- Chrysenes 43 40 50 44 5.0 1 Benzo (b) Fluoranthene 2.0 1.7 2.1 1.9 0.20 11 Benzo (k) Fluoranthene 4.5 4.1 4.2 4.3 0.18 4.4 Benzo (e) Pyrene 1.4 13 13 13 0.82 6. Benzo (a) Pyrene 1.7 1.6 1.5 1.6 0.09 5. Perylene 1.20 1.08 1.06 1.1 0.08 7.1 Indeno (1.2.3 - cd) Pyrene 0.00 0.00 0.00 0.00 0.00 0.00 0.00 Dibenzo (a, h) anthracene 0.00 0.00 0.00 0.00 0.00 0.00 0.00	C1- Chrysenes	96	90	84	90	5.8	6.4
C3- Chrysenes 64 62 66 64 2.1 3.3 C4- Chrysenes 43 40 50 44 5.0 1 Benzo (b) Fluoranthene 2.0 1.7 2.1 1.9 0.20 10 Benzo (b) Fluoranthene 4.5 4.1 4.2 4.3 0.18 4.4 Benzo (e) Pyrene 14 13 13 13 0.82 6. Benzo (a) Pyrene 1.7 1.6 1.5 1.6 0.09 5. Perylene 1.20 1.08 1.06 1.1 0.08 7.1 Indeno (1.2.3 - cd) Pyrene 0.00 0.00 0.00 0.00 0.00 0.00 0.00 Dibenzo (a,h) anthracene 0.00 0.0 0.00 0.00 0.00 0.00 0.00							11
C4- Chrysenes 43 40 50 44 5.0 11 Benzo (b) Fluoranthene 2.0 1.7 2.1 1.9 0.20 10 Benzo (b) Fluoranthene 2.0 1.7 2.1 1.9 0.20 11 Benzo (k) Fluoranthene 4.5 4.1 4.2 4.3 0.18 44 Benzo (a) Pyrene 14 13 13 0.82 6. Benzo (a) Pyrene 1.7 1.6 1.5 1.6 0.09 5. Perylene 1.20 1.08 1.06 1.1 0.08 7. Indeno (1.2.3 - cd) Pyrene 0.00 0.00 0.00 0.00 0.00 0.00 0.00 Dibenzo (a, h) anthracene 0.00 0.0 0.00 0.00 0.00 0.00 0.00							3.3
Benzo (b) Fluoranthene 2.0 1.7 2.1 1.9 0.20 10 Benzo (k) Fluoranthene 4.5 4.1 4.2 4.3 0.18 4.1 Benzo (k) Fluoranthene 4.5 4.1 4.2 4.3 0.18 4.1 Benzo (e) Pyrene 14 13 13 13 0.82 6. Benzo (a) Pyrene 1.7 1.6 1.5 1.6 0.09 5. Perylene 1.20 1.08 1.06 1.1 0.08 7.1 Indeno (1.2.3 - cd) Pyrene 0.00 0.00 0.00 0.00 0.00 0.00 0.00 Dibenzo (a,h) anthracene 0.00 0.0 0.00 0.00 0.00 0.00 0.00							11
Benzo (k) Fluoranthene 4.5 4.1 4.2 4.3 0.18 4.4 Benzo (e) Pyrene 14 13 13 13 0.82 6. Benzo (a) Pyrene 1.7 1.6 1.5 1.6 0.09 5. Perylene 1.20 1.08 1.06 1.1 0.08 7.1 Indeno (1.2.3 - cd) Pyrene 0.00 0.00 0.00 0.00 0.00 0.00 Dibenzo (a,h) anthracene 0.00 0.0 0.00 0.00 0.00 0.00 0.00							10
Benzo (e) Pyrene 14 13 13 13 0.82 6. Benzo (a) Pyrene 1.7 1.6 1.5 1.6 0.09 5. Perylene 1.20 1.08 1.06 1.1 0.08 7. Indeno (1.23 - cd) Pyrene 0.00							4.2
Benzo (a) Pyrene 1.7 1.6 1.5 1.6 0.09 5. Perylene 1.20 1.08 1.06 1.1 0.08 7.1 Indeno (1.2.3 - cd) Pyrene 0.00 <							
Perviene 1.20 1.08 1.06 1.1 0.08 7.1 Indeno (1.2.3 - cd) Pyrene 0.00							
Indeno (1.2.3 - cd) Pyrene 0.00 <th< td=""><td></td><td></td><td></td><td></td><td></td><td></td><td>5.5</td></th<>							5.5
Dibenzo (a,h) anthracene 0.00 0.0 0.0 0.00 0.00 0.00 0.00							7.0
							0.00
Benzo (g,h,i) perviene 1.4 1.3 1.5 1.4 0.15 0.0							0.00
	Benzo (g,h,i) perylene	1.4	1.3	1.5	1.4	0.15	0.00
Total Aromatics 11901 10931 10957 11263 553 4.							

STATIST	NUTRIE ICS FOR SURROG Testing Date	ATE CORRECTED D	ATA			
	Replicate #1	Replicate #2	Replicate #3	-		
Alkane Analyte:	Conc (ng/mg)	Conc (ng/mg)	Conc (ng/mg)	Average	Stdev	%RSD
nC-10 Decane	680	630	590	633	45	7.1
nC-11 Undecane	900	850	910	887	32	3.6
nC-12 Dodecane	1300	1400	1200	1300	100	7.7
nC-13 Tridecane	1400	1500	1300	1400	100	7.1
nC-14 Tetradecane	1500	1400	1300	1400	100.0	7.1
nC-15 Pentadecane	2000	1900	2000	1967	58	2.9
nC-16 Hexadecane	1800	1800	1600	1733	115	6.7
nC-17 Heptadecane	1700	1600	1500	1600	100	6.3
Pristane	1800	1800	1600	1733	115	6.7
nC-18 Octadecane	1700	1500	1500	1567	115	7.4
Phytane	1500	1400	1400	1433	58	4.0
nC-19 Nonadecane	1300	1300	1200	1267	58	4.6
nC-20 Elcosana	1700	1500	1600	1600	100	6.3
nC-21 Heneicosane	1500	1400	1400	1433	58	4.0
nC-22 Docosane	1500	1600	1400	1500	100	6.7
nC-23 Tricosane	1500	1400	1300	1400	100	7.1
nC-24 Tetracosane	1500	1800	1600	1633	153	9.4
nC-25 Pentacosane	2500	2500	2400	2467	58	2.3
nC-26 Hexacosane	2400	2600	2200	2400	200	8.3
nC-27 Heptacosane	2000	2100	2000	2033	58	2.8
nC-28 Octacosane	1400	1600	1400	1467	115	7.9
nC-29 Nonacosane	1500	1500	1500	1500	0.0	0.00
nC-30 Triacontane	1000	1000	1100	1033	58	5.6
nC-31 Hentriacontane	1300	1200	1200	1233	58	4.7
nC-32 Dotriacontane	1100	1200	1100	1133	58	5.1
nC-33 Tritriacontane	1100	920	990	1003	91	9.0
nC-34 Tetratriacontane		1000	890	963	64	6.6
nC-35 Pentatriacontane	1100	1100	1200	1133	58	5.1
Total Alkanes	41680	41500	39380	40853	1279	3.1
Aromatic Analyte:	Conc (ng/mg)	Conc (ng/mg)	Conc (ng/mg)	Average	Stdev	%RSD
Naphthalene	350	370	330	350	20	0.00
C1-Naphthalenes	1300	1200	930	1143	191	0.00
C2-Naphthalenes	1800	2000	1900	1900	100	5.3
C3-Naphthalenes	1300	1500	1500	1433	115	8.1
C4-Naphthalenes	790	810	790	797	12	1.4
Fluorene	93	79	92	88	7.5	8.5
C1-Fluorenes	220	230	220	223	5.8	2.6
C2-Fluorenes	270	270	270	270	0.00	0.00
C3- Fluorenes	280	280	300	287	11	4.0
Dibenzothiophene	220	210	220	217	5.7	2.6
C1-Dibenzothiophenes	390	360	390	380	17	4.6
C2-Dibenzothiophenes	490	460	480	477	15	3.2
C3- Dibenzothiophenes	320	290	320	310	17	5.6
Phenanthrene	290	280	290	287	5.8	2.0
C1-Phenanthrenes	590	510	540	547	40	7.4
C2-Phenanthrenes	610	630	620	620	10	1.6
C3-Phenanthrenes	350	370	380	367	15	4.2
C4-Phenanthrenes	160	150	160	157	5.9	4.2 3.8
Anthracene	1.1	1.1	1.0	1.1	0.07	0.00
Fluoranthene	2.5	2.1	2.4	2.3	0.07	8.1
Pidoranmene Pyrene	7.8	6.6	6.9	7.1	0.65	9.1
C1- Pyrenes	120	120	110	117	5.8	4.9
C1- Pyrenes C2- Pyrenes	130	140	130	133	5.8	4.9
C3- Pyrenes	120	120	110	133	5.8	4.9
C4- Pyrenes	73	67	70	70	2.7	3.9
		48	50	49	0.88	3.9
Naphthobenzothiophene		150	150			
C-1 Naphthobenzothiophenes C-2 Naphthobenzothiophenes		150	180	153	5.8	3.8 3.3
				173	5.8	
C-3 Naphthobenzothiophenes		120	110	117	5.8	4.9
Benzo (a) Anthracene	13	12	13	13	0.42	3.3
Chrysene C1 Chrysene	79	75	75	76	2.0	2.6
C1- Chrysenes C2- Chrysenes	120 160	120 140	110 140	117 147	5.8	4.9 7.9
					12	
	110	95 64	100	102	7.4	7.3
C3- Chrysenes	<u>~</u>		59	63	3.6	5.8
C3- Chrysenes C4- Chrysenes	66		4.0			
C3- Chrysenes C4- Chrysenes Benzo (b) Fluoranthene	1.2	1.3	1.3	1.3	0.04	3.2
C3- Chrysenes C4- Chrysenes Benzo (b) Fluoranthene Benzo (k) Fluoranthene	1.2 2.9	1.3 3.0	2.9	2.9	0.02	0.72
C3- Chrysenes C4- Chrysenes Benzo (b) Fluoranthene Benzo (k) Fluoranthene Benzo (e) Pyrene	1.2 2.9 12	1.3 3.0 13	2.9 13	2.9 13	0.02 0.40	0.72 3.2
C3- Chrysenes C4- Chrysenes Benzo (b) Fluoranthene Benzo (k) Fluoranthene Benzo (e) Pyrene Benzo (a) Pyrene	1.2 2.9 12 2.4	1.3 3.0 13 2.3	2.9 13 2.1	2.9 13 2.2	0.02 0.40 0.13	0.72 3.2 5.6
C3- Chrysenes C4- Chrysenes Benzo (b) Fluoranthene Benzo (k) Fluoranthene Benzo (e) Pyrene Benzo (a) Pyrene Perylene	1.2 2.9 12 2.4 0.38	1.3 3.0 13 2.3 0.43	2.9 13 2.1 0.40	2.9 13 2.2 0.40	0.02 0.40 0.13 0.03	0.72 3.2 5.6 7.0
C3- Chrysenes C4- Chrysenes Benzo (b) Fluoranthere Benzo (k) Fluoranthere Benzo (e) Pyrene Benzo (a) Pyrene Perylene Indeno (1.2.3 - cd) Pyrene	1.2 2.9 12 2.4 0.38 0.00	1.3 3.0 13 2.3 0.43 0.00	2.9 13 2.1 0.40 0.00	2.9 13 2.2 0.40 0.00	0.02 0.40 0.13 0.03 0.00	0.72 3.2 5.6 7.0 0.00
C3- Chrysenes C4- Chrysenes Benzo (b) Fluoranthene Benzo (k) Fluoranthene Benzo (e) Pyrene Benzo (a) Pyrene Perylene Indeno (1.2.3 - cd) Pyrene Dibenzo (a,h) anthracene	1.2 2.9 12 2.4 0.38 0.00 0.00	1.3 3.0 13 2.3 0.43 0.00 0.00	2.9 13 2.1 0.40 0.00 0.00	2.9 13 2.2 0.40 0.00 0.00	0.02 0.40 0.13 0.03 0.00 0.00	0.72 3.2 5.6 7.0 0.00 0.00
C3- Chrysenes C4- Chrysenes Benzo (b) Fluoranthere Benzo (k) Fluoranthere Benzo (e) Pyrene Benzo (a) Pyrene Perylene Indeno (1.2.3 - cd) Pyrene	1.2 2.9 12 2.4 0.38 0.00	1.3 3.0 13 2.3 0.43 0.00	2.9 13 2.1 0.40 0.00	2.9 13 2.2 0.40 0.00	0.02 0.40 0.13 0.03 0.00	0.72 3.2 5.6 7.0 0.00

		DDUCT ATE CORRECTED DA :: Day 7	ATA			
	Replicate #1	Replicate #2	Replicate #3	-		
Alkane Analyte:	Conc (ng/mg)	Conc (ng/mg)	Conc (ng/mg)	Average	Stdev	%RSD
nC-10 Decane	1100	1000	1000	1033	58	5.6
nC-11 Undecane	1600	1500	1400	1500	100	6.7
nC-12 Dodecane	1700	1500	1600	1600	100	6.3
nC-13 Tridecane	1800	1800	1600	1733	115	6.7
nC-14 Tetradecane	2100	1900	1900	1967	115	5.9
nC-15 Pentadecane	1700	1900	1600	1733	153	8.8
nC-16 Hexadecane	1800	2000	1800	1867	115	6.2
nC-17 Heptadecane	2100	2000	1800	1967	153	7.8
Pristane	1700	1700	1600	1667	58	3.5
nC-18 Octadecane	1800	2000	1700	1833	153	8.3
Phytane	1600	1500	1400	1500	100	6.7
nC-19 Nonadecane	1600	1800	1500	1633	153	9.4
nC-20 Elcosane	1800	1900	1700	1800	100	5.6
nC-21 Heneicosane	1700	1800	1600	1700	100	5.9
nC-22 Docosane	1700	1800	1500	1667	153	9.2
nC-23 Tricosane	1800	1900	1600	1767	153	8.6
nC-24 Tetracosane	1500	1700	1500	1567	115	7.4
nC-25 Pentacosane	1800	1800	1800	1800	0.00	0.00
nC-26 Hexacosane	1800	1800	1800	1800	0.00	0.00
nC-27 Heptacosane	1700	1800	1600	1700	100	5.9
nC-28 Octacosane	1700	1800	1600	1700	100	5.9
nC-28 Octacosane nC-29 Nonacosane						
nC-29 Nonacosane nC-30 Triacontane	1800	1800	1600	1733	115 85	6.7
	1100	940	970	1003		8.5
nC-31 Hentriacontane	1100	950	1000	1017	76	7.5
nC-32 Dotriacontane	1000	1000	980	993	12	1.2
nC-33 Tritriacontane	880	870	900	883	15	1.7
nC-34 Tetratriacontane	930	950	840	907	59	6.5
nC-35 Pentatriacontane	1200	1200	1100	1167	58	4.9
Total Alkanes	44110	44610	40990	43237	1962	4.5
Aromatic Analyte:	Conc (ng/mg)	Conc (ng/mg)	Conc (ng/mg)	Average	Stdev	%RSD
Naphthalene	330	290	330	317	23	7.3
C1-Naphthalenes	1300	1200	1200	1233	58	4.7
C2-Naphthalenes	2200	1900	1900	2000	173	8.7
C3-Naphthalenes	1300	1400	1300	1333	58	4.3
C4-Naphthalenes	740	700	660	700	40	5.7
Fluorene	79	83	78	80	2.4	3.0
C1-Fluorenes	240	220	200	220	20	9.1
C2-Fluorenes	290	270	250	270	20	7.4
C3- Fluorenes	270	260	270	267	5.7	2.1
Dibenzothiophene	220	220	200	213	12	5.4
C1-Dibenzothiophenes	350	350	350	350	0.00	0.00
C2-Dibenzothiophenes	480	480	440	467	23	4.9
C3- Dibenzothiophenes	330	300	290	307	21	6.8
Phenanthrene	300	280	270	283	15	5.4
C1-Phenanthrenes	560	550	490	533	38	7.1
C2-Phenanthrenes	630	590	560	593	35	5.9
C3-Phenanthrenes	390	370	360	373	15	4.1
C4-Phenanthrenes	130	130	140	133	5.8	4.3
Anthracene	1.4	1.3	1.5	1.4	0.07	5.0
Fluoranthene	2.3	2.2	2.4	2.3	0.09	4.1
Pyrene	8.8	8.8	8.9	8.8	0.08	0.91
C1- Pyrenes	110	100	94	101	7.9	7.8
C2- Pyrenes	130	120	120	123	5.8	4.7
C3- Pyrenes	110	110	100	107	5.8	5.4
C4- Pyrenes	65	59	63	62	3.1	5.0
Naphthobenzothiophene	47	49	46	48	1.4	2.9
-1 Naphthobenzothiophenes	180	160	160	167	12	6.9
-2 Naphthobenzothiophenes	170	160	160	163	6.1	3.7
-3 Naphthobenzothiophenes	120	100	110	110	10	9.1
Benzo (a) Anthracene	15.3	14.3	13.0	14	1.1	8.0
Chrysene	80	69	68	72	6.6	9.1
C1- Chrysenes	120	100	110	110	10	9.1
C2- Chrysenes	145	130	140	138	7.7	5.5
C3- Chrysenes	100	96	87	94	6.5	6.9
C4- Chrysenes	61	57	64	61	3.4	5.7
						5.7
Benzo (b) Fluoranthene	1.7	1.5	1.5	1.6	0.08	
	2.3	2.3	2.3	2.3	0.04	1.9
Benzo (k) Fluoranthene	12	11	12	12	0.46	3.9
Benzo (e) Pyrene		1.8	2.1	2.1	0.19	9.2
Benzo (e) Pyrene Benzo (a) Pyrene	2.2			0.30	0.02	6.4
Benzo (e) Pyrene Benzo (a) Pyrene Perylene	0.31	0.28	0.31			
Berzo (e) Pyrene Berzo (a) Pyrene Perylene Indeno (1.2.3 - cd) Pyrene	0.31 0.00	0.00	0.00	0.00	0.00	0.00
Berzo (e) Pyrene Benzo (a) Pyrene Perylene Indeno (1.2.3 - cd) Pyrene Dibenzo (a,h) anthracene	0.31 0.00 0.00	0.00 0.00	0.00 0.00	0.00 0.00	0.00 0.00	0.00 0.00
Benzo (e) Pyrene Benzo (a) Pyrene Perylene Indeno (1.2.3 - cd) Pyrene	0.31 0.00	0.00	0.00	0.00	0.00	0.00

CC T Thild	CONTROL, REPLICATE 1 Testing Date: Day 28 Initial Oil Weight: 500 mg Final Extracted Volume: 15 mL	u 5 mL	Final	CONTROL, REPLICATE 2 Testing Date: Day 28 Initial Oil Weight: 500 mg Final Extracted Volume: 15 mL	s mL	o . II	CONTROL, REPLICATE 5 Testing Date: Day 28 Initial Oil Weight: 500 mg Final Extracted Volume: 15 mL	u S mL
	Concentration (ng/mg)	Surrogate Corrected (ng/mg)		Concentration (ng/mg)	Surrogate Corrected (ng/mg)		Concentration (ng/mg)	Surrogate Corrected (ng/mg)
-			Alkane Analyte:			Alkane Analyte:		
IIC-10 Decare	101	104	IIC-10 Decalle		110	IIC-10 Decare		108
nC-12 Dodecare	2084	2149	nc-11 Undecane nC-12 Dodecane	2229	2298	nC-11 Undecane nC-12 Dodecane	1951	2033
nC-13 Tridecare	2256	2326	nC-13 Tridecane	2493	2570	nC-13 Tridecane	2491	2595
nC 15 Pentadecane	2338	2410 2556	nC-14 letradecane	2355	2428	nC-14 Letradecane nC-15 Pentadecane	2222	2318
nC-16 Hexadecane	2469	2545	nC-16 Hexadecane	2407	2481	nC-16 Hexadecane	2304	2400
nC-17 Heptadecane	2325	2397	nC-17 Heptadecane	2355	2427	nC-17 Heptadecane	2160	2250
Pristane	1375	1415	Pristane	1542 2668	1590	Pristane	1488	1550
nucrio Uctatecane Phyrane	1326	2470	Phytane Phytane	1311	1351	nu-16 Uctadecane Phytane	1230	1281
nC-19 Nonadecane	1952	2012	nC-19 Nonadecane	2221	2290	nC-15 Nonadceane	2032	2117
nC-20 Eisoaaro	2045	2108	nC-20 Eleosano	2185	2252	nC-20 Eleosano	1088	2060
nC-21 Heneicosane	2035	2096	nC-21 Heneicosane	2032	2095	nC-21 Heneicosane	1828	1904
nC-22 Uocosare	1855	1912	nC-22 Docosane	1880	1936	nC-22 Decosare	16/1	1/40
nu-23 Incosare	20/L	1/02	nC-23 Incosane	70/1	92/I	nu-23 Incosane	1385	1004
n0-25 Pentacosane	1626	1677	nC-25 Pentacosane	1761	1815	nC-25 Pentacosare	1739	18/1
nC-26 Hexacosane	1565	1614	nC-26 Hexacosane	1764	1819	nC-26 Hexacosane	1704	1775
nC-27 lleptacosane	1204	1203	nC-27 lleptacosane	7001	1040	nC-27 lleptacosane	1220	1277
nC-28 Octacosane	1110	1145	nC-28 Octacosane	1165	1201	nC-28 Octacosane	1010	1052
nC-29 Nonacosane	1140	1175	nC-29 Nonacosane	1195	1232	nC-29 Nonacosane	1090	1136
nC-30 Triacontanc	1037	1069	nC-30 Triacontanc	1030	1062	nC-30 Triacontanc	080	1021
nC-31 Hentriacontane	964	994	nC-31 Hentriacontane	967 267	997 200	nC-31 Hentriacontane	870	906
nL-32 Dotriacontane	8//	802	nC-32 Dotriacontane	208 878	058	nC-32 Dotriacontane	764	801 706
IIC-34 Telratriacontane	752	200	IIC-34 Tetratriacontare	913	94:	IIC-34 Teliatriacontarie	198	897
nC-35 Pentatriacontane	1102	1136	nC-35 Pentatriacontare	1308	1349	nC-35 Pertatriacontane	1303	1357
					10011			
Total Alkanes	43613	44962	Total Alkanes	45808	47225	Total Alkanes	42446	44215
Aromatic Analyte: Naciditation	252	287	Aromatic Analyte: Nachfhalene	247	275	Aromatic Analyte: Natifitialene	245	275
C1-Naphthalenes	992	1127	C1-Naphthalenes	992	1102	C1-Naphthalenes	973	1093
CZ-Naphthalenes	1634	1857	C2-Naphthalenes		9891	C2-Naphthalenes	1437	1614
C3-Naphthalenes	1374	1561	C3-Naphthalenes		1579	C3-Naphthalenes	1360	1528
C4-Naphthalenes	769	/80	C4-Naphthalenes	682	28	C4-Naphmalenes	2	805 105
C1-Fluorenes	234	265	C1-Fluorenes	259	- 100	C1-Fluorenes	231	260
CZ-Fluorenes	357	406	CZ-Fluorenes	3/4	416	CZ-Fluorenes	345	388
C3- Fluorenes	287	326	C3- Fluorenes	336	373	C3- Fluorenes	335	377
Diberzothiophene	219	249	Dibenzothiphene	238	264	Dibenzothiophene	203	228
C2 Dihanzothinhenes	4.0 625		C1-Unenzonninnenes	194 194	705	C.1-Unenzorninpnenes C2 Nikenzothirohense	1	40.5 644
C3- Dibenzothiophenes	391	445	C3- Dibenzothiophenes	200	425	C3- Dibenzothicphenes	351	380
Hhenanthrere	285	324	Phenanthrene	067	322	Phenanthrene	255	206
C1-Phenanthrenes	657	747	C1-Prenanthrenes	999 193	743	C1-Phenarthrenes	595	668 201
C3-Phenanthrenes	477	542	C3-Phenanthrenes	482	100	CZ-Phenarthrenes	432	0 89 88 89
C4-Phenanthrenes	180	204	C4-Phenanthrenes	175	194	C4-Phenarthrenes	181	207
Anthracere	1.2	1.3	Anthracene	1.93	2.1	Anthracene	1.3	
Fluoralliere	2.9	3.2	Pluuranthene Decos	2.8	1.5	Fluoranthene	2.8	n ţ
C1- Pyrenes	°.8	112	C1- Pvrenes	107	119	C1- Pvrenes	105	- 11
C2- Pyrenes	128	145	C2- Pyrenes	136	152	C2- Pyrenes	127	143
C3- Fyrenes	115	131	C3- Pyrenes	129	143	C3- Pyrenes	122	137
C4 Hyrenes Manhthohanzothionhana	رہ حم	98 97	C4 Pyrenec Manhthohanzothionhara	83	82	C4 Pyrenec Manhthohanzothinnhana	2 2	8 8
C-1 Naphthobenzothlophenes	175	5 65	C-1 Naphthobenzothlophenes	8 <u>8</u>	202	C-1 Naphthobenzothlophenes	76 165	185
C-2 Naphthobenzothiophenes	194	220	C-2 Naphthobenzothiophenes	197	219	C-2 Naphthobenzothiophenes	178	200
-3 Naphthobenzothiophenes	128	146	C-3 Naphthobenzothiophenes	138	154	C-3 Naphthobenzothiophenes	119	134
Benzo (a) Anthracene	6.)	0.9 20	Benzo (a) Anthracene	4.9 4 0	10.5 AG	Benzo (a) Anthracene Christene	9.0	10
C1- Chryachea	40 68	101	C1- Chryaenea	8 8	26	C1- Chryaonea	60	88
C2- Chrysenes	102	116	C2- Chrysenes	103	114	C2- Chrysenes	96	107
CO- Chrysenes	55	76	C3- Chrysenes	8 8	£ 1	C3- Chrysenes	9 9	30
C4- Chrysenes Hanzh (h) Fluoranthana	74	80 F	C4- Unitysenes Hanzo (N) -litorethere	00 P	8 5	C4- Chrysenes Henzo (h) Fluoranthene	4 2 2	5 r
Renzn (k) Fluoranthene	4.8	55	Benzo (k) Fluoranthere	5 ¥	5.0	Renzo (k) Fluoranthese	4.5	4 107
Benzo (e) Pyrene	£ (£ 5	Benzo (e) Pyrene	12	t 13	Benzo (e) Pyrene	= ;	£ ;
Lenzo (a) ryrene Perviene	27 U	9.0	Denviene Perviene	180	F 7	Denzo (aj Fyrene Perviene	0.86	10
Inceno (1,2,3 - cd) Pyrene	0.22	0.25	Indeno (1,2,3 - cd) Pyrene	0.25	0.28	Indeno (1,2,5 - cd) Pyrene	0.23	0.3
Dibenzo (a,h) anthracene Benzo (o h i) nerviene	0.59	0.67 2.6	Dibenzo (a,h) anthracere Barzo (a h i) pendana	0.8 2 c	0.84	Cibenzo (a,h) anthracene Benzo (o h i) repulsione	0.68	0.8
Total Aromatics	11264	12800	Total Aromatics	11584	12871	Total Aromatics	10689	12010
% Surrogata Recovery			% Surrogate Recovery			% Surrogate Recovery		
5 Alpha Androstane	0.97	1.00	5 Alpha Androstane	0.97	1.00	5 Alpha Androstane	0.96	1.00
Phenalthrane c- 10	0.00	1.00	Phenanthrene d- 10		1.00	Phenanthrene G-10		0.1

NUTR Test Initial Final Extr	NUTRIENT, REPLICATE 1 Testing Date: Day 28 Initial Oil Weight: 500 mg Final Extracted Volume: 15 mL	15 mL	rinal Tinal	NUTRIENT, REPLICATE 2 Testing Date: Day 28 Initial Oil Weight: 500 mg Final Extracted Volume: 15 mL	ig ML	Ini Ini Final of	NUTRIENT, REPLICATE 3 Testing Date: Day 28 Initial Oil Weight: 500 mg Final Extracted Volume: 15 mL	9 8 10 10
Can	Concentration (no/mo)	Surroate Corrected (nalma)		Concentration (na/ma)	Surmaate Corrected (na/mal		Concentration (na/ma)	Surronate Corrected Innima)
-	ì		Alkane Analyte:	i n		Alkane Analyte:	Î	5
nC-10 Decare	0.00	0 2	nC-10 Decane nC-11 Undecane	0000	0 5	nC-10 Decane nC-11 Undecane	0.00	0 8
nC-12 Dodecare	115.00	116	nC-12 Dodecane	132.00	132	nC-12 Dodecane	128.00	128
nc-14 Tetradecare	148.00	149	nC-14 Tetradecane	158.00	155	nc-14 Tetradecane	178.00	8 E
n0-15 Pentadecane	169.00	171	nC-15 Pentadecane	179.00	179	nC-15 Pentadecane	192.00	192
nu-10 received in number	248	294	nc-10 resourceme	967	907	nu-romexercane nu-1/ Heptadecane	241	241
Priskane	1631	1848 200	Prislare	1624	1824 107	Priskne of 18 October	1621	1621 180
Phytane Phytane	1745	1760	Phytane	134	13/2	nu-lo uctadecane Phytane	1728	108
nC-19 Nonadecane	289	292	nC-19 Nonadecane	315	315	nC-19 Nonadecane	340	340
nC-20 Elosare nC 21 Heneicosane	149.0	151	nC-20 Ecosane nC 21 Henebosane	125.0	c12 125	nC-20 Elcosane nC 21 Heneicosane	215.0	112
nC-22 Doceand	135	136	nC-22 Docosanc	142	142	nC-22 Decosanc	140	140
nC-23 Tricosare	116.0	21:1	nC-23 Tricosane	149.0	149	nC-23 Tricosane	153.0	<u>1</u>
nC-25 Pentacosane	228	230	nC-25 Pentacosane	229	229	nC-25 Pentacosane	227	227
nC-26 Hexacosane	196	198	nC-26 Hexacosane		188	nC-26 Hexacosane	178	178
nC-27 Heptacosane	142	143	nC-27 Heptacosane	143	143	nC-27 Heptacosane	144	144
nC-29 Nonacosane	147	148	nC-29 Nonacosane	146	146	nC-29 Nonacosane	15	5
nC-30 Irracontane	711	118	nC-30 Inacontane	12/	12/	nC-30 Inacontane	152	152
nC-31 Hentriacontane nC-32 Dotriacontane	112	113	nC-31 Hentriacontane nC-32 Dotriacontane	121	121 135	nC-31 Hentriacontane nC-32 Dotriacontane	111 130	111
nC-33 Tritriacontone	0.00	£	n0-33 Tritrineontono	101	100	nC-33 Tritrianantano	0.50	8
nC-34 Tetratriacontane nC-35 Pentatriacontane	50.00 41.00	5 1	nC-34 Tetratriacontare nC-35 Pentatriacontare	62.00 45.00	62 45	nC-34 Tetratriacontane nC-35 Pertatriacontane	49.00	84 00
Total Alkanes	7091	7162	Total Alkanes	7188	7188	Total Alkanas	7267	7267
Assessed in South Asses			Assessed in the base			Assessing Assesses		
Aromatic Analyte: Naphthaere	0.00	0	Aromatic Analyte: Naphthalene	0:00	0	Aromatic Analyte: Naphthalene	0.00	0
C1-Naphilhalenes	85	115	C1-Raphthalenes		66	C1-Maphilitalenes	82 00 201	109
C2-Naphthalenes C3-Naphthalenes	644 844	958 1140	C2-Naphthalenes C3-Naphthalenes		1034	C2-Naphthalenes C3-Naphthalenes	805 856	10/3
C4-Naphthalenes	701	948	C4-Naphthalenes		912	C4-Naphthalenes	712	949
Fluorene C1 Fluorenee	21.0	28 273	Fluorene C1 Fluorenee	22.0 886	29	Fluorene C1 Fliorenee	26.0 237	35 346
C2-Fluorenes	301	407	C2-Fluorenes	312	405	C2-Fluorenes	315	420
C3- Fluorenes	234	316	C3- Fluorenes	265	344	C3- Fluorenes	265	353
Ulberzothiophene C1-Dibenzothiophenes	225	305	Ulbenzothophene C1-Dibenzothiophenes	69 264	90 343	Ulbenzothiophene C1-Dibenzothicphenes	73 249	9/ 332
C2-Dibenzothiophenes	503	680	C2-Dibenzothiophenes	492	639	C2-Dibenzothicphenes	498	664
C3 Dibenzothiophenes Phonosthrond	319	431 68	C3 Dibenzothiophenee	349	453 74	C3 Dibenzothicphense Phonenthrond	346 61	461 8
C1-Phenanthrenes	327	441	C1-Phenanthrenes	325	422	C1-Phenarthrenes	325	433
C2-Phenanthrenes	293	181	U2-Phenanthrenes	225	6/8	UZ-Phenarthrenes	547	R7. \$
C4-Phenanthrenes	162	219	C4-Phenanthrenes	171	222	C4-Phenarthrenes	179	239
Anthracere	66	30	Anthracene	24	31	Anthracene	21	28
Pyrene	2.0	2.2	Pyrene Pyrene	21	0.4	Pyrene	210	10.1
C1- Pyrenes	06	121	C1- Pyrenes	119	155	C1- Pyrenes	109	145
C3 Fyrenee	3 5	153	C.2 Pyrenes C.3 Pyrenes	135	10	C.3 Pyrenes	146	\$ \$
C4- Fyrenes	67	91	C4- Pyrenes	14	92	C4- Pyrenes	58 F	113
C-1 Naphthobenzothiophenes	165	223	C-1 Naphthobenzothiophenes		229	C-1 Naphthobenzothiophenes	178	237
C-2 Naphthoben zothrophenes	100	162	C-2 Naphthobenzollhophenes		62	C-2 Naphthobencothiophenes	196	204
U-3 Naphtmobenzotniophenes Benzo (a) Anthracene	120	9.3	C-3 Naprimopenzomiophenes Benzo (a) Anthracene	9 9	57	C-3 respiritobenzotniopnenes Benzo (a) Anthracene	5 S	7.2
Chrysene	42	57	Chrysene Chrysene	43	8 ç	Chrysene	47	63 16
CI- CIII yadita C2 Chryeenee	135	182	C1 Chrysenes C2 Chrysenes	142	184	C2 Chrysenes	152	203
C5- Chrysenes C4- Chrysenes	105 79	142	C3- Chrysenes C4- Chrysenes	116 69	150 90	C3- Chrysenes C4- Chrysenes	122 91	163
Benzo (b) Fluoranthene	0.63	6.0	Benzo (b) Fluoranthene	0.75	1.0	Benzo (b) Fluoranthene	0.68	6.0
Benzo (k) Fluoranthene	3.1	4.2	Benzo (k) Fluoranthere	3.9	5.0	Benzo (k) Fluoranthene	3.7	4.9
Benzo (a) Pyrene	1.5	2.0	Benzo (a) Pyrene	o 1:	2.3	Benzo (a) Pyrene	e 1.6	2.2
Peryland Indona (1.2.3 od). Suman	0.25	0.3	Perylene Indeep (1.2.2 ed) Breese	0.21	0.3	Peryland	0.23	0.3
Ubenzo (a,h) anthracane Benzo (a,h) nandene	2.0	9720	Utbenzo (a,h) anthracere Berzo (a,h) anthracere	0.20	970	Libenzo (a,h) anthracene Barzo (a,h) reordene	0.21	0.28
Total Aromatics	7225	9764	Total Aromatics	7540	9793	Total Aromatics	7670	10227
% Surrugate Recovery	000	90 1	Surrugate Recovery	1.00	4.00	% Surrugate Recovery		6
5 Alpha Androstane Phenanthrene d-10	0.99 0.74	1.00	5 Alpha Androstane Phenanthrene d-10	0.77	1.00	5 Alpha Ancrostane Phenanthrene d-10	1.00 0.75	1.00
					1 - 1	-	l	

SARV 	SARVA PRODUCT, REPLICATE 1 Testing Date: Day 28 Initial OII Weight: 500 mg Final Extracted Volume: 15 mL	LTE1 1 mg 15 mL	SAR\ In Final	SARVA PRODUCT, REPLICATE 2 Testing Date: Day 28 Initial Oil Weight: 500 mg final Extracted Volume: 15 ml	ATE 2 8 mg 15 mL	SARVA T Init	SARVA PRODUCT, REPLICATE 3 Testing Date: Nay 28 Initial OII Weight: 500 mg Final Extracted Volume: 15 mL	TE 3 9 5 mL
Alkane Analvte:	Concentration (ng/mg)	Surrogate Corrected (rg/mg)	Alkane Analyte:	Concentration (ng/mg)	Surrogate Corrected (ag/mg)	Alkane Analyte:	Concentration (ng/mg)	Surrogate Corrected (ng/mg)
nC-10 Decare	0.00	0	nC-10 Decane	0.00	0	nC-10 Decane	0.00	0
nC-12 Dodecare	0.00		nC-11 Undecane nC-12 Dodecane	0.00		nC-11 Undecane nC-12 Dodecane	0.0	
nC-13 Tridecare	0.00	0 1	nC-13 Tridecane	0.00	0 .	nC-13 Tridecane	0.00	0 1
nc-15 Pentadecane	00.0		nc-15 Pentadecane	00.0		nc-15 Pentadecane	0.0	
nC-16 Hexadecane	00:00	0	nC-16 Hexadecane	00:0	0	nC-16 Hexadecane	0.00	0
nu-17 heptadecane Prislane	324	391 391	nu-17 neptadecane Pristare	9 g	32 423	nu17 Heptadecane Pristane	5 8	32 431
nC-18 Octadecane	11	13	nC-18 Octadecane	12	15	nC-18 Octadecane	13	16
Phytane nC.19 Nonadecane	315 11	380 13	Phytane nC-19 Nonadecane	297 10	380	Phytene nC.19 Nonadecane	286 8.6	362
nc-20 Elcosare	7.9	2 თ	nc-19 revitadedate nc-20 Ecosane	72	7 6	nc-remuneredate nc-20 Elcosane	6.7	- 60
nC 21 Heneicocane	5 5	13	nC 21 Henebosane	9	13	nC 21 Heneicocane	9	£ :
nC-22 Doposare	Q Q	4 7	nC-23 Tricosane	= =	4 S	nC-22 Dicosanc nC-23 Tricosane	2 :-	5 <u>5</u>
nC-24 Tetracosane	18	21	nC-24 Tetracosane	16	21	nC-24 Tetracosane	16	20
nC-25 Pentacosane	107	129	nC-25 Pentacosane	8	12	nC-25 Pentacosane	8	112
nC-27 Heptacosane	35	86	nu-zo nexacosane nC-27 Hentacosane	30	101	nC-20 Hexacosane nC-27 Heptacosane	28	35
nC-28 Octacoanno	5	30	nC-28 Octacosanc	24	30	nC-28 Octacosanc	53	5
nC-29 Nonacosane	8	46	nC-29 Nonacosane	35	45	nC-29 Nonacosane	8	42
nC-30 Irriacontane	59 E	32	nC-30 Inacontane	5 5	2/	nC-30 Iriacontane	2 €	22 16
nC-32 Dotriacontane	6.9	2 00	nC-32 Dotriacontane	6.9	2 თ	nC-32 Dotriacontane	7.6	10
n0-33 Tritriacontane n0-34 Tetratriacontane	8	¢ c	n0-33 Tritrineontone n0-34 Tetratrineontone	° 000	5 c	n0-33 Tritriacontano n0-34 Tatratriacontana	8 00	¢ c
nC-35 Pentatriacontane	00.0	, o	nC-35 Pentatriacontare	00.0	0.00	nC-35 Pertatriacontane	0.0	0.0
Total Alkanes	1086	1308	Total Alkanes	1040	1333	Total Alkanss	1018	1289
Aromatic Analyte:			Aromatic Analyte:			Aromatic Analyte:		
Naphthalere C1 Monthlatere	0.0	0.0	C1 Northfultrate	0.00	00.0	C1 Monthered	0.0	0.0
C2-Naphthalenes	216	288	C2-Naphthalenes	208	293	C2-Naphthalenes	204	287
C3-Naphthalenes	819	1092	C3-Naphthalenes	744	1047	C3-Naphthalenes	692	975
Fluorene	12	16.5	Fluorene	11	16	Fluorene	- 	14.4
C1 Fluorenes	8	131	C1 Fluorenes	96	133	C1 Fluorenee	83	131
C2-Fluorenes C3- Fluorenes	228 187	304 248	C2-Fluorenes C3- Fluorenes	253 203	357 206	C2-Fluorenes C3- Fluorenes	279 221	393 311
Diberzothiophene	14	18.3	Dibenzothiphene	13	18	Dibenzothiophene	12	16.9
C1-Ulbenzothiophenes C2-Dihenzothiophenes	415	240	C1-Dibenzothiophenes C2-Dibenzothiophenes	165	233	C1-Ulbenzotnicpnenes C2-Dihenzothicnhenes	152	214 505
C3 Dibenzothiophenee	262	350	C3 Dibenzothiophenee	261	367	C3 Dibenzothicphense	261	372
C1-Phenanthrenes	14 263	350	C1-Phenanthrenes	13 245	345	C1-Phenarthrenes	11	16 328
UZ-Phenanthrenes	455	606	C2-Phenanthrenes	484	99	C2-Phenarthrenes	517	1.28
C3-Phenanthrenes C4-Phenanthrenes	344 139	458 185	C3-Phenanthrenes C4-Phenanthrenes	362 144	510 203	C3-Phenarthrenes C4-Phenarthrenes	385 151	542 213
Anthracere	1.5	2 U	Anthracene	15	2.1	Anthracene	14	0 0
Fluoranthere	1.4 6.2	1.8	Fluoranthene Pvrene	1.3	1.8	Fluoranthene	1.3	1.8 9.8
C1- Pyrenes	72	26	C1- Pyrenes	67	95	C1- Pyrenes	64	ò
C2- Fyrenes	102	136	C2- Pyrenes	88	134	C2- Pyrenes	91	129
C4- Fyrenes	62	82	C4- Pyrenes	8	74	C4- Pyrenes	46	65
Naphthobenzothiophene C_1 Manhthobenzothionhene	41	55 180	Naphthobenzothiphene C-1 Nanhthohenzothionhenee	44	53 198	Naphthobenzothiophene C.1 Nanhthohenzothionhenee	47	99 UUC
C-2 Naphthoben zuthophenes	2	222	C-2 Naphiliubenzolimphenes	20	577	C-2 Naphiliobenzollinphenes	4	207
C-3 Naphthobenzothiophenes Renzo (a) Anthracene	100	133	C-3 Naphthobenzothiophenes Renzo (a) Anthracene	95 8	134 10.8	C-3 Naphthobenzothiophenes Benzo (a) Anthracene	93	131 103
Chrysene	48	64	Chrysene	50	70	Chrysene	52	73
C1- Chrysenes	88 105	118	C1- Chrysenes	93 101	132	C1- Chrysenes	66 8	140
C3- Chrysenes	84	112	C3- Chrysenes	8	125	C3- Chrysenes	94	133
C4- Chrysenes	5 F	69	C4- Chrysenes	8 2	74	C4- Chrysenes	55	77
Benzo (b) Fluoranthene Benzo (k) Fluoranthene	3.0	4.0	Benzo (b) Fluoranthere	2.8	8'0 3'3	Benzo (k) Fluoranthene	2.7	3.7
Benzo (e) Pyrene Renzn (a) Pyrene	8.1	11 2.0	Benzo (e) Pyrene Benzo (a) Pyrene	8	12 2.0	Benzo (e) Pyrene Benzo (a) Pyrene	8 1	12 2.0
Peryland	0.65	0.73	Perylene	0.5	0.77	Parylana	0.55	0.78
Incieno (1,2,3 - cd) Pyrene Dihenza (a h) anthracene	0.00	0.00	Indeno (1,2,3 - cd) Pyrene Uthenzo (a h) anthracere	00.0	00.0	Indeno (1,2,5 - cd) Pyrene Libenzo (a h) anthracene	0.0	0.0
Benzo (g,h,i) perylene	1.8	2.40	Berzo (g,h,i) perylene	22	3.05	Benzo (g,h,i) perylene	2.5	3.57
Total Aromatics	5416	7222	Total Aromatics	5347	7531	Total Aromatics	5379	7576
			0					
5 Alpha Androstane	0.83	1.00	5 Alpha Androstane	0.78	1.00	5 Alpha Androstane		1.00
Phenanthrene c-10	0.75	1.00	Phenanthrene d-10	0.71	1.00	Phenanthrene d-10	0.71	1.00

STATIST	CONTRO ICS FOR SURROGAT Testing Date:	E CORRECTED DATA]		
		-				
Alkane Analyte:	Replicate #1	Replicate #2	Replicate #3	A	Stdev	%RSD
nC-10 Decane	Conc (ng/mg) 104	Conc (ng/mg) 110	Conc (ng/mg) 106	Average 107	3.3	%RSD 3
nC-11 Undecane	1074	1101	1084	1087	- 3.3 14	1
nC-12 Dodecane	2149	2298	2033	2160	133	6
nC-13 Tridecane	2326	2570	2595	2497	149	6
nC-14 Tetradecane	2410	2428	2318	2386	59	2
nC-15 Pentadecane	2556	2626	2377	2520	129	5
nC-16 Hexadecane	2545	2481	2400	2475	72.7	3
nC-17 Heptadecane	2397	2427	2250	2358	95	4
Pristane	1415	1590	1550	1519	92	6
nC-18 Octadecane Phylane	2470 1368	2635 1351	2355 1281	2487 1333	141 45.88	6 3
nC-19 Nonadecane	2012	2290	2117	2140	140	7
nC-20 Eicosane	2108	2252	2069	2143	97	5
nC-21 Heneicosane	2096	2095	1904	2032	110	5
nC-22 Docosane	1912	1938	1/40	1864	107	6
nC-23 Tricosane	1762	1755	1664	1727	54	3
nC-24 Tetracosane	1709	1777	1543	1676	121	7
nC-25 Pentacosane	1677	1815	1811	1768	79	4
nC-26 Hexacosane	1614	1819	1775	1736	108	6
nC-27 Heptacosane nC-28 Octacosane	1303 1145	1348 1201	1277	1309	36 75	3
nC 28 Octacosane nC-29 Nonacosane	1145 1175	1201	1052 1136	1133 1181	75 48	4
nC-29 Nonacosane nC-30 Triacontane	1069	1232	1021	1050	48 26	2
nC-31 Hentriacontane	994	997	906	966	52	5
nC-32 Dotriacontane	802	830	801	811	16	2
nC-33 Tritriacontane	859	903	795	853	54	6
nC-34 Tetratriacontane	776	941	897	871	86	10
nC-35 Pentatriacontane	1136	1349	1357	1281	126	10
Total Alkanes	44962	47225	44215	45467	1568	3
Aromatic Analyte:	Conc (ng/mg)	Conc (ng/mg)	Conc (ng/mg)	Average	Stdev	%RSD
Naphthalene	287	275	275	279	7.0	3
C1-Naphthalenes	1127	1102	1093	1108	18	2
C2-Naphthalenes	1857	1895	1614	1789	152	9
C3-Naphthalenes	1561	1579	1528	1556	26	2
C4-Naphthalenes	786	758	805	783	24	3
Fluorene C1 Fluorenes	107 265	114 287	103 260	108 271	5.68 15	5
C2-Fluorenes	406	416	388	403	14	3
C3- Fluorenes	326	373	377	359	28	8
Dibenzothiophene	249	264	228	247	18.3	7
C1-Dibenzothiophenes	495	495	453	481	24	5
C2-Dibenzothiophenes	721	705	644	690	41	6
C3- Dibenzothiophenes	445	425	395	421	25	6
Phenanthrene	324	322	286	311	21	7
C1-Phenanthrenes C2-Phenanthrenes	747	743 768	668	719	44	6 6
C2-Phenanthrenes	770 542	535	695 486	745 521	43 31	6
C3-Phenanthrenes	204	194	207	202	6.7	3
Anthracene	1.3	2.1	1.4	1.6	0.4	0
Fluoranthene	3.2	3.1	3.2	3.2	0.06	2
Pyrene	11.2	13.3	13.4	12.6	1.25	10
C1- Pyrenes	112	119	118	116	4.0	3
C2- Pyrenes	145	152	143	146	45	3
C3 Pyronos	131	143	137	137	6.3	5
C4- Pyrenes Naphthobenzothiophene	86 64	82	83 59	83 63	2.1 3.4	2 5
C-1 Naphthobenzothiophenes	199	65 203	185	196	3.4 9	5
C-2 Naphthobenzothiophenes	220	203	200	213	11	5
C-3 Naphthobenzothiophenes	146	154	134	144	10	7
Benzo (a) Anthracene		10.5	10.1	9.8	0.81	8
Chrysene	59	65	56	60	4.6	8
C1- Chrysenes			90	96	5.4	6
C2- Chrysenes	101	97				
	101 116	114	107	112	5	4
C3- Chrysenes	101 116 92	114 95	107 85	91	5	6
C3- Chrysenes C4- Chrysenes	101 116 92 59	114 95 55	107 85 54	91 56	5 3	6 5
C3- Chrysenes C4- Chrysenes Benzo (b) Fluoranthene	101 116 92 59 1.9	114 95 55 2.1	107 85 54 2.1	91 56 2.0	5 3 0.11	6 5 6
C3- Chrysenes C4- Chrysenes Benzo (b) Fluoranthene Benzo (k) Fluoranthene	101 116 92 59 1.9 5.5	114 95 55 2.1 5.0	107 85 54 2.1 5.1	91 56 2.0 5.2	5 3 0.11 0.26	6 5 6 5
C3- Chrysenes C4- Chrysenes Benzo (b) Fluoranthene Benzo (k) Fluoranthene Benzo (e) Pyrene	101 116 92 59 1.9 5.5 13	114 95 55 2.1 5.0 13	107 85 54 2.1 5.1 13	91 56 2.0 5.2 13	5 3 0.11 0.26 0.3	6 5 5 2
C3- Chrysenes C4- Chrysenes Benzo (b) Fluoranthene Benzo (k) Fluoranthene Benzo (e) Pyrene Benzo (a) Pyrene	101 116 92 59 1.9 5.5 13 2.5	114 95 55 2.1 5.0 13 2.4	107 85 54 2.1 5.1 13 2.5	91 56 2.0 5.2 13 2.5	5 3 0.11 0.26 0.3 0.07	6 5 5 2 3
C3- Chrysenes C4- Chrysenes Benzo (b) Fluoranthene Benzo (k) Fluoranthene Benzo (e) Pyrene Benzo (a) Pyrene Perylene	101 116 92 59 1.9 5.5 13 2.5 0.8	114 95 55 2.1 5.0 13	107 85 54 2.1 5.1 13	91 56 2.0 5.2 13	5 3 0.11 0.26 0.3	6 5 5 2
C3- Chrysenes C4- Chrysenes Benzo (b) Fluoranthene Benzo (k) Fluoranthene Benzo (e) Pyrene Benzo (a) Pyrene	101 116 92 59 1.9 5.5 13 2.5 0.8 € 0.25	114 95 55 2.1 5.0 13 2.4 0.9	107 85 54 2.1 5.1 13 2.5 1.0	91 56 2.0 5.2 13 2.5 0.9	5 3 0.11 0.26 0.3 0.07 0.07	6 5 6 5 2 3 8
C3- Chrysenes C4- Chrysenes Benzo (b) Fluoranthene Benzo (k) Fluoranthene Benzo (e) Pyrene Benzo (a) Pyrene Perylene Indeno (1,2,3 - cd) Pyrene	101 116 92 59 1.9 5.5 13 2.5 0.8 0.25 0.67	114 95 55 2.1 5.0 13 2.4 0.9 0.28	107 85 54 2.1 5.1 13 2.5 1.0 0.26	91 56 2.0 5.2 13 2.5 0.9 0.26	5 3 0.11 0.26 0.3 0.07 0.07 0.07	6 5 5 2 3 8 6

Alkane Analyte: nC-10 Decane nC-11 Undecane nC-12 Dodecane	Replicate #1			1		
nC-11 Undecane nC-12 Dodecane		Replicate #2 Conc (ng/mg)	Replicate #3 Conc (ng/mg)	Average	Stdev	%RSD
nC-11 Undecane nC-12 Dodecane	Conc (ng/mg) 0	0	0	Average 0	0.0	0
nC-12 Dodecane	65	65	68	66	1.7	3
1	132	132	128	131	2.3	2
nC-13 Tridecane	145	145	155	148	5.8	4
nC-14 Tetradecane	149	158	178	162	14.6	9
nC-15 Pentadecane	1/1	1/9	192	181	11	6
nC-16 Hexadecane	207	221	235	221	14	6
nC-17 Heptadecane	251	236	241	243	7	3
Pristane	1648	1624	1621	1631	15	1
nC-18 Octadecane	200	192	189	194	6	3
Phylane	1760	1728	1728	1739	19	1
nC-19 Nonadecane	292	315	310	316	21.0	8
nC-20 Eicosane	227	215	215	219	7	3
nC-21 Heneicosane	151	125	142	139	13	9
nC-22 Docosane	136	142	140	139	3	2
nC-23 Tricosane	117	149	153	140	20	14
nC-24 Tetracosane	129	122	125	125	4	3
nC-25 Pentacosane	230	229	227	229	2	1
nC-26 Hexacosane	198	188	178	188	10	5
nC-27 Heptacosane	143	143	144	143	1	ŏ
nC 28 Octacosano	147	144	140	144	4	3
nC-29 Nonacosane	148	146	154	149	4	3
nC-30 Triacontane	118	140	154	132	18	13
nC-31 Hentriacontane	113	121	111	115	5	5
nC-32 Dotriacontane	113	135	130	126	11	9
nC-33 Tritriacontane	91	100	93	95	5	5
nC-34 Tetratriacontane	51	62	49	54	7	13
nC-35 Pentatriacontane	41	45	39	42	3	1
Total Alkanes	7175	7188	7267	7210	50	1
Anomalia Analuta	Construction a)	Conc (ng/mg)	Constructions)	A	Stdev	%RSD
Aromatic Analyte:	Conc (ng/mg) ()	O Conc (ng/mg)	Conc (ng/mg) 0	Average 0	0.0	0
Naphthalene	-					
C1-Naphthalenes	115 959	99 1034	109 10/3	108 1022	8 58	06
C2-Naphthalenes C3-Naphthalenes	1140	1131	1141	1138	6	0
C4-Naphthalenes	948	912	949	936	21	2
E4-Naphthalenes	28	29	35	31	4	12
C1 Fluorenes	322	345	316	328	4 16	5
C1 Fluorenes	407	405	420	411	8	2
C3- Fluorenes	316	344	353	338	19	6
Dibenzothiophene	99	90	97	95	5	5
C1-Dibenzothiophenes	305	343	332	327	20	6
C2-Dibenzothiophenes	680	639	664	661	20	3
C3- Dibenzothiophenes	431	453	461	448	16	4
Phenanthrene	68	71	81	73	7	10
C1-Phenanthrenes	441	422	433	432	10	2
C2-Phenanthrenes	761	678	729	723	42	6
C3-Phenanthrenes	565	568	585	572	11	2
C4-Phenanthrenes	219	222	239	226	11	5
Anthracene	3.0	3.1	2.8	3.0	0.16	0
Fluoranthene	2.2	2.1	2.0	2.1	0.06	3
Pyrene	1.8	8.4	10.1	8.8	1.19	14
C1- Pyrenes	121	155	145	140	17.1	12
C2- Pyrenes	167	181	184	1//	9	5
C3 Pyrenes	153	175	195	174	21	12
C4- Pyrenes	91	92	113	99	12.7	13
Naphthobenzothiophene	100	92	97	97	4.0	4
C-1 Naphthobenzothiophenes	223	229	237	229	7	3
C-2 Naphthobenzothiophenes	251	255	264	257	7	3
C-3 Naphthobenzothiophenes	173	177	193	181	11	6
Benzo (a) Anthracene	9.3	7.5	7.2	8.0	1.14	14
Chrysene	57	56	63	58	4	6
C1- Chrysenes	153	132	185	157	27	17
C2- Chrysenes	182	184	203	190	11	6
C3- Chrysenes	142	150	163	152	10.2	7
C4- Chrysenes	107	90	121	106	16	15
Benzo (b) Fluoranthene	0.9	1.0	0.9	0.9	0.06	6
Benzo (k) Fluoranthene	1.2	5.0	1.9	4.7	0.45	10
Benzo (e) Pyrene	9	10	12	11	1.29	12
Benzo (a) Pyrene	2.0	2.3	2.2	2.2	0.13	6
Perylene	0.3	0.3	0.3	0.3	0.03	11
Indeno (1,2,3 - cd) Pyrene	0.22	0.17	0.16	0.18	0.03	0
Dibenzo (a,h) anthracene	0.26	0.26	0.28	0.27	0.01	0
	0.2	0.3	0.4	0.3	0.07	0
Benzo (g,h,i) perylene					·····	r

STATIST	SARVA PROE ICS FOR SURROGATI Testing Date:	E CORRECTED DATA				
Alkane Analyte:	Replicate #1	Replicate #2	Replicate #3		-	
-	Conc (ng/mg)	Conc (ng/mg)	Conc (ng/mg)	Average	Stdev	%RSD
nC-10 Decane	0	0	0	0	0.0	0
nC-11 Undecane	0	0	0	0	0.0	0
nC-12 Dodecane	0	0	0	0	0.0	0
nC-13 Tridecane	0	0	0	0	0.0	0
nC-14 Tetradecane	0	0	0	0	0.00	0
nC-15 Pentadecane	0	0	0	0	0	0
nC-16 Hexadecane	0	0	0	0	0.0	0
nC-17 Heptadecane	31	32	32	32	1	2
Pristane	391	423	431	415	21	5
nC-18 Octadecane	13	15	16	15	1.2	8
Phylane	380	380	362	374	10	3
nC-19 Nonadecane	13	12	11	12	1.0	8
nC-20 Eicosane	9	9	8	9	0.5	6
nC-21 Heneicosane	13	13	12	13	0.3	2
nC-22 Docosane	14	14	13	14	0.6	4
nC-23 Tricosane	14	15	15	14	0.4	3
nC-24 Tetracosane	21	21	20	21	0.7	4
nC-25 Pentacosane	129	122	112	121	8.66	7
nC-26 Hexacosane	98	100	93	97	4.0	4
nC-27 I leptacosane	42	38	35	38	3.3	9
nC 28 Octacosane	30	30	29	30	0.6	2
nC-29 Nonacosane	46	45	42	44	2.0	5
nC-30 Triacontane	32	27	22	27	4.8	18
nC-31 Hentriacontane	15	16	16	16	0.5	3
nC-32 Dotriacontane	8	9	10	9	1.0	12
nC-33 Tritriacontane	10	11	10	10	0.2	2
nC-34 Tetratriacontane	0	0	0	0	0.0	Ō
nC-35 Pentatriacontane	0.00	0.00	0.00	0.00	0.00	Ŭ
Total Alkanes	1308	1333	1289	1310	22	2
Aromatic Analyte:	Conc (ng/mg)	Conc (ng/mg)	Conc (ng/mg)	Average	Stdev	%RSD
Naphthalene	0.00	0.00	0.00	0.00	0.00	0
C1-Naphthalenes	0.00	0.00	0.00	0.00	0.00	0
C2-Naphthalenes	288	293	287	289	2.8	1
C3-Naphthalenes	1092	1047	975	1038	59	6
C4-Naphthalenes	747	798	820	788	37	5
Fluorene	16.5	16	14.4	15.5	1.1	1
C1 Fluorenes	131	133	131	132	1	
C2-Fluorenes	304	357	393	352	45	13
C3- Fluorenes	249	286	311	282	31	11
Dibenzothiophene	18.3	18	16.9	18	0.7	4
C1-Dibenzothiophenes	246	233	214	231	16	7
C2-Dibenzothiophenes	553	585	595	578	22	4
C3- Dibenzothiophenes	350	367	372	363	12	3
Phenanthrene	18	18	16	17	1.4	8
C1-Phenanthrenes	350	345	328	341	12	3
C2-Phenanthrenes	606	681	728	672	62	9
C3-Phenanthrenes	458	510	542	503	42	8
C4-Phenanthrenes	185	203	213	200	13.79	7
Anthracene	2.0	2.1	2.0	2.0	0.02	
Fluoranthene	1.8	1.8	1.8	1.8	0.02	
Providence	8.2	9.2	9.8	9.1	0.02	9
C1- Pyrenes	97	95	91	94	3.0	3
C2- Pyrenes	136	134	129	133	3.0	3
C3 Pyrenes	136	134	121	129	7.3	6
C4- Pyrenes	82	74	65	74	8.4	11
Naphthobenzothiophene	55	62	66	61	5.7	9
C-1 Naphthobenzothiophenes	189	198	200	196	6	3
C-2 Naphthobenzothiophenes	233	223	200	221	13	6
C-3 Naphthobenzothiophenes	133	134	131	133	15	1
Benzo (a) Anthracene	133	10.8	10.3	10.7	0.36	3
Chrysene	64	70	73	69	4.4	6
C1- Chrysenes	118	132	140	130	11.2	9
C2- Chrysenes	141	142	139	141	2	1
C2- Chrysenes C3- Chrysenes	141	142	139	141	10.3	8
C3- Chrysenes C4- Chrysenes		74				6
64- Chrysenes Benzo (b) Fluoranthene	69 0.9	0.9	77 0.8	73 0.9	4.4 0.08	9
Benzo (k) Fluoranthene Benzo (k) Fluoranthene	1.0	3.9	3.7	3.9	0.08	3
	1.0	12	12			5
	11		12	11	0.5	
Benzo (e) Pyrene Benzo (a) Pyrene	11 2.0		20	20	0.02	
Benzo (a) Pyrene	2.0	2.0	2.0	2.0	0.02	1
Benzo (a) Pyrene Perylene	2.0 0.73	2.0 0.//	0.78	0.76	0.02	3
Benzo (a) Pyrene Perylene Indeno (1,2,3 - cd) Pyrene	2.0 0.73 • 0.00	2.0 0.77 0.00	0.78 0.00	0.76 0.00	0.02 0.00	3 0
Benzo (a) Pyrene Perylene	2.0 0.73	2.0 0.//	0.78	0.76	0.02	3

APPENDIX II

General Linear Model: Alkanes versus Day, Treatment

Day f	Type Levels V ixed 3 ixed 3 C	0 7 28	utrient	Sarva		
Analysis of	Variance for	Alkanes,	using A	Adjusted SS	for Test	s
Day*Treatme Error Total	2 504690 2 142538 n 4 213292 18 4803	4846 5046 4113 1425 4217 2132 1825 48 5001	904846 2 384113	533231054	945.67 267.08	0.000
Comparisons Day =	riable Alkane with Control O Control subtr	Level	m:			
Level	Differe	nce	SE of		Adjusted	
Day*Treatme O Nutrien O Sarva 7 Control 7 Nutrien 7 Sarva 28 Control 28 Nutrien	n of Me t -2 3 -1 t -8 -5 -3	ans Diff 000 100 357 110 727 496 753	erence 1334 1334 1334 1334 1334 1334 1334		P-Value 0.5604 0.1598 0.8693 0.0001 0.0029 0.0924	

General Linear Model: Alkanes versus Day, Treatment

-47653

1334

-35.73 0.0000

28 Sarva

Type Levels Values Factor fixed 3 0 7 28 Day 3 Control Nutrient Sarva Treatmen fixed Analysis of Variance for Alkanes, using Adjusted SS for Tests Seq SS Adj SS Adj MS F \mathbf{P} Source DFDay 2 5046904846 5046904846 2523452423 945.67 0.000 2 1425384113 1425384113 712692056 267.08 0.000 Treatmen Day*Treatmen 4 2132924217 2132924217 533231054 199.83 0.000 18 48031825 48031825 2668435 Error Total 26 8653245001 Dunnett Simultaneous Tests Response Variable Alkanes Comparisons with Control Level Day = 7Treatmen = Control subtracted from: SE of Adjusted Difference Level of Means Difference T-Value Day*Treatmen P-Value 0.8693 1.02 0 Control 1357 1334 0 Nutrient -643 1334 -0.48 0.9978 4457 0 Sarva 3.34 0.0218 1334 7 Nutrient -6753 1334 -5.06 0.0006 0.0249 7 Sarva -4370 1334 -3.28 0.4921 1334 -1.60 28 Control -2139 1334 -30.29 0.0000 28 Nutrient -40397 28 Sarva -46297 1334 -34.71 0.0000

General Linear Model: Alkanes versus Day, Treatment

Type Levels Values Factor Day fixed 3 0 7 28 Treatmen fixed 3 Control Nutrient Sarva Analysis of Variance for Alkanes, using Adjusted SS for Tests DF Source Seq SS Adj SS Adj MS E Ρ
 Day
 2 5046904846 5046904846 2523452423
 945.67
 0.000

 Treatmen
 2 1425384113 1425384113
 712692056 267.08
 0.000

 Day*Treatmen
 4 2132924217 2132924217
 533231054 199.83
 0.000
 Error 18 48031825 Total 26 8653245001 18 48031825 48031825 2668435 Dunnett Simultaneous Tests Response Variable Alkanes Comparisons with Control Level Day = 28 Treatmen = Control subtracted from: Difference SE of Adjusted Level of Means Difference T-Value P-Value Day*Treatmen Day-TreatmenOf MeansDifferenceT-ValueP-Value0Control349613342.620.09240Nutrient149613341.120.81100Sarva659613344.950.00077Control213913341.600.49217Nutrient-46141334-3.460.01707Sarva-22311334-1.670.449428Nutrient-382571334-28.680.000028Sarva-441571334-33.110.0000)ay*Treatmen O Control O Nutrient

General Linear Model: Rank_Alkanes versus Day, Treatment

Day			307	28	rient	Sarva	÷	
Anal	ysis of Va	ariance	for Ran	k_Alk,	using	Adjusted S	S for Te	sts
Sour	ce	DF	Seq SS	Ac	tj SS	Adj MS	F	P
Day		2	904.22	90	4.22	452.11	68.97	0.000
Trea	tmen	2	349.56	34	9.56	174.78	26.66	0.000
Day*	Treatmen	4	266.22	26	6.22	66.56	10.15	0.000
Erro	r	18	118.00	11	8.00	6.56		
Tota	l	26	1638.00					
Resp Comp Day	ett Simuli onse Varia arisons wi = 0 tmen = Con	able Ran ith Cont	k_Alk rol Lev					
Leve	1	Diff	erence	5	E of		Adjuste	d
Day*	Treatmen	of	Means	Differ	ence	T-Value	P-Valu	e
0	Nutrient		-4.00			-1.913		8
0	Sarva		4.33	2	.091	2.073	0.245	7
	Control		-1.00		.091			
	Nutrient		-12.67			-6.059		
	Sarva		-10.67			-5.102		
	Control		-6.33			-3.030		-
	Nutrient		-16.33			-7.813		
28	Sarva		-19.33	2	.091	-9.248	0.000	0

General Linear Model: Rank_Alkanes versus Day, Treatment

Type Levels Values Factor Day fixed 3 0 7 28 Treatmen fixed 3 Control Nutrient Sarva Analysis of Variance for Rank_Alk, using Adjusted SS for Tests
 Source
 DF
 Seq SS
 Adj SS
 Adj MS
 F
 P

 Day
 2
 904.22
 904.22
 452.11
 68.97
 0.000

 Treatmen
 2
 349.56
 349.56
 174.70
 000.000
 Treatmen2349.56349.56174.7826.660.000Day*Treatmen4266.22266.2266.5610.150.000Error18118.00118.006.56Total261638.00 Dunnett Simultaneous Tests Response Variable Rank Alk Comparisons with Control Level Day = 7Treatmen = Control subtracted from: Difference SE of Level Adjusted of Means Difference T-Value P-Value Day*Treatmen 0 Control 0 Nutrient 0 Sarva 7 Nutrient 7 Sarva 28 Control 28 Nutrient 28 Sarva 1.00 2.091 0.478 0.9979

General Linear Model: Rank_Alkanes versus Day, Treatment

Day	fi	xed	ls Values 3 0 7 2 3 Control	28 1 Nutrient	Sarva	
Ana	lysis of	Variance	for Rank	Alk, using	Adjusted SS	for Tests
Sou	rce	DF	Seq SS	Adj SS		
Day		2	904.22	904.22	452.11	68.97 0.000
Tre	atmen	2	349.56	349.56	174.78	26.66 0.000
Day	*Treatmen	. 4	266.22	266.22	66.56	10.15 0.000
Err	or	18	118.00	118.00	6.56	
Tot	al	26	1638.00			
Res Com Day	= 2	iable Ran with Cont 8				
Lev	el	Dif	ference	SE of		Adjusted
Day	*Treatmen	0	f Means I	Difference	T-Value	P-Value
0	Control		6.33	2.091	3.030	0.0413
0	Nutrient		2.33	2.091	1.116	0.8143
0	Sarva		10.67		5,102	0.0005
	Control		5.33		2.551	
7	Nutrient		-6.33		-3.030	
	Sarva		-4.33		-2.073	
	Nutrient		-10.00		-4.783	
28	Sarva		-13.00	2.091	-6.218	0.0001

General Linear Model: Aromatics versus Day, Treatment

Factor Type Levels Values Day fixed 3 0 7 28 Treatmen fixed 3 Control Nutrient Sarva Analysis of Variance for Aromatic, using Adjusted SS for Tests Source DF Seq SS Adj SS Adj MS F
 Day
 2
 68510796
 68510796
 34255398
 114.47
 0.000

 Treatmen
 2
 10785102
 10785102
 5392551
 18.02
 0.000

 Day*Treatmen
 4
 29262580
 29262580
 7315645
 24.45
 0.000

 Error
 18
 5386510
 5386510
 299251
 Error 18 5386510 Total 26 113944989 Dunnett Simultaneous Tests Response Variable Aromatic Comparisons with Control Level Day = 0Treatmen = Control subtracted from: Difference SE of Adjusted LevelDifferenceSE ofAdjustedDay*Treatmenof MeansDifferenceT-ValueP-Value0Nutrient295446.70.660.98430Sarva659446.71.480.57667Control-2221446.7-4.970.00077Nutrient-2155446.7-4.820.00097Sarva-2407446.7-5.390.000328Control-924446.7-7.960.000028Sarva-6041446.7-13.530.0000 Level

Ρ

General Linear Model: Aromatics versus Day, Treatment

±.	pe Levels Value ed 3 0 7 ed 3 Contr		Sarva	
Analysis of V	ariance for Aro	matic, using	Adjusted S	S for Tests
Error Total Dunnett Simul Response Vari Comparisons w Day = 7	2 10785102 4 29262580 18 5386510 26 113944989 taneous Tests	68510796 10785102 29262580 5386510	34255398 5392551 7315645	114.47 0.000 18.02 0.000
Level Day*Treatmen 0 Control 0 Nutrient 0 Sarva 7 Nutrient 7 Sarva 28 Control 28 Nutrient 28 Sarva	Difference of Means 2221 2516 2880 66 -186 1297 -1335 -3820	Difference 446.7 446.7 446.7 446.7 446.7 446.7 446.7 446.7	T-Value 4.973 5.634 6.449 0.149 -0.416 2.905 -2.989	0.0002 0.0000 1.0000 0.9992 0.0532 0.0449

General Linear Model: Aromatics versus Day, Treatment

Factor Type Levels Values fixed 3 0 7 28 Day Treatmen fixed 3 Control Nutrient Sarva Analysis of Variance for Aromatic, using Adjusted SS for Tests SourceDFSeq SSAdj SSAdj MSFPDay2685107966851079634255398114.470.000Treatmen21078510210785102539255118.020.000Day*Treatmen42926258029262580731564524.450.000Error18538651053865102992511000Total26113944989113944989113944989 Dunnett Simultaneous Tests Response Variable Aromatic Comparisons with Control Level Day = 28 Treatmen = Control subtracted from: Difference SE of Adjusted Level LevelDifferenceSE ofAdjustedDay*Treatmenof MeansDifferenceT-ValueP-Value0Control924446.72.070.24740Nutrient1219446.72.730.07510Sarva1583446.73.540.01427Control-1297446.7-2.900.05327Nutrient-1231446.7-2.760.07137Sarva-1483446.7-3.320.022728Nutrient-2632446.7-5.890.000128Sarva-5117446.7-11.460.0000

General Linear Model: Aromatics versus Day, Treatment

Factor Typ Day fixe Treatmen fixe		s 28 ol Nutrient	Sarva					
Analysis of Va	riance for Arc	matic, using	Adjusted S	S for Tests				
Source Day Treatmen Day*Treatmen Error Total	DF Seq SS 2 68510796 2 10785102 4 29262580 18 5386510 26 113944989	68510796 10785102 29262580 5386510	Adj MS 34255398 5392551 7315645 299251	18.02 0.000				
Tukey Simultan Response Varia All Pairwise C	ble Aromatic	ng Levels of	Day*Treatm	nen				
Day = 0 Treatmen = Con	trol subtracte	d from:						
Level Day*Treatmen 0 Nutrient 0 Sarva 7 Control 7 Nutrient 7 Sarva 28 Control 28 Nutrient 28 Sarva	Difference of Means 295 659 -2221 -2155 -2407 -924 -3556 -6041	SE of Difference 446.7 446.7 446.7 446.7 446.7 446.7 446.7 446.7	T-Value 0.66 1.48 -4.97 -4.82 -5.39 -2.07 -7.96 -13.53	Adjusted P-Value 0.9988 0.8527 0.0025 0.0034 0.0011 0.5207 0.0000 0.0000				
Day = 0 Treatmen = Nut:	rient subtract	ed from:						
Level Day*Treatmen 0 Sarva 7 Control 7 Nutrient 7 Sarva 28 Control 28 Nutrient 28 Sarva	Difference of Means 364 -2516 -2450 -2702 -1219 -3851 -6336	SE of Difference 446.7 446.7 446.7 446.7 446.7 446.7 446.7	T-Value 0.81 -5.63 -5.49 -6.05 -2.73 -8.62 -14.19	Adjusted P-Value 0.9948 0.0007 0.0009 0.0003 0.2049 0.0000 0.0000				
Day = 0 Treatmen = Sarva subtracted from:								
Level Day*Treatmen 7 Control 7 Nutrient 7 Sarva 28 Control 28 Nutrient 28 Sarva	Difference of Means -2880 -2814 -3066 -1583 -4215 -6700	SE of Difference 446.7 446.7 446.7 446.7 446.7 446.7	T-Value -6.45 -6.30 -6.86 -3.54 -9.44 -15.00	Adjusted P-Value 0.0001 0.0002 0.0001 0.0462 0.0000 0.0000				
Day = 7 Treatmen = Con	trol subtracte	d from:						
Level Day*Treatmen 7 Nutrient 7 Sarva	Difference of Means 66 -186	SE of Difference 446.7 446.7	T-Value 0.149 -0.416	Adjusted P-Value 1.0000 1.0000				

28 Control 28 Nutrient 28 Sarva	1297 -1335 -3820	446.7 446.7 446.7	-2.989	0.1523 0.1313 0.0000					
Day = 7 Treatmen = Nutr	ient subtract	ed from:							
Level Day*Treatmen 7 Sarva 28 Control 28 Nutrient 28 Sarva	Difference of Means -252 1231 -1401 -3886	Difference 446.7 446.7 446.7	-0.564 2.756 -3.137	0.9996 0.1960 0.1004					
Day = 7 Treatmen = Sarva	a subtracted	from:							
Level Day*Treatmen 28 Control 28 Nutrient 28 Sarva	1483 -1149	Difference 446.7	3.320 -2.573	0.0713 0.2629					
Day = 28 Treatmen = Control subtracted from:									
Level Day*Treatmen 28 Nutrient 28 Sarva		Difference 446.7	T-Value -5.89	Adjusted P~Value 0.0004 0.0000					
Day = 28 Treatmen = Nutr:	ient subtract	ed from:							
Level Day*Treatmen 28 Sarva	Difference of Means -2485	Difference	T-Value -5,564						

General Linear Model: Alkanes versus Day, Treatment

Factor Type Levels Values fixed 3 0 7 28 Day Treatmen fixed 3 Control Nutrient Sarva Analysis of Variance for Alkanes, using Adjusted SS for Tests
 Source
 DF
 Seq SS
 Adj SS
 Adj MS
 F
 P

 Day
 2
 5046904846
 5046904846
 2523452423
 945.67
 0.000

 Treatmen
 2
 1425384113
 1425384113
 712692056
 267.08
 0.000
 2 1425384113 1425384113 712692056 267.08 0.000 Day*Treatmen 4 2132924217 2132924217 533231054 199.83 0.000 Error 18 48031825 48031825 2668435 Total 26 8653245001 Tukey Simultaneous Tests Response Variable Alkanes All Pairwise Comparisons among Levels of Day*Treatmen Dav = 0Treatmen = Control subtracted from: Adjusted LevelDifferenceSE ofAdjustedDay*Treatmenof MeansDifferenceT-ValueP-Value0Nutrient-20001334-1.500.84190Sarva310013342.320.37807Control-13571334-1.020.97907Nutrient-81101334-6.080.00037Sarva-57271334-4.290.010128Control-34961334-2.620.243928Nutrient-417531334-31.300.000028Sarva-476531334-35.730.0000 Difference SE of Level Dav = 0Treatmen = Nutrient subtracted from: Adjusted Difference SE of Level
 Day*Treatmen
 of Means
 Difference
 T-Value
 P-Value

 0
 Sarva
 5100
 1334
 3.82
 0.0265

 7
 Control
 643
 1334
 0.48
 0.9999

 7
 Nutrient
 -6110
 1334
 -4.58
 0.0056

 7
 Sarva
 -3727
 1334
 -2.79
 0.1840
 7 Control 7 Nutrient 7 Sarva 28 Control 28 Nutrient 28 Sarva Day = 0Treatmen = Sarva subtracted from: Difference SE of Adjusted Level Day*Treatmen of Means Difference T-Value P-Value 7 Control -4457 1334 -3.34 0.0685
 Day*Treatmen
 OI Means
 Difference
 T-value
 P-value

 7
 Control
 -4457
 1334
 -3.34
 0.0685

 7
 Nutrient
 -11210
 1334
 -8.40
 0.0000

 7
 Sarva
 -8827
 1334
 -6.62
 0.0001

 28
 Control
 -6596
 1334
 -4.95
 0.0026

 28
 Nutrient
 -44853
 1334
 -33.63
 0.0000

 28
 Sarva
 -50753
 1334
 -38.05
 0.0000
 Day = 7Treatmen = Control subtracted from: LevelDifferenceSE ofAdjustedDay*Treatmenof MeansDifferenceT-ValueP-Value7Nutrient-67531334-5.060.00217Sarva-43701334-3.280.0775

28 Control 28 Nutrient 28 Sarva	-2139 -40397 -46297	1334 1334 1334	-30.29	
Day = 7 Treatmen = Nutr	ient subtract	ed from:		
Level Day*Treatmen 7 Sarva 28 Control 28 Nutrient 28 Sarva Day = 7	Difference of Means 2383 4614 -33643 -39543	Difference 1334 1334	T-Value 1.79 3.46 -25.22 -29.65	0.0546 0.0000
Treatmen = Sarv	a subtracted	from:		
Level Day*Treatmen 28 Control 28 Nutrient 28 Sarva	Difference of Means 2231 -36027 -41927	Difference 1334	T-Value 1.67 -27.01 -31.43	Adjusted P-Value 0.7547 0.0000 0.0000
Day = 28 Treatmen = Cont	rol subtracte	d from:		
Level Day*Treatmen 28 Nutrient 28 Sarva	Difference of Means -38257 -44157	Difference	T-Value -28.68 -33.11	Adjusted P-Value 0.0000 0.0000
Day = 28 Treatmen = Nutr	ient subtract	ed from:		
Level Day*Treatmen 28 Sarva	Difference of Means -5900	SE of Difference 1334	T-Value -4.424	Adjusted P-Value 0.0078

General Linear Model: Rank_Alkanes versus Day, Treatment

Factor Ty Day fix Treatmen fix		28	Sarva	
Analysis of V	ariance for Ran	k_Alk, using	Adjusted S	SS for Tests
Source Day Treatmen Day*Treatmen Error Total	DF Seq SS 2 904.22 2 349.56 4 266.22 18 118.00 26 1638.00	904.22 349.56 266.22 118.00	Adj MS 452.11 174.78 66.56 6.56	F P 68.97 0.000 26.66 0.000 10.15 0.000
	neous Tests able Rank_Alk Comparisons amo	ong Levels of	Day*Treatm	nen
Day = 0 Treatmen = Co	ntrol subtracte	d from:		
Level Day*Treatmen 0 Nutrient 0 Sarva 7 Control 7 Nutrient 7 Sarva 28 Control 28 Nutrient 28 Sarva	Difference of Means -4.00 4.33 -1.00 -12.67 -10.67 -6.33 -16.33 -19.33	SE of Difference 2.091 2.091 2.091 2.091 2.091 2.091 2.091 2.091	T-Value -1.913 2.073 -0.478 -6.059 -5.102 -3.030 -7.813 -9.248	Adjusted P-Value 0.6140 0.5183 0.9999 0.0003 0.0019 0.1221 0.0000 0.0000
Day = 0 Treatmen = Nu	trient subtract	ed from:		
Level Day*Treatmen 0 Sarva 7 Control 7 Nutrient 7 Sarva 28 Control 28 Nutrient 28 Sarva	Difference of Means 8.33 3.00 -8.67 -6.67 -2.33 -12.33 -15.33	SE of Difference 2.091 2.091 2.091 2.091 2.091 2.091 2.091	T-Value 3.986 1.435 -4.146 -3.189 -1.116 -5.900 -7.335	Adjusted P-Value 0.0191 0.8699 0.0138 0.0912 0.9639 0.0004 0.0000
Day = 0 Treatmen = Sa	rva subtracted	from:		
Level Day*Treatmen 7 Control 7 Nutrient 7 Sarva 28 Control 28 Nutrient 28 Sarva	Difference of Means -5.33 -17.00 -15.00 -10.67 -20.67 -23.67	SE of Difference 2.091 2.091 2.091 2.091 2.091 2.091	T-Value -2.55 -8.13 -7.18 -5.10 -9.89 -11.32	Adjusted P-Value 0.2719 0.0000 0.0001 0.0019 0.0000 0.0000
Day = 7 Treatmen = Control subtracted from:				
Level Day*Treatmen 7 Nutrient 7 Sarva	Difference of Means -11.67 -9.67	SE of Difference 2.091 2.091	T-Value -5.581 ~4.624	Adjusted P-Value 0.0007 0.0051

28 Control 28 Nutrient 28 Sarva	-5.33 -15.33 -18.33	2.091 2.091 2.091	-7.335	
Day = 7 Treatmen = Nutr	ient subtract	ed from:		
Level Day*Treatmen 7 Sarva 28 Control 28 Nutrient 28 Sarva	Difference of Means 2.000 6.333 -3.667 ~6.667	Difference 2.091 2.091 2.091		
Day = 7 Treatmen = Sarv	a subtracted	from:		
Level Day*Treatmen 28 Control 28 Nutrient 28 Sarva	Difference of Means 4.333 -5.667 -8.667	Difference 2.091	T-Value 2.073 -2.711 -4.146	Adjusted P-Value 0.5183 0.2112 0.0138
Day = 28 Treatmen = Cont	rol subtracte	d from:		
Level Day*Treatmen 28 Nutrient 28 Sarva	Difference of Means -10.00 -13.00	Difference	T-Value -4.783 -6.218	Adjusted P-Value 0.0037 0.0002
Day = 28 Treatmen = Nutr	ient subtract	ed from:		
Level Day*Treatmen 28 Sarva	Difference of Means -3.000	SE of Difference 2.091	T-Value -1.435	Adjusted P-Value 0.8699

APPENDIX III

*	* * * * * * * * * * * * * * * * * * * *	۰*			
*	MOST PROBABLE NUMBER CALCULATOR	*			
*	Version 4.04	*			
*	UNITED STATES ENVIRONMENTAL PROTECTION AGENCY	*			
*	RISK REDUCTION ENGINEERING LABORATORY	*			
*	CINCINNATI, OHIO	*			
*	* * * * * * * * * * * * * * * * * * * *				

DATA SET SARVA NUMBER OF DILUTIO	•	Day 0
VOLUME, ml	# TUBES	# POSITIVE TUBES
.1	24	24
.01	24	24
.001	24	24
.0001	24	10
.00001	24	4
.10000E-05	24	0
.10000E-06	24	0
.10000E-07	24	0
.10000E-08	24	0
.10000E-09	24	0

MOST PROBABLE NUMBER, MPN (per ml) = 6,794 MPN CORRECTED FOR BIAS (Salama) = 6,526 SPEARMAN-KARBER ESTIMATE = 12,115 LOWER 95% CONFIDENCE LIMIT (Cornish & Fisher) = 3,958

UPPER	95%	CONFIDENCE	LIMIT	(Cornish	&	Fisher)	=	10,808
OTTDIC	200	COLL TDRUCH		(COTITIOII	~	I IDIICI /		±0,000

DATA SET SAR NUMBER OF DILU		Day U		
VOLUME, ml	# TUBES	# POSITIVE TUBES		
.1	24	24		
.01	24	24		
.001	24	24		
.0001	24	9		
.00001	24	4		
.10000E-05	24	0		
.10000E-06	24	0		
.10000E-07	24	0		
.10000E-08	24	0		
.10000E-09	24	0		
MOS	T PROBABLE NU	MBER,,MPN (per ml)	= 6,195	
MPN	CORRECTED FC	R BIAS (Salama)	= 5,912	
		ESTIMATE		
LOWER 95% CONF	IDENCE LIMIT	(Cornish & Fisher)	 = 3,648	
		(Cornish & Fisher)	-	

DATA SET 1 SA	RVA Control	3, Day 0
NUMBER OF DILUT	IONS = 10	
VOLUME, ml	# TUBES	# POSITIVE TUBES
.1	24	24
.01	24	24
.001	24	24
.0001	24	10
.00001	24	3
.10000E-05	24	0
.10000E-06	24	0
.10000E-07	24	0
.10000E-08	24	0
.10000E-09	24	0

MOST PROBABLE NUMBER, MPN (per ml) = 6,330 MPN CORRECTED FOR BIAS (Salama) = 6,051 SPEARMAN-KARBER ESTIMATE = 11,007 LOWER 95% CONFIDENCE LIMIT (Cornish & Fisher) = 3,718 UPPER 95% CONFIDENCE LIMIT (Cornish & Fisher) = 10,149

/OLUME, ml	# TUBES	# POSITIVE TUBES		
1				
.1	24	24		
.01	24	24		
.001	24	24		
.0001	24	12		
.00001	24	2		
10000E-05	24	0		
10000E-06	24	0		
10000E-07	24	0		
10000E-08	24	0		
.10000E-09	24	0		
M	OST PROBABLE NU	MBER,,MPN (per ml)	=	7.129
		R BIAS (Salama)		
		ESTIMATE		
اد				12,113
	VETDENCE LIMIT	(Cornish & Fisher)	=	4 132
		(Cornish & Fisher)		-

DATA SET SARVA		Day 7
NUMBER OF DILUTI VOLUME, ml		# POSITIVE TUBES
.1	24	24
.01	24	24
.001	24	24
.0001	24	11
.00001	24	3
.10000E-05	24	0
.10000E-06	24	0
.10000E-07	24	0
.10000E-08	24	0
.10000E-09	24	0

MOST PROBABLE NUMBER, MPN (per ml) = 6,956 MPN CORRECTED FOR BIAS (Salama) = 6,692 SPEARMAN-KARBER ESTIMATE = 12,115 LOWER 95% CONFIDENCE LIMIT (Cornish & Fisher) = 4,042 UPPER 95% CONFIDENCE LIMIT (Cornish & Fisher) = 11,038

	SARVA Control 3, LUTIONS = 10	Day 7		
VOLUME, ml	# TUBES	# POSITIVE	TUBES	
.1	24	24		
.01	24	24		
.001	24	24		
.0001	24	12		
.00001	24	3		
.10000E-05	24	0		
.10000E-06	24	0		
.10000E-07	24	0		
.10000E-08	24	0		
.10000E-09	24	0		
M	IOST PROBABLE NUM IPN CORRECTED FOR SPEARMAN-KARBER E	BIAS (Sala	.ma) =	7,393
	ONFIDENCE LIMIT (ONFIDENCE LIMIT (-

DATA SET SARVA NUMBER OF DILUTI	-	Day 28
VOLUME, ml		# POSITIVE TUBES
.1	24	24
.01	24	24
.001	24	24
.0001	24	13
.00001	24	5
.10000E-05	24	0
.10000E-06	24	0
.10000E-07	24	0
.10000E-08	24	0
.10000E-09	24	0

MOST PROBABLE NUMBER,,MPN (per ml) = 9,549 MPN CORRECTED FOR BIAS (Salama) = 9,282 SPEARMAN-KARBER ESTIMATE = 17,783 LOWER 95% CONFIDENCE LIMIT (Cornish & Fisher) = 5,472

UPPER 95% CONFIDENCE LIMIT (Cornish & Fisher) = 14,711

	SARVA Control 2, DILUTIONS = 10	Day 28	
		# POSITIVE TUBES	
.1	24	24	
.01	24	24	
.001	24	24	
.0001	24	12	
.00001	24	5	
.10000E-05	24	0	
.10000E-06	24	0	
.10000E-07	24	0	
.10000E-08	24	0	
.10000E-09	24	0	
	MPN CORRECTED FO	MBER,,MPN (per ml) R BIAS (Salama) ESTIMATE	= 8,470
		(Cornish & Fisher) (Cornish & Fisher)	•

DATA SET SAR NUMBER OF DILU	-	Day 28		
		# POSITIVE TUBE	S	
			-	
.1	24	24		
.01	24	24		
.001	24	24		
.0001	24	12		
.00001	24	б		
.10000E-05	24	0		
.10000E-06	24	0		
.10000E-07	24	0		
.10000E-08	24	0		
.10000E-09	24	0		
MOG		MDED NDN (non m	- X -	0 070
		MBER,,MPN (per m	,	
		R BIAS (Salama)		
SPE	ARMAN-KARBER	ESTIMATE	=	11,183
	TDENCE I IMTE	(Compich C Eicher		E 210
		(Cornish & Fisher		-
UPPER 95% CONF	IDENCE LIMIT	(Cornish & Fishe	r) =	14,327

*	* * * * * * * * * * * * * * * * * * * *	* *			
*	MOST PROBABLE NUMBER CALCULATOR	*			
*	Version 4.04	*			
*	UNITED STATES ENVIRONMENTAL PROTECTION AGENCY	*			
*	RISK REDUCTION ENGINEERING LABORATORY	*			
*	CINCINNATI, OHIO	*			
*	* * * * * * * * * * * * * * * * * * * *				

DATA SET SARVA NUMBER OF DILUTIO		Day 0
VOLUME, ml	# TUBES	# POSITIVE TUBES
.1	24	24
.01	24	24
.001	24	24
.0001	24	15
.00001	24	11
.10000E-05	24	0
.10000E-06	24	0
.10000E-07	24	0
.10000E-08	24	0
.10000E-09	24	0

MOST PROBABLE NUMBER,,MPN (per ml) = 15,565 MPN CORRECTED FOR BIAS (Salama) = 15,051 SPEARMAN-KARBER ESTIMATE = 38,312

LOWER	95%	CONFIDENCE	LIMIT	(Cornish	&	Fisher) =	9,482
UPPER	95%	CONFIDENCE	LIMIT	(Cornish	&	Fisher) =	23,556

	SARVA Product 2, DILUTIONS = 10	Day 0		
VOLUME, ml	# TUBES	# POSITIVE TUBES		
.1	24	24		
.01	24	24		
.001	24	24		
.0001	24	14		
.00001	24	13		
.10000E-05	5 24	0		
.10000E-06	5 24	0		
.10000E-07	24	0		
.10000E-08	3 24	0		
.10000E-09	24	0		
	MPN CORRECTED FOR	MBER,,MPN (per ml) R BIAS (Salama) ESTIMATE	=	15,165
		(Cornish & Fisher) (Cornish & Fisher)		•

	SARVA Product 3 DILUTIONS = 10	8, Day 0	
	# TUBES	# POSITIVE TUBES	
.1	24	24	
.01	24	24	
.001	24	24	
.0001	24	16	
.00001	24	11	
.10000E-05	24	0	
.10000E-06	24	0	
.10000E-07	24	0	
.10000E-08	24	0	
.10000E-09	24	0	
	MOST PROBABLE N	NUMBER,,MPN (per ml)	= 16,966
	MPN CORRECTED F	OR BIAS (Salama)	= 16,374

LOWER 95% CONFIDENCE LIMIT (Cornish & Fisher) = 10,433	EARMAN-KARBER ESTIMATE= 10,374= 42,170
UPPER 95% CONFIDENCE LIMIT (Cornish & Fisher) = 25,724	

	SARVA Product 1, DILUTIONS = 10	, Day 7	
		# POSITIVE TUBES	
.1	24	24	
.01	24	24	
.001	24	24	
.0001	24	24	
.00001	24	24	
.10000E-05	24	15	
.10000E-06	24	7	
.10000E-07	24	4	
.10000E-08	24	0	
.10000E-09	24	0	
	MOST PROBABLE N	JMBER,,MPN (per ml)	= 1,535,074
	MPN CORRECTED FO	DR BIAS (Salama)	= 1,485,041
	SPEARMAN-KARBER	ESTIMATE	= 3,831,187
		(Cornish & Fisher)	
UPPER 95% C	CONFIDENCE LIMIT	(Cornish & Fisher)	= 2,322,000

DATA SET S NUMBER OF DI	ARVA Product 2	, Day 7
		# POSITIVE TUBES
.1	24	24
.01	24	24
.001	24	24
.0001	24	24
.00001	24	24
.10000E-05	24	15
.10000E-06	24	б
.10000E-07	24	3
.10000E-08	24	0
.10000E-09	24	0
M	OSI PROBABLE N	UMBER,,MPN (per ml) =

	=	1,399,693 1,356,483 3,162,278
LOWER 95% CONFIDENCE LIMIT (Cornish & Fisher) UPPER 95% CONFIDENCE LIMIT (Cornish & Fisher)		

DATA SET SAN NUMBER OF DIL	RVA Product 3, UTIONS = 10	Day 7		
VOLUME, ml	# TUBES		JBES	
.1	24	24		
.01	24	24		
.001	24	24		
.0001	24	24		
.00001	24	24		
.10000E-05	24	14		
.10000E-06	24	10		
.10000E-07	24	2		
.10000E-08	24	0		
.10000E-09	24	0		
MPI	ST PROBABLE NU N CORRECTED FO EARMAN-KARBER	R BIAS (Salama	a) =	1,441,015
	FIDENCE LIMIT FIDENCE LIMIT			-

DATA SET SARVA	Product 1,	Day 28
NUMBER OF DILUTI	ONS = 10	
VOLUME, ml	# TUBES	# POSITIVE TUBES
.1	24	24
.01	24	24
.001	24	24
.0001	24	24
.00001	24	24
.10000E-05	24	24
.10000E-06	24	24
.10000E-07	24	12
.10000E-08	24	9
.10000E-09	24	5

MOST PROBABLE NUMBER,,MPN (per ml) MPN CORRECTED FOR BIAS (Salama) SPEARMAN-KARBER ESTIMATE	=	138,190,440 133,939,339 383,118,679
LOWER 95% CONFIDENCE LIMIT (Cornish & Fisher) UPPER 95% CONFIDENCE LIMIT (Cornish & Fisher)		

NUMBER OF 1	SARVA Product 2 DILUTIONS = 10	· •	
VOLUME, ml	# TUBES	# POSITIVE TUBES	3
.1	24	24	-
.01	24	24	
.001	24	24	
.0001	24	24	
.00001	24	24	
.10000E-05	24	24	
.10000E-06	24	24	
.10000E-07	24	13	
.10000E-08	24	8	
.10000E-09	24	7	
	MOST PROBABLE N	UMBER,,MPN (per m]	L) = 154,866,275
	MPN CORRECTED F	OR BIAS (Salama)	= 149,771,570
	SPEARMAN-KARBER	ESTIMATE	= 464,158,876

LOWER	95%	CONFIDENCE	LIMIT	(Cornish	&	Fisher)	=	94,282,566
				•				234,367,731
OFFER	0.00	COMPTDENCE	птытт	(COLIIISII	QL.	L'TPITET)	_	234,307,731

NUMBER OF	SARVA Product 3, DILUTIONS = 10	-	
VOLUME, MI	# TUBES	# POSITIVE TUBES	
.1	24	24	
.01	24	24	
.001	24	24	
.0001	24	24	
.00001	24	24	
.10000E-05	24	24	
.10000E-06	24	24	
.10000E-07		12	
.10000E-08		11	
.10000E-09	24	б	
	MPN CORRECTED FO	JMBER,,MPN (per ml) DR BIAS (Salama) ESTIMATE	= 151,227,253

LOWER	95%	CONFIDENCE	LIMIT	(Cornish	&	Fisher)	=	95,332,383
UPPER	95%	CONFIDENCE	LIMIT	(Cornish	&	Fisher)	=	236,720,299

DATA SET SARVA NUMBER OF DILUTIO	ONS = 10	-
VOLUME, ml	# IUBES	# POSITIVE TOBES
.1	24	24
.01	24	24
.001	24	24
.0001	24	11
.00001	24	1
.10000E-05	24	0
.10000E-06	24	0
.10000E-07	24	0
.10000E-08	24	0
.10000E-09	24	0

MOST PROBABLE NUMBER, MPN (per ml) = 6,014 MPN CORRECTED FOR BIAS (Salama) = 5,727 SPEARMAN-KARBER ESTIMATE = 10,000 LOWER 95% CONFIDENCE LIMIT (Cornish & Fisher) = 3,554

LOWER	900	CONLIDENCE		(COLIIISII	œ	FISHEL)	-	5,554
UPPER	95%	CONFIDENCE	LIMIT	(Cornish	&	Fisher)	=	9,696

NUMBER OF DI	CARVA Nutrient 2 LUTIONS = 10 # TUBES	2, Day 0 # POSITIVE TUBES		
	# 10865	# FOSTITVE TOBES		
.1	24	24		
.01	24	24		
.001	24	24		
.0001	24	11		
.00001	24	2		
.10000E-05	24	0		
.10000E-06	24	0		
.10000E-07	24	0		
.10000E-08	24	0		
.10000E-09	24	0		
М	IPN CORRECTED FC	JMBER,,MPN (per ml) DR BIAS (Salama) ESTIMATE	= 6,199	
		(Cornish & Fisher) (Cornish & Fisher)		

DATA SET SARVA	Nutrient	3, Day 0
NUMBER OF DILUTI	ONS = 10	
VOLUME, ml	# TUBES	# POSITIVE TUBES
.1	24	24
.01	24	24
.001	24	24
.0001	24	10
.00001	24	3
.10000E-05	24	0
.10000E-06	24	0
.10000E-07	24	0
.10000E-08	24	0
.10000E-09	24	0

MOST PROBABLE NUMBER, MPN (per ml) = 6,330 MPN CORRECTED FOR BIAS (Salama) = 6,051 SPEARMAN-KARBER ESTIMATE = 11,007 LOWER 95% CONFIDENCE LIMIT (Cornish & Fisher) = 3,718 UPPER 95% CONFIDENCE LIMIT (Cornish & Fisher) = 10,149

	SARVA Nutrient 1 DILUTIONS = 10	, Day 7		
VOLUME, ml	# TUBES	# POSITIVE TUBES		
.1	24	24		
.01	24	24		
.001	24	24		
.0001	24	20		
.00001	24	10		
.10000E-05	24	4		
.10000E-06	24	0		
.10000E-07	24	0		
.10000E-08	24	0		
.10000E-09	24	0		
	MPN CORRECTED FO	MBER,,MPN (per ml) R BIAS (Salama) ESTIMATE	=	27,089
		(Cornish & Fisher) (Cornish & Fisher)		

DATA SET SARVA		2, Day 7
NUMBER OF DILUTI	ONS = 10	
VOLUME, ml	# TUBES	# POSITIVE TUBES
.1	24	24
.01	24	24
.001	24	24
.0001	24	20
.00001	24	11
.10000E-05	24	4
.10000E-06	24	0
.10000E-07	24	0
.10000E-08	24	0
.10000E-09	24	0

MOST PROBABLE NUMBER, MPN (per ml) = 29,933 MPN CORRECTED FOR BIAS (Salama) = 28,330 SPEARMAN-KARBER ESTIMATE = 90,852 LOWER 95% CONFIDENCE LIMIT (Cornish & Fisher) = 18,795 UPPER 95% CONFIDENCE LIMIT (Cornish & Fisher) = 48,388

	SARVA Nutrient 3, DILUTIONS = 10	Day 7		
VOLUME, ml	# TUBES #	POSITIVE TUBES		
.1	24	24		
.01	24	24		
.001	24	24		
.0001	24	19		
.00001	24	10		
.10000E-05	24	4		
.10000E-06	24	0		
.10000E-07	24	0		
.10000E-08	24	0		
.10000E-09	24	0		
	MOST PROBABLE NUME MPN CORRECTED FOR SPEARMAN-KARBER ES	BIAS (Salama)	=	24,239
	CONFIDENCE LIMIT (C CONFIDENCE LIMIT (C	,		

DATA SET SARVA	Nutrient	1, Day 28			
NUMBER OF DILUTIONS = 10					
VOLUME, ml	# TUBES	# POSITIVE TUBES			
.1	24	24			
.01	24	24			
.001	24	24			
.0001	24	24			
.00001	24	24			
.10000E-05	24	24			
.10000E-06	24	9			
.10000E-07	24	4			
.10000E-08	24	2			
.10000E-09	24	0			

MOST PROBABLE NUMBER, MPN (per ml) = 7,068,457 MPN CORRECTED FOR BIAS (Salama) = 6,806,931 SPEARMAN-KARBER ESTIMATE = 13,335,214 LOWER 95% CONFIDENCE LIMIT (Cornish & Fisher) = 4,101,489 UPPER 95% CONFIDENCE LIMIT (Cornish & Fisher) = 11,194,367

DATA SET SA	RVA Nutrient 2 UTIONS = 10	, Day 28	
		# POSITIVE TUBES	
.1	24	24	
.01	24	24	
.001	24	24	
.0001	24	24	
.00001	24	24	
.10000E-05	24	24	
.10000E-06	24	10	
.10000E-07	24	3	
.10000E-08	24	2	
.10000E-09	24	0	
MP	N CORRECTED FOR	MBER,,MPN (per ml) R BIAS (Salama) ESTIMATE	= 6,976,899
		(Cornish & Fisher) (Cornish & Fisher)	

NUMBER OF DILU		Day 28 # POSITIVE TUBES		
.1 .01	24 24 24 24	24 24 24 24		
.001 .0001 .00001	24 24	24 24		
.10000E-05 .10000E-06 .10000E-07		24 10 4		
.10000E-08 .10000E-09	24 24	1 0		
MPI	I CORRECTED FOR	BER,,MPN (per ml) BIAS (Salama) STIMATE	=	6,992,750
		Cornish & Fisher) Cornish & Fisher)		

APPENDIX IV

Two-Sample T-Test and CI: Con_0, Sarva_0

Two-sample T for Con_0 vs Sarva_0 SE Mean StDev N Mean Con 0 3 0.10067 0.00379 0.0022 Sarva 0 3 0.10833 0.00289 0.0017 Difference = mu Con_0 - mu Sarva_0 Estimate for difference: -0.00767 95% CI for difference: (-0.01641, 0.00108) T-Test of difference = 0 (vs not =): T-Value = -2.79 P-Value = 0.068 DF = 3

Two-Sample T-Test and Cl: Con_7, Sarva_7

Two-sample T for Con_7 vs Sarva_7

 N
 Mean
 StDev
 SE Mean

 Con_7
 3
 0.09800
 0.00200
 0.0012

 Sarva<7</td>
 3
 0.09600
 0.00265
 0.0015

Difference = mu Con_7 - mu Sarva_7 Estimate for difference: 0.0020095% CI for difference: (-0.00409, 0.00809) T-Test of difference = 0 (vs not =): T-Value = 1.04 P-Value = 0.373 DF = 3

Two-Sample T-Test and CI: Con_28, Sarva_28

Two-sample T for Con_28 vs Sarva_28

 N
 Mean
 StDev
 SE
 Mean

 Con_28
 3
 0.09533
 0.00321
 0.0019

 Sarva 28
 3
 0.015667
 0.000577
 0.00033

Difference = mu Con_28 - mu Sarva_28 Estimate for difference: 0.07967 95% CI for difference: (0.07155, 0.08778) T-Test of difference = 0 (vs not =): T-Value = 42.25 P-Value = 0.001 DF = 2

Two-Sample T-Test and CI: Con_0, Nutr_0

Two-sample T for Con_0 vs Nutr_0

 N
 Mean
 StDev
 SE Mean

 Con_0
 3
 0.10067
 0.00379
 0.0022

 Nutr 0
 3
 0.097667
 0.000577
 0.00033

Difference = mu Con_0 - mu Nutr_0 Estimate for difference: $0.003\overline{0}0$ 95% CI for difference: (-0.00651, 0.01251) T-Test of difference = 0 (vs not =): T-Value = 1.36 P-Value = 0.308 DF = 2

Two-Sample T-Test and CI: Con_7, Nutr_7

Two-sample T for Con_7 vs Nutr_7

N Mean StDev SE Mean

Con_730.098000.002000.0012Nutr_730.085000.001000.00058

Difference = mu Con_7 - mu Nutr_7 Estimate for difference: 0.01300 95% CI for difference: (0.00745, 0.01855) T-Test of difference = 0 (vs not =): T-Value = 10.07 P-Value = 0.010 DF = 2

Two-Sample T-Test and CI: Con_28, Nutr_28

Two-sample T for Con_28 vs Nutr_28 Ν StDev SE Mean Mean Con 28 3 0.09533 0.00321 0.0019 Nutr_28 3 0.0010 0.02400 0.00173 Difference = mu Con_28 - mu Nutr_28 Estimate for difference: $0.0713\overline{3}$ 95% CI for difference: (0.06462, 0.07804) T-Test of difference = 0 (vs not =): T-Value = 33.84 P-Value = 0.000 DF = 3

Two-Sample T-Test and CI: Nutr_28, Sarva_28

Two-sample T for Nutr_28 vs Sarva_28 SE Mean Ν Mean StDev Nutr 28 3 0.02400 0.00173 0.0010 Sarva_28 3 0.015667 0.000577 0.00033 Difference = mu Nutr 28 - mu Sarva_28 Estimate for difference: 0.00833 95% CI for difference: (0.00380, 0.01287) T-Test of difference = 0 (vs not =): T-Value = 7.91 P-Value = 0.016 DF = 2 APPENDIX V