Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120221\

Data File : BG051310.D

Acq On : 2 Dec 2021 19:04

Operator : CG/JU Sample : PB141106BS

Misc

ALS Vial : 15 Sample Multiplier: 1

Quant Time: Dec 03 00:01:44 2021

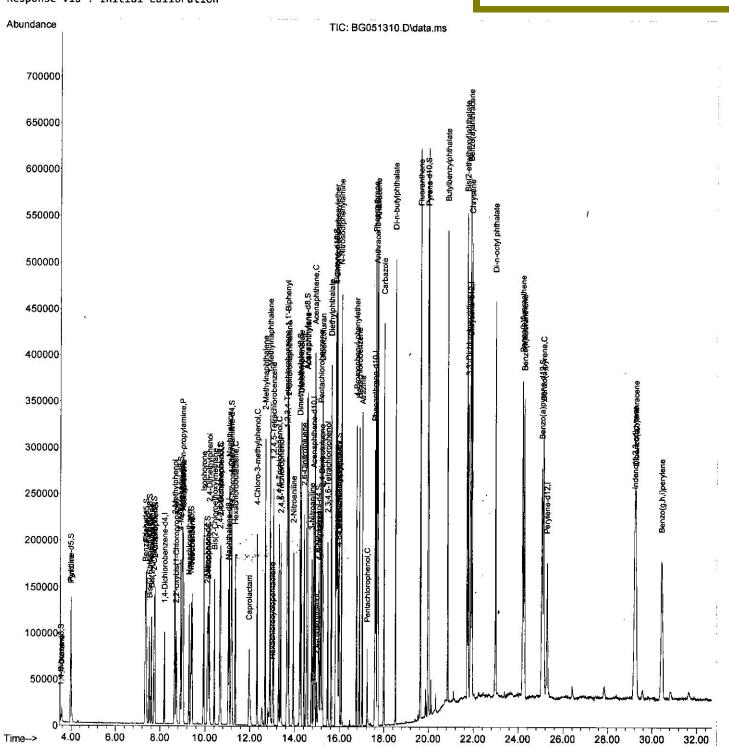
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M

Quant Title : SVOA CALIBRATION
QLast Update : Wed Nov 24 06:04:50 2021
Response via : Initial Calibration

Instrument :
BNA_G
ClientSampleId :

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/03/2021 Supervised By :mohammad ahmed 12/05/2021



SFAM-EPA-BG112321.M Fri Dec 03 00:25:07 2021

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120221\

Data File : BG051310.D

Acq On : 2 Dec 2021 19:04

Operator : CG/JU Sample : PB141106BS

Misc

ALS Vial : 15 Sample Multiplier: 1

Quant Time: Dec 03 00:01:44 2021

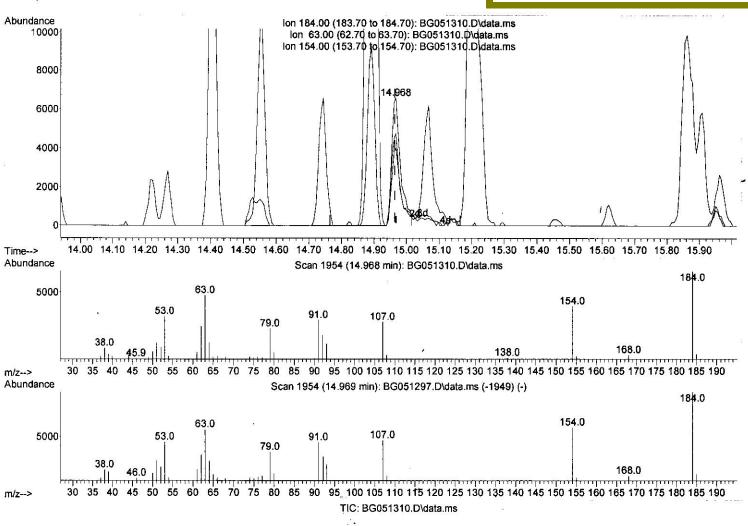
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M

Quant Title : SVOA CALIBRATION
QLast Update : Wed Nov 24 06:04:50 2021
Response via : Initial Calibration

Instrument : BNA_G ClientSampleId :

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/03/2021 Supervised By :mohammad ahmed 12/05/2021



(53) 2,4-Dinitrophenol

14.968min (+ 0.002) 16.67 ng/ul

response	12652			
Ion	Ежр%	Act%		
184.00	100.00	100.00		
63.00	82.70	72.52		
154.00	67.00	60.79		
0.00	0.00	0.00		

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120221\

Data File : BG051310.D

Acq On : 2 Dec 2021 19:04

Operator : CG/JU Sample : PB141106BS

Misc

ALS Vial : 15 Sample Multiplier: 1

Quant Time: Dec 03 00:01:44 2021

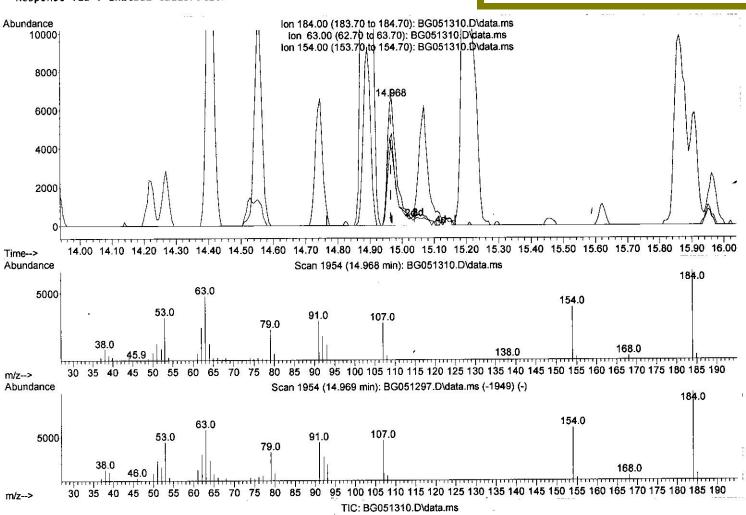
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M

Quant Title : SVOA CALIBRATION
QLast Update : Wed Nov 24 06:04:50 2021
Response via : Initial Calibration

Instrument : BNA_G ClientSampleId :

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/03/2021 Supervised By :mohammad ahmed 12/05/2021



(53) 2,4-Dinitrophenol

14.968min (+ 0.002) 17.61 ng/ul m 3 2/ 20 2

response	13364	
Ion	Exp%	Act%
184.00	100.00	100.00
63.00	82.70	72.52
154.00	67.00	60.79
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120221\

Data File : BG051310.D

Acq On : 2 Dec 2021 19:04

Operator : CG/JU Sample : PB141106BS

Misc

ALS Vial : 15 Sample Multiplier: 1

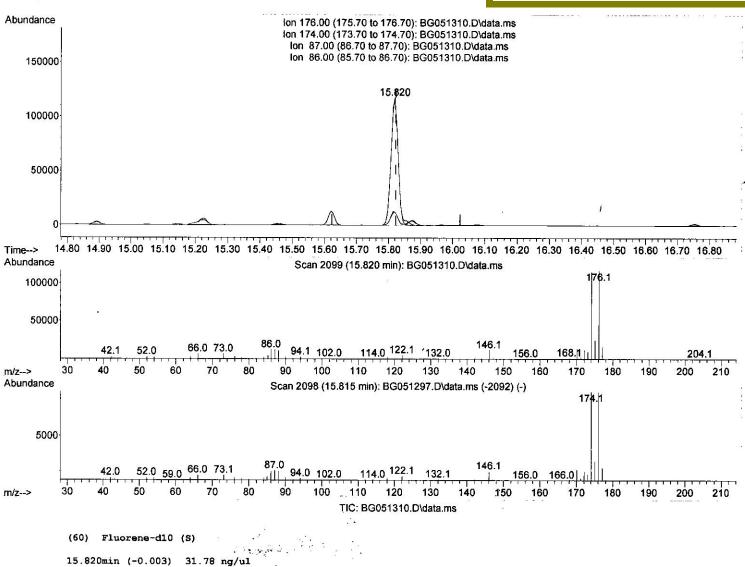
Quant Time: Dec 03 00:01:44 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M

Quant Title : SVOA CALIBRATION QLast Update : Wed Nov 24 06:04:50 2021 Response via : Initial Calibration Instrument :
BNA_G
ClientSampleId :

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/03/2021 Supervised By :mohammad ahmed 12/05/2021



184778

Exps

100.00

Act%

100.00

response Ion

176.00

86.00 10.90 10.57

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120221\

Data File : BG051310.D

: 2 Dec 2021 19:04 Acq On

Operator : CG/JU : PB141106BS Sample

Misc

ALS Vial : 15 Sample Multiplier: 1

Quant Time: Dec 03 00:01:44 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M

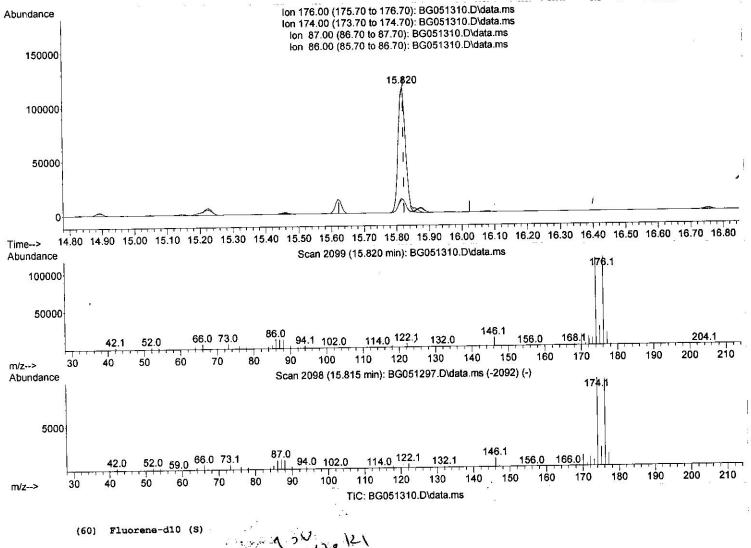
Quant Title : SVOA CALIBRATION

QLast Update : Wed Nov 24 06:04:50 2021 Response via : Initial Calibration

Instrument: BNA_G ClientSampleId:

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/03/2021 Supervised By:mohammad ahmed 12/05/2021



31.17 ng/ul m 3 (2) 15.820min (-0.003)

response	181232	
Ion	Ехр%	Act%
176.00	100.00	100.00
174.00	97.50	98.50
87.00	10.60	10.29
86,00	10.90	10.57

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120221\

Data File : BG051310.D

Acq On : 2 Dec 2021 19:04

Operator : CG/JU Sample : PB1411068S

Misc

ALS Vial : 15 Sample Multiplier: 1

Quant Time: Dec 03 00:01:44 2021

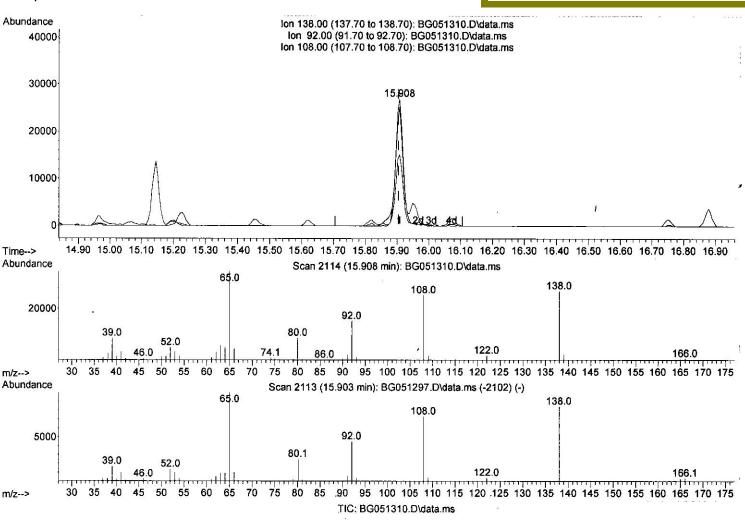
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-8G112321.M

Quant Title : SVOA CALIBRATION
QLast Update : Wed Nov 24 06:04:50 2021
Response via : Initial Calibration

Instrument :
BNA_G
ClientSampleld :

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/03/2021 Supervised By :mohammad ahmed 12/05/2021



(63) 4-Nitroaniline

15.908min (+ 0.002) 35.40 ng/ul

response	46745	
Ion	Ежр%	Act%
138.00	100.00	100.00
92.00	61.60	56.57
108.00	90.70	94.06
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120221\

Data File : BG051310.D

Acq On : 2 Dec 2021 19:04

Operator : CG/JU Sample : PB141106BS

Misc

ALS Vial : 15 Sample Multiplier: 1

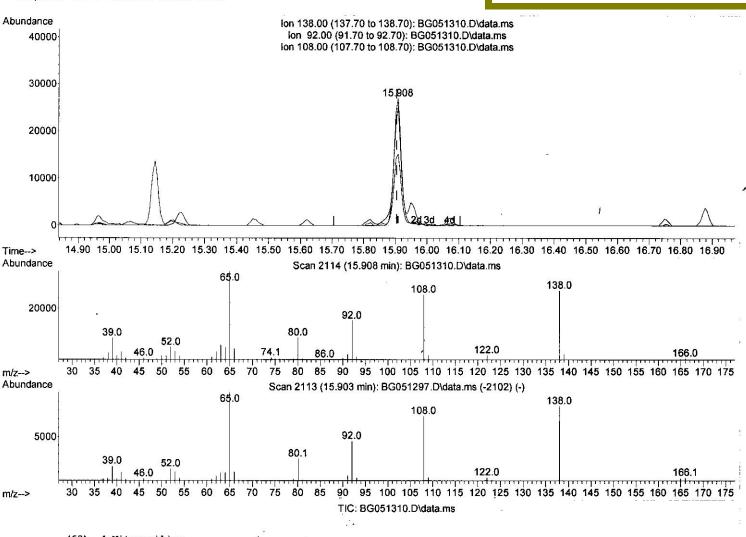
Quant Time: Dec 03 00:01:44 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M

Quant Title : SVOA CALIBRATION QLast Update : Wed Nov 24 06:04:50 2021 Response via : Initial Calibration Instrument : BNA_G ClientSampleId :

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/03/2021 Supervised By :mohammad ahmed 12/05/2021



(63) 4-Nitroaniline

15.908min (+ 0.002) 35.56 ng/ul m

response	46955			
Ion	Exp%	Act%		
138.00	100.00	100.00		
92.00	61.60	56.57		
108.00	90.70	94.06		
0.00	0.00	0.00		

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120221\

Data File : BG051310.D

Acq On : 2 Dec 2021 19:04

Operator : CG/JU Sample : PB141106BS

Misc :

ALS Vial : 15 Sample Multiplier: 1

Quant Time: Dec 03 00:01:44 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M

Quant Title : SVOA CALIBRATION

QLast Update : Wed Nov 24 06:04:50 2021 Response via : Initial Calibration Instrument : BNA_G ClientSampleId : SLCS106

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/03/2021 Supervised By :mohammad ahmed 12/05/2021

Compound	R.T.	QIon	Response	Conc Uni	ts Dev	(Min)	
Internal Standards	· 						
1) 1,4-Dichlorobenzene-d4	8.194	152	27710	20.000	กฮ/มไ	0.00	
20) Naphthalene-d8	11.020	136	126205	20.000		0.00	
38) Acenaphthene-d10	14.827		83917	20.000		0.00	
64) Phenanthrene-d10	17.577		184993	20.000		0.00	
79) Chrysene-d12	21.878	240	151768	20.000		0.00	
88) Perylene-d12	25.280	264	152947	20.000		0.00	
oo, respiese uiz	23,200	204	132347	20.000	iig/ui	0.00	
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.529	96	4328	5.428	ng/uL	-0.02	
4) Pyridine-d5	3.963	84	63482	27.131	10 10 10 10 10 10 10 10 10 10 10 10 10 10 10	-0.02	
7) Phenol-d5	7.354	99	87097	31.803		0.00	
9) Bis-(2-Chloroethyl)eth	7.512	67	53598	31.161		0.00	
11) 2-Chlorophenol-d4	7.724	132	62099	31.489		0.00	
15) 4-Methylphenol-d8	8.911	113	69469	31.434		0.00	
21) Nitrobenzene-d5	9.375	128	32599	30.599		0.00	
24) 2-Nitrophenol-d4	10.097	143	38776	32.266		0.00	
28) 2,4-Dichlorophenol-d3	10.650	165	64565	31.665		0.00	
31) 4-Chloroaniline-d4	11.161	131	114819	38.485		0.00	
46) Dimethylphthalate-d6	14.222	166	204877	31.730		0.00	
49) Acenaphthylene-d8	14.528	160	259138	31.827		0.00	
54) 4-Nitrophenol-d4	15.050	143	29428	28.156	()	0.00 70 /20121	
60) Fluorene-d10	15.820	176	181232m	31.169		0.00 12/20/21	
65) 4,6-Dinitro-2-methylph	15.949	200	30989	27.147		0.00	
73) Anthracene-d10	17.677		278022	31.424		0.00	
81) Pyrene-d10	19.956	212	311425	33.913		0.00	
92) Benzo(a)pyrene-d12	25.050	264	260948	31.946		0.00	
Target Compounds					Ov	alue	
2) 1,4-Dioxane	3.570	88	9329	10.374			
5) Pyridine	3.987	79	67937	27.903	THE RESERVE THE PROPERTY.	91	
6) Benzaldehyde	7.330	77	54097	31.017		92	
8) Phenol	7.383	94	91055	32.094		99	
10) Bis(2-Chloroethyl)ether	7.606	93	67906	31.637		96 .	
12) 2-Chlorophenol	7.759	128	64666	32.177		95	
13) 2-Methylphenol	8.646	108	67796 .			96	
14) 2,2'-oxybis(1-Chloropr	8.711		99212	32.032	1770 - 1879 July 1870 (1870)	98	
16) Acetophenone	9.028		107885	31.560		98	
17) N-Nitroso-di-n-propyla	8.999	70	64072	32.617		99	
18) 4-Methylphenol	8.975	-	72255	31.975		95	
19) Hexachloroethane	9.281	117	26383	31.081		94.	
22) Nitrobenzene	9.416	77	88262	31.596		96	
23) Isophorone	9.933	82	174306	32.117		98	
25) 2-Nitrophenol	10.133	139	38866	31.223		99	
26) 2,4-Dimethylphenol	10.186	107	81502	32.025		97	
27) Bis(2-Chloroethoxy)met	10.409	93	94726	31.616		99	
29) 2,4-Dichlorophenol	10.673	162	64933	32.351		100	
30) Naphthalene	11.073	128	215843	31.432		100	
32) 4-Chloroaniline	11.184	127	85742	28.627		97	
33) Hexachlorobutadiene	11.337	225	41134	29.712		97	
34) Caprolactam	11.954	113	25076	31.779		92	
35) 4-Chloro-3-methylphenol	12.307	107	79490	32.968		96	
					· · · · · ·		

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120221\

Data File : BG051310.D

Acq On : 2 Dec 2021 19:04

Operator : CG/JU Sample : PB141106BS

Misc :

ALS Vial : 15 Sample Multiplier: 1

Quant Time: Dec 03 00:01:44 2021

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M

Quant Title : SVOA CALIBRATION

QLast Update : Wed Nov 24 06:04:50 2021 Response via : Initial Calibration

nstrument :	
BNA_G	
ClientSampleId:	
01.004.00	

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/03/2021 Supervised By :mohammad ahmed 12/05/2021

Compound	R.T.	QIon	Response	Conc Units Dev	(Min)
36) 2-Methylnaphthalene	12.665	142	146233	31.307 ng/ul	99
<pre>37) 1-Methylnaphthalene</pre>	12.882	142	148516	30.906 ng/ul	99
39) 1,2,4,5-Tetrachloroben	13.029	216	83420	31.665 ng/ul	96
40) Hexachlorocyclopentadiene	12,994	237	9860	9.260 ng/ul#	
41) 2,4,6-Trichlorophenol	13.276	196	53837	32.565 ng/ul	99
42) 2,4,5-Trichlorophenol	13.358	196	57137	33.003 ng/ul	97
43) 1,1'-Biphenyl	13.658	154	199071	31.761 ng/ul	98
44) 2-Chloronaphthalene	13.711	162	156553	31.400 ng/ul	98
45) 2-Nitroaniline	13.922	65	58503	33.904 ng/ul	93
47) Dimethylphthalate	14.269	163	205947	31.511 ng/ul	100
48) 2,6-Dinitrotoluene	14.404	165	44973	32.759 ng/ul	95
50) Acenaphthylene	14.557	152	257517	32.013 ng/ul	98
51) 3-Nitroaniline	14.745	138	44999	33.160 ng/ul	98
52) Acenaphthene	14.892	153	169852	32.017 ng/ul	96
53) 2,4-Dinitrophenol	14.968	184	13364m	17.611 ng/ul	
55) 4-Nitrophenol	15.068	109	26107	28.795 ng/ul	96
56) Dibenzofuran	15.227	168	242167	∖31.647 ng/ul	99
57) 2,4-Dinitrotoluene	15.197	165	62891	32.074 ng/ul	99
58) 2,3,4,6-Tetrachlorophenol	15.456	232	42920	31.570 ng/ul	97
59) Diethylphthalate	15.620	149	219899	32.054 ng/ul	99
61) Fluorenė	15.873	166	192787	31.453 ng/ul	97
62) 4-Chlorophenyl-phenyle	15.855	204	102243	30.953 ng/ul	96
63) 4-Nitroaniline	15.908	138	46955m	35.557 ng/ul	
66) 4,6-Dinitro-2-methylph	15.967	198	29371 -	J 26.679 ng/ul#	96
67) N-Nitrosodiphenylamine	16.073	169	173037	32.673 ng/ul	99
68) 4-Bromophenyl-phenylether	16.754	248	63478	32.016 ng/ul	93
69) Hexachlorobenzene	16.878	284	65740	32.517 ng/ul	98
70) Atrazine	17.019	200	64720	29.078 ng/ul	98
71) Pentachlorophenol	17.236	266	17729	19.790 ng/ul	99
72) Phenanthrene	17.624	178	328869	32.197 ng/ul	99
74) Anthracene	17.712	178	323589	31.899 ng/ul	99
75) 1,2,3,4-Tetrachloroben	13.634	216	85640	31.738 ng/uL	96
76) Pentachlorobenzene	15.144	250	79234	31.515 ng/uL	99
77) Carbazole	17.988	167	294864	33.115 ng/ul	98 .
78) Di-n-butylphthalate	18.505	149	371514	32.358 ng/ul	99
80) Fluoranthene	19.622	202	385249		98
82) Pyrene	19.986		376015	34.081 ng/ul	99
83) Butylbenzylphthalate	20.844		153094	33.377 ng/ul	96
84) 3,3'-Dichlorobenzidine	21.766	252	112598	31.865 ng/ul	98
85) Benzo(a)anthracene	21.860		338086	32.844 ng/ul	99 -
86) Bis(2-ethylhexyl)phtha	21.719	149	219537	33.261 ng/ul	99.
87) Chrysene	21.931	228	321268	32.488 ng/ul	100
89) Di-n-octyl phthalate	22.976	149	373268	33.687 ng/ul	100
90) Benzo(b)fluoranthene	24.193	252	335245	32.479 ng/ul	98
91) Benzo(k)fluoranthene	24.263	252	314209	32.439 ng/ul	99
93) Benzo(a)pyrene	25.121	252	319222	32.417 ng/ul	98
94) Indeno(1,2,3-cd)pyrene	29.210	276	361507	32.806 ng/ul	98
95) Dibenzo(a,h)anthracene	29.257	278	310159	33.178 ng/ul	98
96) Benzo(g,h,i)perylene	30.432	276	303573	32.744 ng/ul	95

^{(#) =} qualifier out of range (m) = manual integration (+) = signals summed