Data File : BM033368.D

: 09 Dec 2021 21:36 Acq On

Operator : CG/JU : PB141265BS Sample

Misc

ALS Vial : 22 Sample Multiplier: 1

Quant Time: Dec 10 01:16:05 2021

Quant Method: Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM120921.M

Quant Title : SVOA CALIBRATION

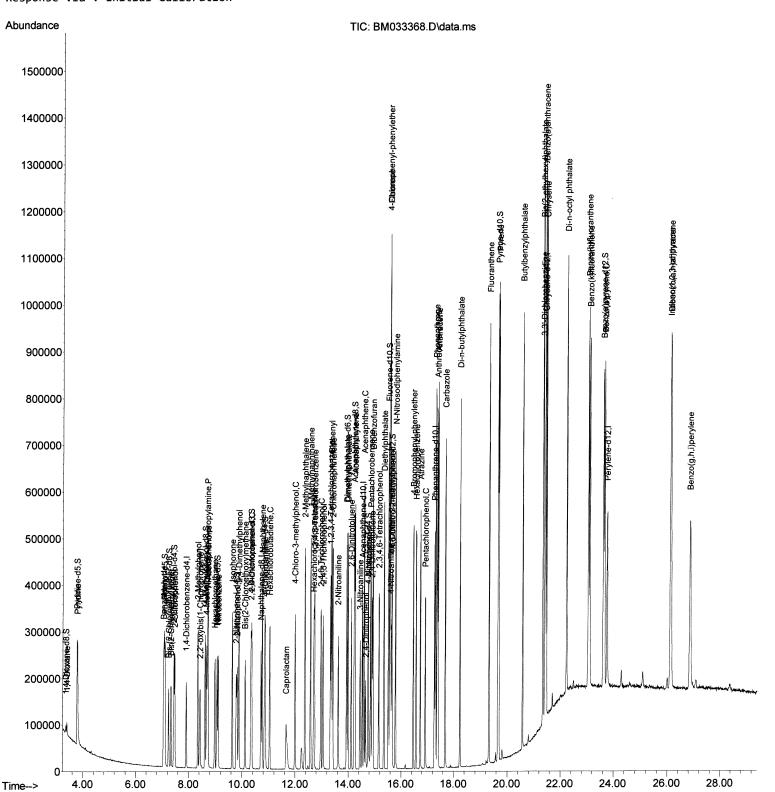
QLast Update : Thu Dec 09 13:25:37 2021

Response via: Initial Calibration



Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/10/2021 Supervised By :mohammad ahmed 12/15/2021



Data File: BM033368.D

Acq On : 09 Dec 2021 21:36

Operator : CG/JU Sample : PB141265BS

Misc ALS Vial

: 22 Sample Multiplier: 1

Quant Time: Dec 10 01:16:05 2021

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM120921.M

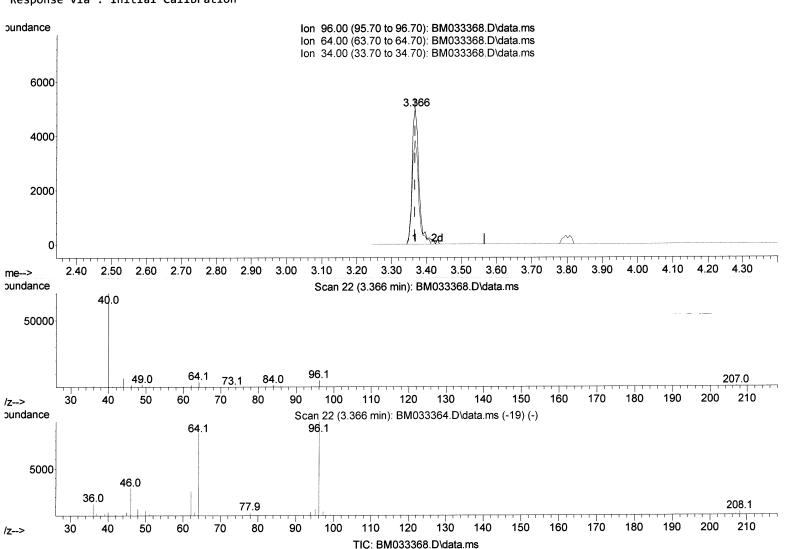
Quant Title : SVOA CALIBRATION

QLast Update : Thu Dec 09 13:25:37 2021 Response via : Initial Calibration

Instrument: BNA_M ClientSampleId: SLCS265

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/10/2021 Supervised By :mohammad ahmed 12/15/2021



(3) 1,4-Dioxane-d8 (S)

3.366min (+ 0.000) 5.86 ng/uL

response	6764	
Ion	Exp%	Act%
96.00	100.00	100.00
64.00	74.20	76.60
34.00	0.00	0.00
0.00	0.00	0.00

Data File : BM033368.D

Acq On : 09 Dec 2021 21:36

Operator : CG/JU Sample : PB141265BS

Misc ALS Vial

: 22 Sample Multiplier: 1

Quant Time: Dec 10 01:16:05 2021

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM120921.M

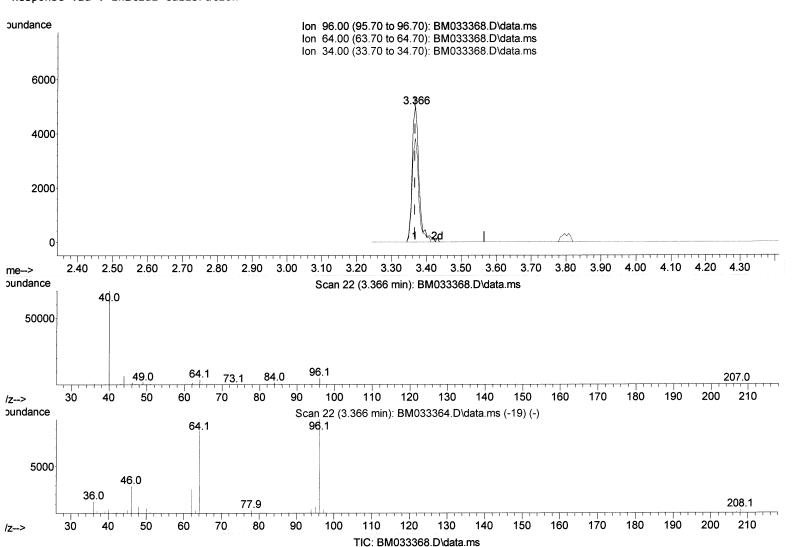
Quant Title : SVOA CALIBRATION

QLast Update : Thu Dec 09 13:25:37 2021 Response via : Initial Calibration

Instrument: BNA_M ClientSampleId: SLCS265

Manual IntegrationsAPPROVED

Reviewed By: Jagrut Upadhyay 12/10/2021 Supervised By: mohammad ahmed 12/15/2021



(3) 1,4-Dioxane-d8 (S)

3.366min	(+ 0.000)	5.98 ng/u	ıL m	p/28/21
response	6903		J4	101.01
Ion	Exp%	Act%		
96.00	100.00	100.00		
64.00	74.20	76.60		
34.00	0.00	0.00		
0.00	0.00	0.00		

Data File : BM033368.D

Acq On : 09 Dec 2021 21:36

Operator : CG/JU Sample : PB141265BS

Misc

ALS Vial : 22 Sample Multiplier: 1

Quant Time: Dec 10 01:16:05 2021

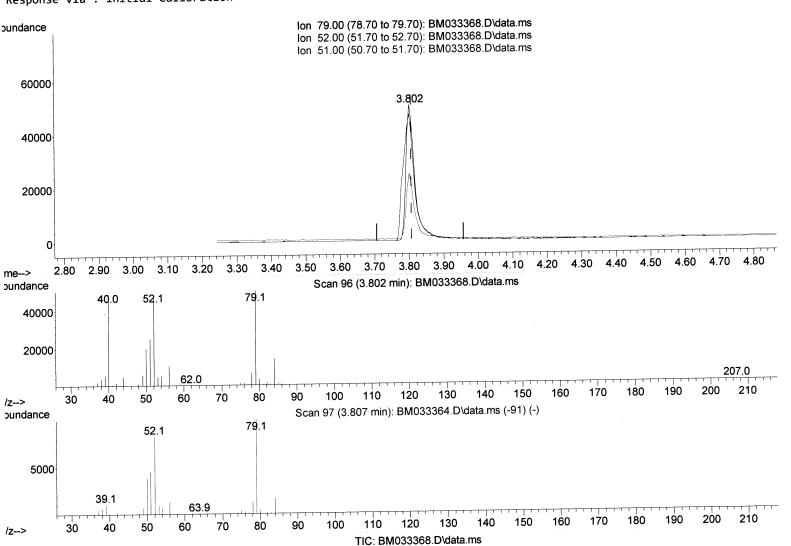
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM120921.M

Quant Title : SVOA CALIBRATION

QLast Update : Thu Dec 09 13:25:37 2021 Response via : Initial Calibration Instrument :
BNA_M
ClientSampleId :
SLCS265

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/10/2021 Supervised By :mohammad ahmed 12/15/2021



(5) Pyridine

3.802min (-0.006) 25.52 ng/ul

response	87926	
Ion	\mathbf{Exp} %	Act%
79.00	100.00	100.00
52.00	97.60	93.36
51.00	45.00	49.51
0.00	0.00	0.00

Data File : BM033368.D

Acq On : 09 Dec 2021 21:36

Operator : CG/JU Sample : PB141265BS

Misc

ALS Vial : 22 Sample Multiplier: 1

Quant Time: Dec 10 01:16:05 2021

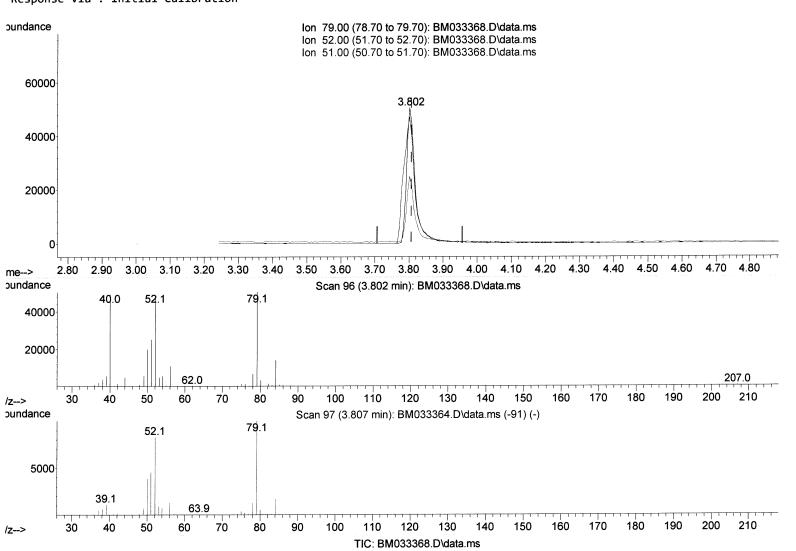
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM120921.M

Quant Title : SVOA CALIBRATION

QLast Update : Thu Dec 09 13:25:37 2021 Response via : Initial Calibration Instrument :
BNA_M
ClientSampleId :
SLCS265

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/10/2021 Supervised By :mohammad ahmed 12/15/2021



(5) Pyridine

3.802min (-0.006) 25.94 ng/ul m

response	89361		JU!
Ion	Exp%	Act%	
79.00	100.00	100.00	
52.00	97.60	93.36	
51.00	45.00	49.51	
0.00	0.00	0.00	

Data File: BM033368.D

Acq On : 09 Dec 2021 21:36

Operator : CG/JU Sample : PB141265BS

Misc :

ALS Vial : 22 Sample Multiplier: 1

Quant Time: Dec 10 01:16:05 2021

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM120921.M

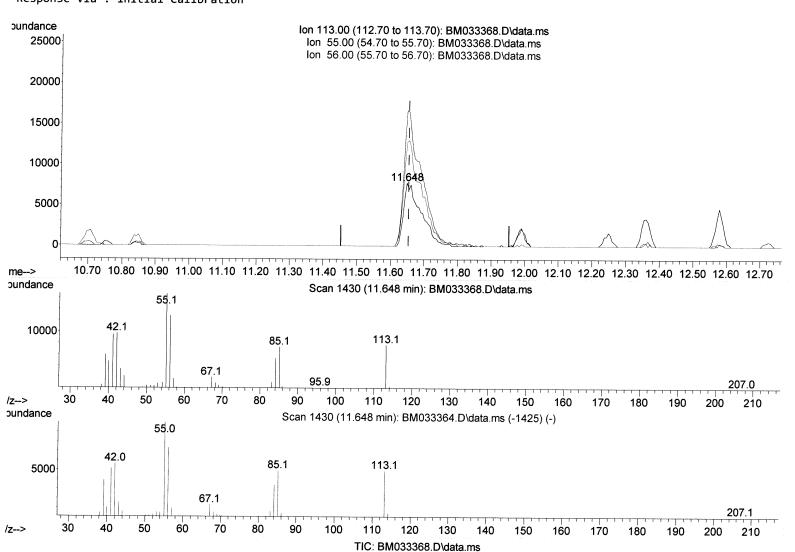
Quant Title : SVOA CALIBRATION

QLast Update : Thu Dec 09 13:25:37 2021 Response via : Initial Calibration

Instrument :
BNA_M
ClientSampleId :
SLCS265

Manual Integrations APPROVED

Reviewed By :Jagrut Upadhyay 12/10/2021 Supervised By :mohammad ahmed 12/15/2021



(34) Caprolactam

11.648min (-0.006) 28.47 ng/ul

response	28219	
Ion	Ехр%	Act%
113.00	100.00	100.00
55.00	197.40	215.10
56.00	164.70	165.21
0.00	0.00	0.00

Data File: BM033368.D

Acq On : 09 Dec 2021 21:36

Operator : CG/JU Sample : PB141265BS

Misc

ALS Vial : 22 Sample Multiplier: 1

Quant Time: Dec 10 01:16:05 2021

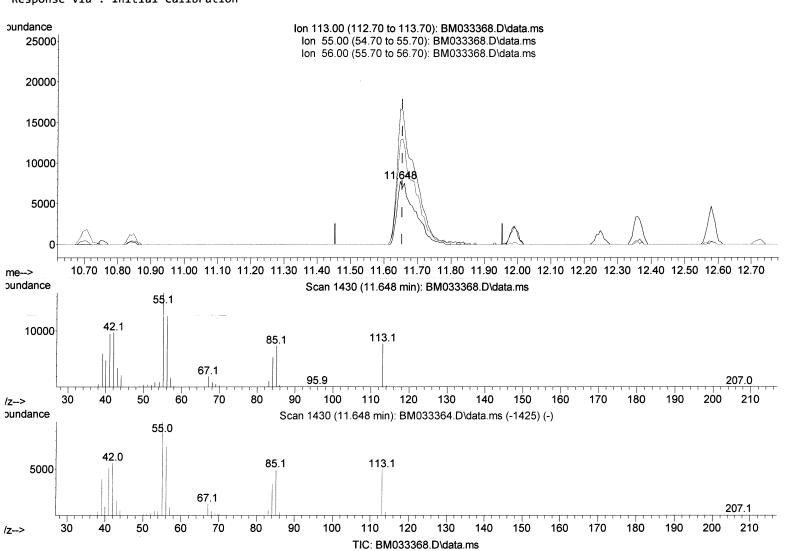
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM120921.M

Quant Title : SVOA CALIBRATION

QLast Update : Thu Dec 09 13:25:37 2021 Response via : Initial Calibration Instrument: BNA_M ClientSampleId: SLCS265

Manual Integrations APPROVED

Reviewed By :Jagrut Upadhyay 12/10/2021 Supervised By :mohammad ahmed 12/15/2021



(34) Caprolactam

11.648min	(-0.006)	28.63 ng/ul	Tun 20/21
response	28382		Juin .
Ion	Exp %	Act%	
113.00	100.00	100.00	
55.00	197.40	215.10	
56.00	164.70	165.21	
0.00	0.00	0.00	

Data File : BM033368.D

Acq On : 09 Dec 2021 21:36

Operator : CG/JU
Sample : PB141265BS

4isc :

ALS Vial : 22 Sample Multiplier: 1

Quant Time: Dec 10 01:16:05 2021

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM120921.M

Quant Title : SVOA CALIBRATION

QLast Update : Thu Dec 09 13:25:37 2021 Response via : Initial Calibration

Instrument : BNA_M ClientSampleld : SLCS265

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/10/2021 Supervised By :mohammad ahmed 12/15/2021

Internal Standards 1) 1,4-Dichlorobenzene-d4 7.907 152 43352 20.000 ng/ul 0.00 20) Naphthalene-d8 10.701 136 184578 20.000 ng/ul 0.00 38) Acenaphthene-d10 14.536 164 127293 20.000 ng/ul 0.00 64) Phenanthrene-d10 17.271 188 275152 20.000 ng/ul 0.00 79) Chrysene-d12 21.436 240 284358 20.000 ng/ul 0.00	2)
1) 1,4-Dichlorobenzene-d4 7.907 152 43352 20.000 ng/ul 0.00 20) Naphthalene-d8 10.701 136 184578 20.000 ng/ul 0.00 38) Acenaphthene-d10 14.536 164 127293 20.000 ng/ul 0.00 64) Phenanthrene-d10 17.271 188 275152 20.000 ng/ul 0.00 79) Chrysene-d12 21.436 240 284358 20.000 ng/ul 0.00	2)
20) Naphthalene-d8 10.701 136 184578 20.000 ng/ul 0.00 38) Acenaphthene-d10 14.536 164 127293 20.000 ng/ul 0.00 64) Phenanthrene-d10 17.271 188 275152 20.000 ng/ul 0.00 79) Chrysene-d12 21.436 240 284358 20.000 ng/ul 0.00	2)
38) Acenaphthene-d10 14.536 164 127293 20.000 ng/ul 0.00 64) Phenanthrene-d10 17.271 188 275152 20.000 ng/ul 0.00 79) Chrysene-d12 21.436 240 284358 20.000 ng/ul 0.00	2)
64) Phenanthrene-d10 17.271 188 275152 20.000 ng/ul 0.00 79) Chrysene-d12 21.436 240 284358 20.000 ng/ul 0.00	2)
79) Chrysene-d12 21.436 240 284358 20.000 ng/ul 0.00	2)
	2)
88) Perylene-d12 23.759 264 263706 20.000 ng/ul 0.00	2)
System Monitoring Compounds 3) 1,4-Dioxane-d8 4) Pyridine-d5 3.366 40 87908 3.384 3.366 3.384	
3) 1,4-Dioxane-d8 3.366 96 6903m 5.984 ng/uL $\supset 0.00 > 1/(1)$	
4) Pyridine-d5 3.784 84 87908 26.282 ng/ul 0.00	
7) Phenol-d5 7.078 99 116233 28.314 ng/ul 0.00	
9) Bis-(2-Chloroethyl)eth 7.243 67 76855 28.597 ng/ul 0.00	
11) 2-Chlorophenol-d4 7.443 132 84699 29.469 ng/ul 0.00	
15) 4-Methylphenol-d8 8.619 113 92137 28.675 ng/ul 0.00	
21) Nitrobenzene-d5 9.072 128 44471 29.693 ng/ul 0.00	
24) 2-Nitrophenol-d4 9.789 143 46786 30.448 ng/ul 0.00	
28) 2,4-Dichlorophenol-d3 10.331 165 89748 30.906 ng/ul 0.00	
31) 4-Chloroaniline-d4 10.842 131 127047 29.509 ng/ul 0.00	
46) Dimethylphthalate-d6 13.942 166 290373 30.530 ng/ul 0.00	
49) Acenaphthylene-d8 14.230 160 360008 30.534 ng/ul 0.00	
54) 4-Nitrophenol-d4 14.742 143 51587 29.873 ng/ul 0.00	
60) Fluorene-d10 15.524 176 261714 30.763 ng/ul 0.00	
65) 4,6-Dinitro-2-methylph 15.642 200 48294 29.080 ng/ul 0.00	
73) Anthracene-d10 17.371 188 418518 30.777 ng/ul 0.00	
81) Pyrene-d10 19.654 212 499333 31.419 ng/ul 0.00	
92) Benzo(a)pyrene-d12 23.612 264 453047 31.696 ng/ul 0.00	
Target Compounds Qvalue	
2) 1,4-Dioxane 3.402 88 13635 10.619 ng/uL 95 h 2/3/	
5) Pyridine 3.802 79 89361mシ 25.940 ng/ulシ バルグ へ	
2) 1,4-Dioxane 3.402 88 13635 10.619 ng/uL 95 5) Pyridine 3.802 79 89361m 25.940 ng/ul 93 6) Benzaldehyde 7.054 77 79683 34.841 ng/ul 93	
8) Phenol 7.107 94 115070 27.213 ng/ul 97	
10) Bis(2-Chloroethyl)ether 7.337 93 87803 27.602 ng/ul 97	
12) 2-Chlorophenol 7.478 128 86044 28.951 ng/ul 95	
13) 2-Methylphenol 8.354 108 84220 27.472 ng/ul 94	
14) 2,2'-oxybis(1-Chloropr 8.437 45 148828 27.114 ng/ul 98	
16) Acetophenone 8.737 105 146428 27.510 ng/ul 97	
17) N-Nitroso-di-n-propyla 8.713 70 85393 29.350 ng/ul 100	
18) 4-Methylphenol 8.678 108 90727 27.049 ng/ul 97	
19) Hexachloroethane 8.984 117 41862 27.880 ng/ul 86	
22) Nitrobenzene 9.113 77 126319 28.808 ng/ul 98	
23) Isophorone 9.636 82 218156 29.055 ng/ul 99	
25) 2-Nitrophenol 9.819 139 48117 29.537 ng/ul 98	
26) 2,4-Dimethylphenol 9.878 107 107214 27.144 ng/ul 95	
27) Bis(2-Chloroethoxy)met 10.113 93 120630 28.566 ng/ul 99	
29) 2,4-Dichlorophenol 10.354 162 86511 29.430 ng/ul 93	
30) Naphthalene 10.754 128 289063 28.103 ng/ul 99	
32) 4-Chloroaniline 10.866 127 108951 25.178 ng/ul 99	2.1
33) Hexachlorobutadiene 11.031 225 61830 28.735 ng/ul 98	- 1
32) 4-Chloroaniline 10.866 127 108951 25.178 ng/ul 99 33) Hexachlorobutadiene 11.031 225 61830 28.735 ng/ul 98 34) Caprolactam 11.648 113 28382m 28.630 ng/ul 98 35) 4-Chloro-3-methylphenol 11.989 107 105814 30.666 ng/ul	
35) 4-Chloro-3-methylphenol 11.989 107 105814 30.666 ng/ul 91	

Data File : BM033368.D

Acq On : 09 Dec 2021 21:36

Dperator : CG/JU
Sample : PB141265BS

4isc :

ALS Vial : 22 Sample Multiplier: 1

Quant Time: Dec 10 01:16:05 2021

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM120921.M

Quant Title : SVOA CALIBRATION

QLast Update : Thu Dec 09 13:25:37 2021 Response via : Initial Calibration

Instrument :
BNA_M
ClientSampleId :
SLCS265

Manual IntegrationsAPPROVED

Reviewed By: Jagrut Upadhyay 12/10/2021 Supervised By: mohammad ahmed 12/15/2021

Compound	R.T.	QIon	Response	Conc Units Dev(Min)
36) 2-Methylnaphthalene	12.360	142	198701	28.485 ng/ul	99
37) 1-Methylnaphthalene	12.577		204118	28.157 ng/ul	99
39) 1,2,4,5-Tetrachloroben	12.725	216	106683	27.816 ng/ul	98
40) Hexachlorocyclopentadiene	12.695	237	79477	31.751 ng/ul	97
41) 2,4,6-Trichlorophenol	12.972	196	68452	30.322 ng/ul	95
42) 2,4,5-Trichlorophenol	13.048	196	74660	30.527 ng/ul	92
43) 1,1'-Biphenyl	13.366	154	276638	28.407 ng/ul	98
44) 2-Chloronaphthalene	13.413	162	212374	28.345 ng/ul	99
45) 2-Nitroaniline	13.619	65	84234	31.524 ng/ul	99
47) Dimethylphthalate	13.989		273870	29.001 ng/ul	100
48) 2,6-Dinitrotoluene	14.113	165	55774	30.553 ng/ul	95
50) Acenaphthylene	14.254	152	357023	29.083 ng/ul	99
51) 3-Nitroaniline	14.448	138	53280	29.652 ng/ul	99
52) Acenaphthene	14.595	153	233023	28.601 ng/ul	97
53) 2,4-Dinitrophenol	14.654	184	30759	28.672 ng/ul	92
55) 4-Nitrophenol	14.760	109	56138	30.077 ng/ul	92
56) Dibenzofuran	14.930	168	340388	28.811 ng/ul	98
57) 2,4-Dinitrotoluene	14.901	165	84690	31.580 ng/ul#	93
58) 2,3,4,6-Tetrachlorophenol	15.160	232	63557	30.570 ng/ul#	99
59) Diethylphthalate	15.348	149	287860	29.370 ng/ul	100
61) Fluorene	15.577	166	287734	29.619 ng/ul	99
62) 4-Chlorophenyl-phenyle	15.571	204	142362	29.378 ng/ul	96
63) 4-Nitroaniline	15.607	138	62677	33.968 ng/ul	99
66) 4,6-Dinitro-2-methylph	15.654	198	46875	28.335 ng/ul	93
67) N-Nitrosodiphenylamine	15.789	169	239147	29.560 ng/ul	99
68) 4-Bromophenyl-phenylether	16.465	248	83730	30.212 ng/ul	96
69) Hexachlorobenzene	16.571	284	91171	28.552 ng/ul	96
70) Atrazine	16.736	200	89643	27.925 ng/ul	99
71) Pentachlorophenol	16.924	266	57779	32.269 ng/ul	97
72) Phenanthrene	17.312	178	466702	29.402 ng/ul	98
74) Anthracene	17.407	178	471112	29.274 ng/ul	99
75) 1,2,3,4-Tetrachloroben	13.330	216	111090	27.994 ng/uL	97
76) Pentachlorobenzene	14.848	250	108466	27.678 ng/uL	95
77) Carbazole	17.677	167	421031	28.946 ng/ul	99
78) Di-n-butylphthalate	18.230	149	503197	30.831 ng/ul	100
80) Fluoranthene	19.324	202	555654	29.662 ng/ul	98
82) Pyrene	19.683	202	589359	30.052 ng/ul	100
83) Butylbenzylphthalate	20.571	149	231717	30.846 ng/ul	98
84) 3,3'-Dichlorobenzidine	21.353	252	169384	26.208 ng/ul	96
85) Benzo(a)anthracene	21.418	228	552094	29.551 ng/ul	99
86) Bis(2-ethylhexyl)phtha	21.336	149	334774	31.007 ng/ul	98
87) Chrysene	21.471	228	533413	29.050 ng/ul	99
89) Di-n-octyl phthalate	22.242	149	573347	29.654 ng/ul	100
90) Benzo(b)fluoranthene	23.053	252	552278	30.567 ng/ul	99
91) Benzo(k)fluoranthene	23.100	252	508446	30.393 ng/ul	100
93) Benzo(a)pyrene	23.659	252	525122	30.290 ng/ul	100
94) Indeno(1,2,3-cd)pyrene	26.130	276	547607 491101	29.045 ng/ul	99
95) Dibenzo(a,h)anthracene 96) Benzo(g,h,i)perylene	26.141	278 276	481101 471114	29.336 ng/ul	98 98
oo, benzo(g,n,1)perytene	26.853 	276	471114 	29.296 ng/ul	<i>3</i> 0

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