

(QT Reviewed)

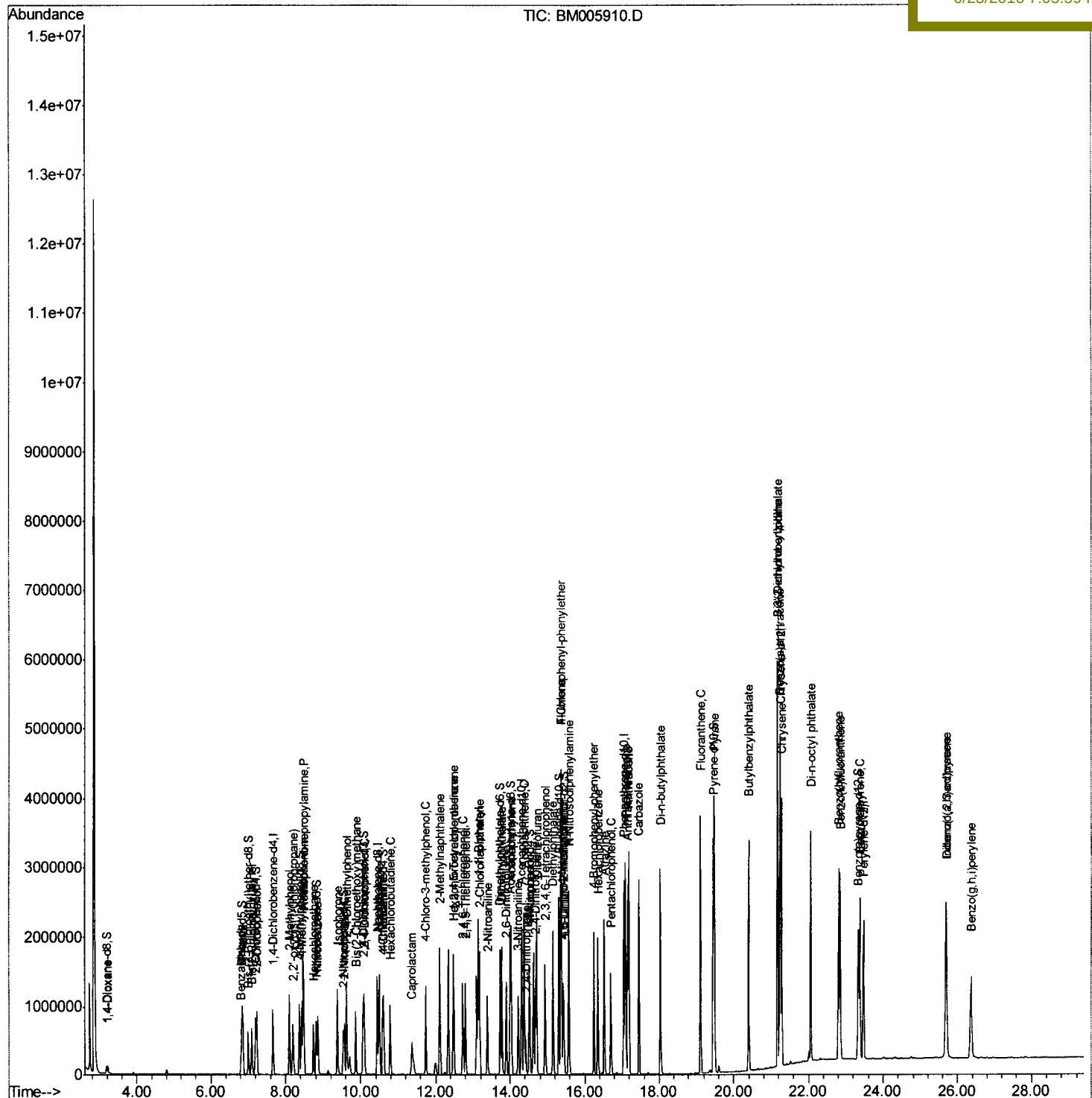
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Data Path   : Z:\HPCHEM1\BNA M\DATA\BM062216\  
Data File  : BM005910.D  
Acq On     : 22 Jun 2016   17:46  
Operator   : UM/SJ  
Sample     : SSTD02045  
Misc       :  
ALS Vial   : 5      Sample Multiplier: 1
```

Quant Time: Jun 23 01:28:51 2016
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM062216.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jun 23 01:14:46 2016
Response via : Initial Calibration

Instrument :
BNA_M
ClientSampleId :
SSTD02045

Manual Integrations APPROVED

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Quantitation Report (Qedit)

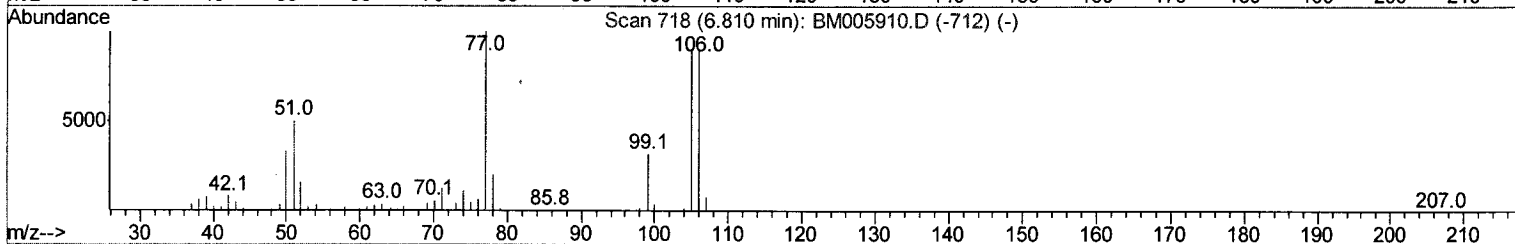
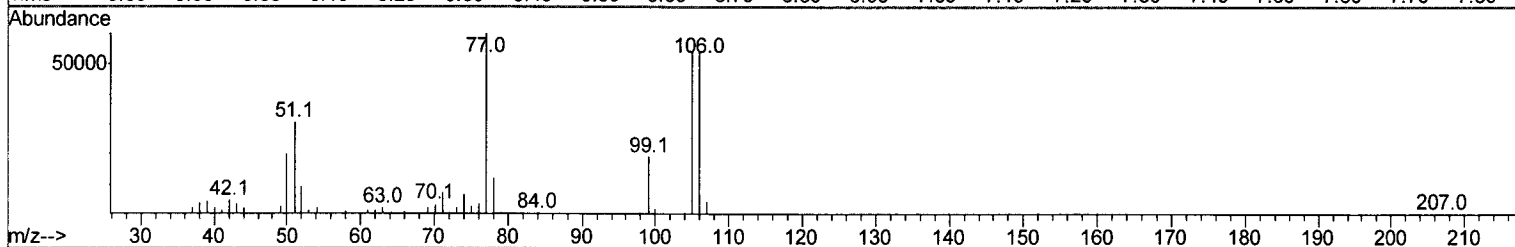
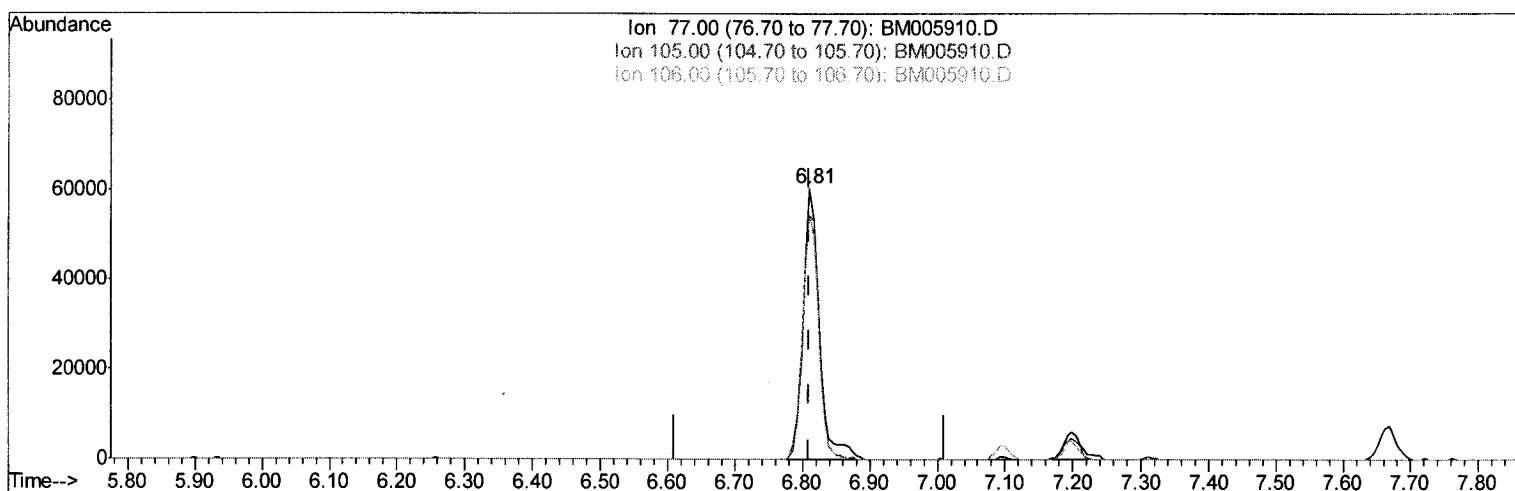
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Quant Time: Jun 23 01:18:06 2016
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 Quant Title : SVOA CALIBRATION
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 Response via : Initial Calibration



TIC: BM005910.D

(4) Benzaldehyde

6.810min (0.000) 28.00ng/ul m

response 100673

U.M
 06/28/16

Ion	Exp%	Act%
77.00	100	100
105.00	90.10	90.14
106.00	90.30	90.28
0.00	0.00	0.00

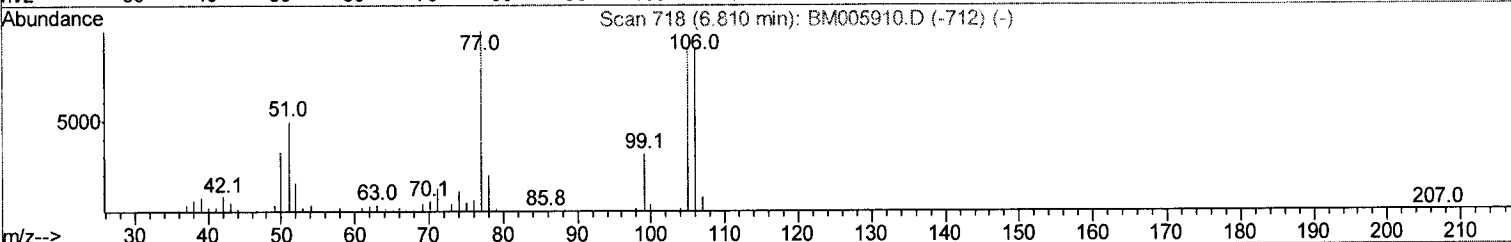
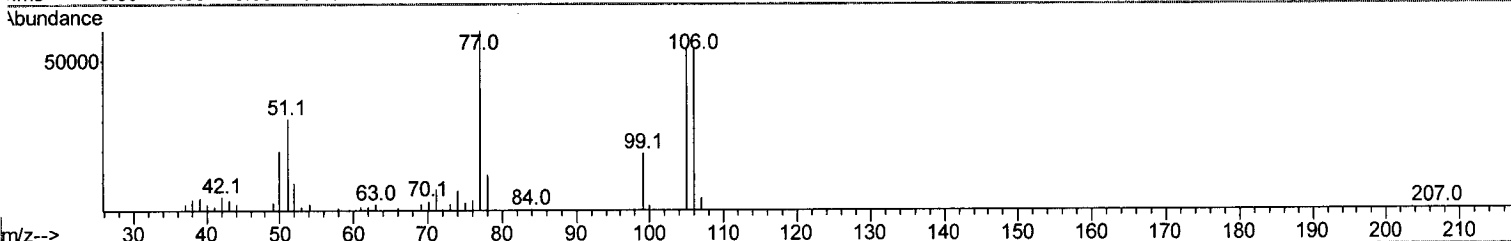
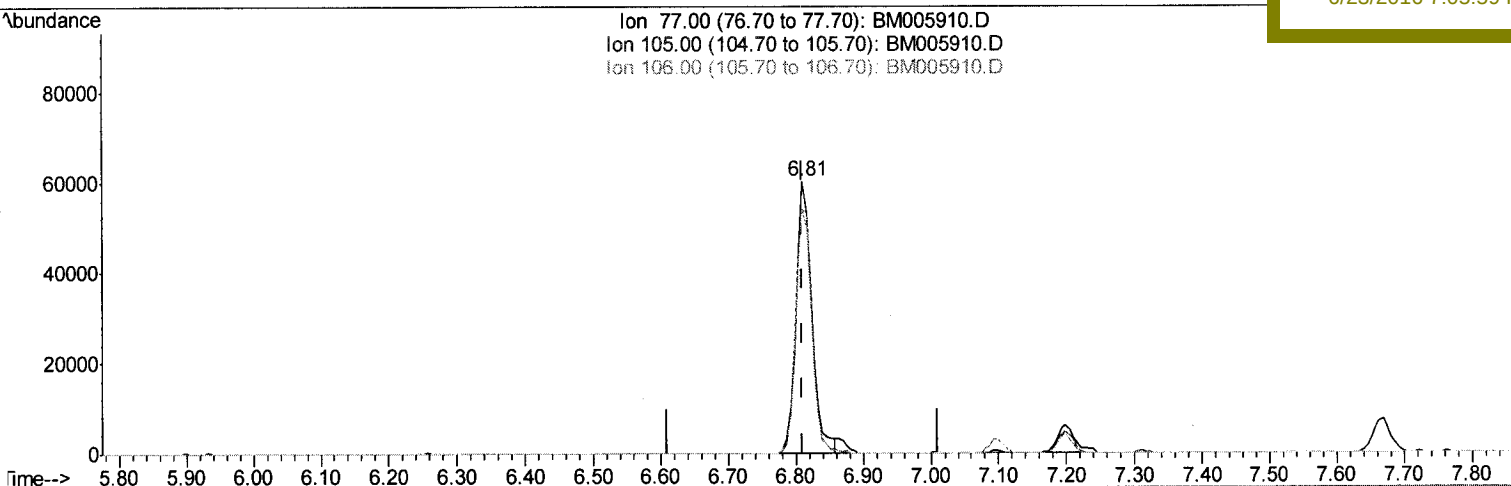
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Manual Integrations
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TIC: BM005910.D

(4) Benzaldehyde

6.810min (0.000) 27.09ng/ul

response 97390

Ion	Exp%	Act%
77.00	100	100
105.00	90.10	90.14
106.00	90.30	90.28
0.00	0.00	0.00

Data Path : Z:\HPCHEM1\BNA M\DATA\BM062216\
 Data File : BM005910.D
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 Operator : UM/SJ
 Sample : SSTD02045
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA M
 ClientSampleId :
 SSTD02045

Quant Time: Jun 23 01:28:51 2016
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 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jun 23 01:14:46 2016
 Response via : Initial Calibration

Manual Integrations
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	249367	20.00	ng/ul	0.00
18) Naphthalene-d8	10.45	136	1170668	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.32	164	696334	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.06	188	1675430	20.00	ng/ul	0.00
75) Chrysene-d12	21.26	240	1724205	20.00	ng/ul	0.00
83) Perylene-d12	23.47	264	1474614	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.18	96	49522	7.93	ng/uL	0.00
5) Phenol-d5	6.83	99	463372	17.65	ng/ul	0.00
7) Bis-(2-Chloroethyl) ether-d	7.00	67	287465	18.81	ng/ul	0.00
9) 2-Chlorophenol-d4	7.20	132	354884	18.16	ng/ul	0.00
13) 4-Methylphenol-d8	8.37	113	377425	17.36	ng/ul	0.00
19) Nitrobenzene-d5	8.82	128	174764	20.67	ng/ul	0.00
22) 2-Nitrophenol-d4	9.53	143	184833	21.49	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.07	165	363300	18.65	ng/ul	0.00
29) 4-Chloroaniline-d4	10.59	131	453103	23.73	ng/ul	0.00
43) Dimethylphthalate-d6	13.73	166	1180777	18.61	ng/ul	0.00
46) Acenaphthylene-d8	14.00	160	1430025	19.01	ng/ul	0.00
51) 4-Nitrophenol-d4	14.51	143	222751	19.83	ng/ul	0.00
57) Fluorene-d10	15.31	176	999471	18.95	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.43	200	165603	22.85	ng/ul	0.00
70) Anthracene-d10	17.16	188	1563311	19.06	ng/ul	0.00
76) Pyrene-d10	19.46	212	1741305	19.35	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.33	264	1397969	19.18	ng/ul	0.00

Target Compounds

					Ovalue
2) 1,4-Dioxane	3.22	88	54389	5.40 ng/uL	100
4) Benzaldehyde	6.81	77	100673m	28.00 ng/ul	100
6) Phenol	6.86	94	511450	19.04 ng/ul	100
8) Bis(2-Chloroethyl) ether	7.10	93	378590	18.79 ng/ul	100
10) 2-Chlorophenol	7.23	128	370012	18.51 ng/ul	100
11) 2-Methylphenol	8.10	108	374317	17.97 ng/ul	100
12) 2,2'-oxybis(1-Chloropropan	8.20	45	557585	18.86 ng/ul	100
14) Acetophenone	8.48	105	603231	18.79 ng/ul	100
15) N-Nitroso-di-n-propylamine	8.47	70	330453	17.97 ng/ul	100
16) 4-Methylphenol	8.43	108	414160	18.02 ng/ul	100
17) Hexachloroethane	8.74	117	145671	19.16 ng/ul	100
20) Nitrobenzene	8.86	77	458873	20.69 ng/ul	100
21) Isophorone	9.38	82	895491	18.16 ng/ul	100
23) 2-Nitrophenol	9.56	139	208874	21.71 ng/ul	100
24) 2,4-Dimethylphenol	9.63	107	459695	18.75 ng/ul	100
25) Bis(2-Chloroethoxy)methane	9.87	93	541590	18.52 ng/ul	100
27) 2,4-Dichlorophenol	10.09	162	366094	18.98 ng/ul	100
28) Naphthalene	10.50	128	1182054	18.88 ng/ul	100
30) 4-Chloroaniline	10.61	127	486018	24.22 ng/ul	100
31) Hexachlorobutadiene	10.79	225	213344	19.88 ng/ul	100
32) Caprolactam	11.36	113	141408	17.31 ng/ul	100
33) 4-Chloro-3-methylphenol	11.73	107	443711	18.05 ng/ul	100
34) 2-Methylnaphthalene	12.12	142	899597	18.69 ng/ul	100

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 506/28/16

Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA M\DATA\BM062216\
 Data File : BM005910.D
 Acq On : 22 Jun 2016 17:46
 Operator : UM/SJ
 Sample : SST02045
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SST02045

Manual Integrations
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 Quant Title : SVOA CALIBRATION
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 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
36) 1,2,4,5-Tetrachlorobenzene	12.49	216	437652	19.71	ng/ul	100
37) Hexachlorocyclopentadiene	12.47	237	256793	22.17	ng/ul	100
38) 2,4,6-Trichlorophenol	12.73	196	309788	20.17	ng/ul	100
39) 2,4,5-Trichlorophenol	12.80	196	326141	19.46	ng/ul	100
40) 1,1'-Biphenyl	13.14	154	1162658	19.38	ng/ul	100
41) 2-Chloronaphthalene	13.18	162	890773	19.64	ng/ul	100
42) 2-Nitroaniline	13.39	65	314687	22.25	ng/ul	100
44) Dimethylphthalate	13.77	163	1199524	19.27	ng/ul	100
45) 2,6-Dinitrotoluene	13.89	165	239134	22.21	ng/ul	100
47) Acenaphthylene	14.03	152	1524414	19.68	ng/ul	100
48) 3-Nitroaniline	14.22	138	259463	20.87	ng/ul	100
49) Acenaphthene	14.37	153	954695	18.81	ng/ul	100
50) 2,4-Dinitrophenol	14.42	184	105351	23.64	ng/ul	100
52) 4-Nitrophenol	14.53	109	208977	20.32	ng/ul	100
53) Dibenzofuran	14.72	168	1394233	19.34	ng/ul	100
54) 2,4-Dinitrotoluene	14.68	165	352425	22.19	ng/ul	100
55) 2,3,4,6-Tetrachlorophenol	14.94	232	295294	20.26	ng/ul	100
56) Diethylphthalate	15.15	149	1245300	18.61	ng/ul	100
58) Fluorene	15.36	166	1092641	19.24	ng/ul	100
59) 4-Chlorophenyl-phenylether	15.36	204	524763	19.49	ng/ul	100
60) 4-Nitroaniline	15.39	138	286206	20.57	ng/ul	100
63) 4,6-Dinitro-2-methylphenol	15.44	198	184253	23.34	ng/ul	100
64) N-Nitrosodiphenylamine	15.58	169	997860	18.86	ng/ul	100
65) 4-Bromophenyl-phenylether	16.26	248	359068	19.54	ng/ul	100
66) Hexachlorobenzene	16.36	284	390928	19.38	ng/ul	100
67) Atrazine	16.53	200	387998	21.43	ng/ul	100
68) Pentachlorophenol	16.71	266	235480	20.90	ng/ul	100
69) Phenanthrene	17.10	178	1847003	18.85	ng/ul	100
71) Anthracene	17.19	178	1903143	19.25	ng/ul	100
72) Carbazole	17.46	167	1732269	19.90	ng/ul	100
73) Di-n-butylphthalate	18.04	149	2236739	19.09	ng/ul	100
74) Fluoranthene	19.12	202	2209493	21.03	ng/ul	100
77) Pvrene	19.49	202	2280118	20.07	ng/ul	100
78) Butylbenzylphthalate	20.40	149	1006714	19.24	ng/ul	100
79) 3,3'-Dichlorobenzidine	21.17	252	658256	20.54	ng/ul	100
80) Benzo(a)anthracene	21.24	228	2099696	19.05	ng/ul	100
81) Bis(2-ethylhexyl)phthalate	21.19	149	1399088	19.12	ng/ul	100
82) Chrysene	21.29	228	1932739	18.67	ng/ul	100
84) Di-n-octyl phthalate	22.06	149	2252065	22.33	ng/ul	100
85) Benzo(b)fluoranthene	22.81	252	1954294	20.29	ng/ul	100
86) Benzo(k)fluoranthene	22.86	252	1785966	20.06	ng/ul	100
88) Benzo(a)pyrene	23.38	252	1747062	19.80	ng/ul	100
89) Indeno(1,2,3-cd)pyrene	25.69	276	1763674	19.37	ng/ul	100
90) Dibenzo(a,h)anthracene	25.71	278	1457418	19.55	ng/ul	100
91) Benzo(g,h,i)perylene	26.37	276	1485651	19.32	ng/ul	100

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