

```

Data Path   : Z:\HPCHEM1\BNA_G\Data\BG101417\
Data File   : BG029190.D
Acq On      : 13 Oct 2017   11:18
Operator    : SJ/JU
Sample      : SSTD00557
Misc        :
ALS Vial    : 2      Sample Multiplier: 1

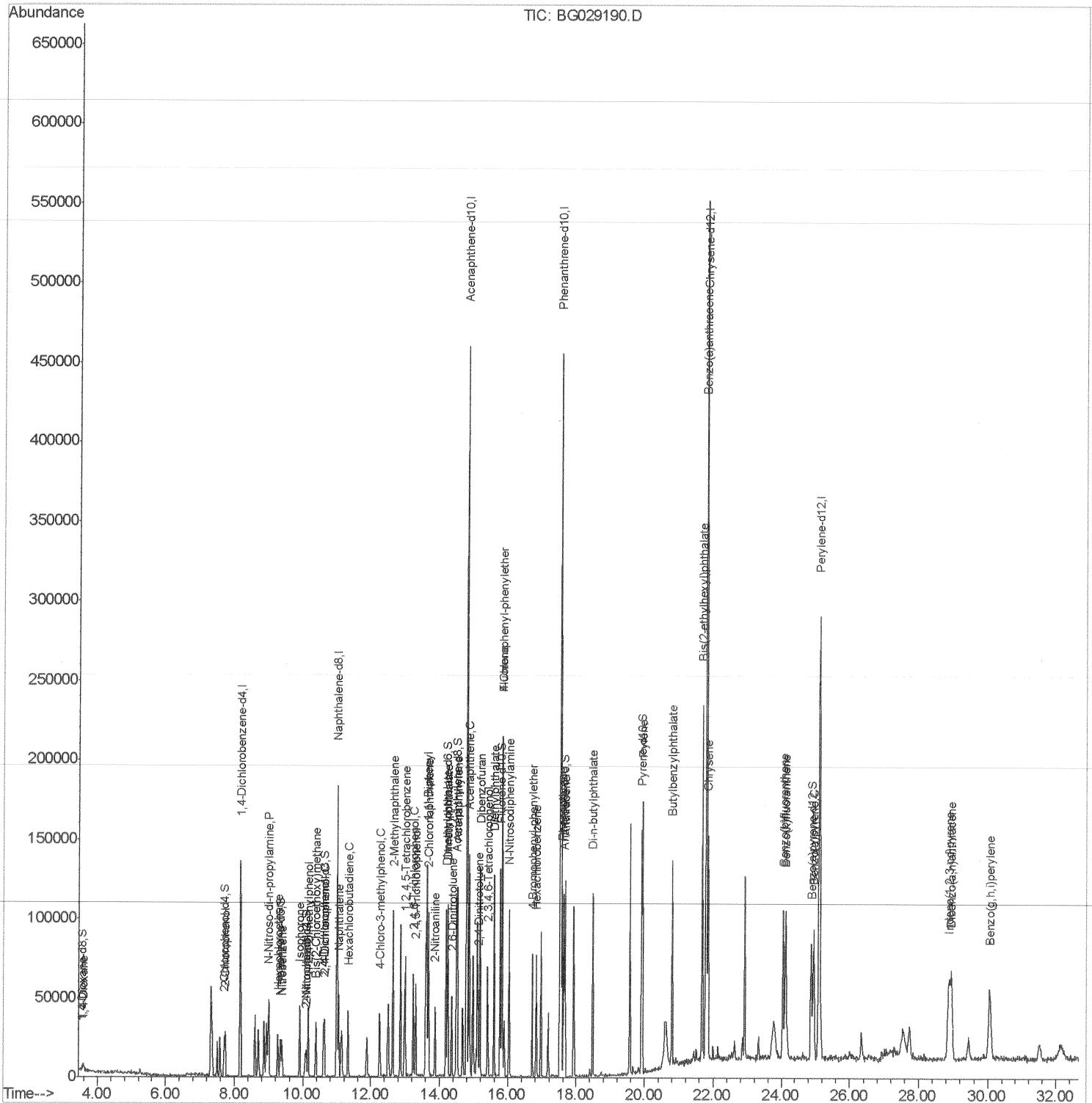
```

Instrument :
BNA_G
ClientSampleId :
SSTD00557

Manual Integrations

Sohil
10/15/2017 12:32:12 PM

Quant Time: Oct 13 17:06:27 2017
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG101417.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Oct 13 16:52:26 2017
Response via : Initial Calibration



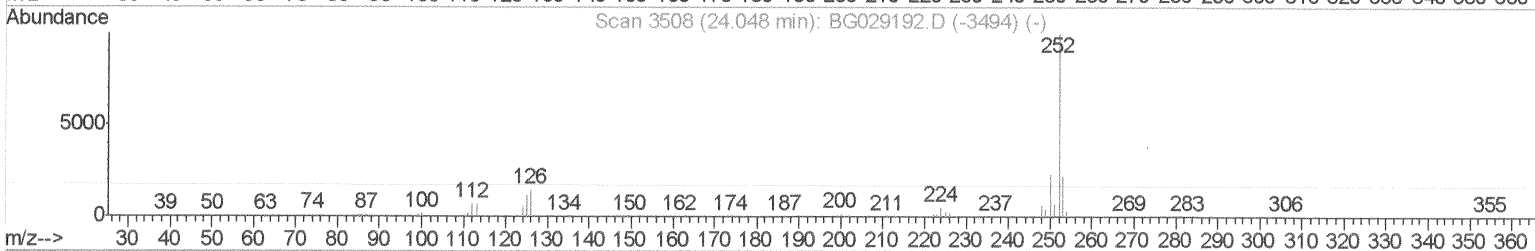
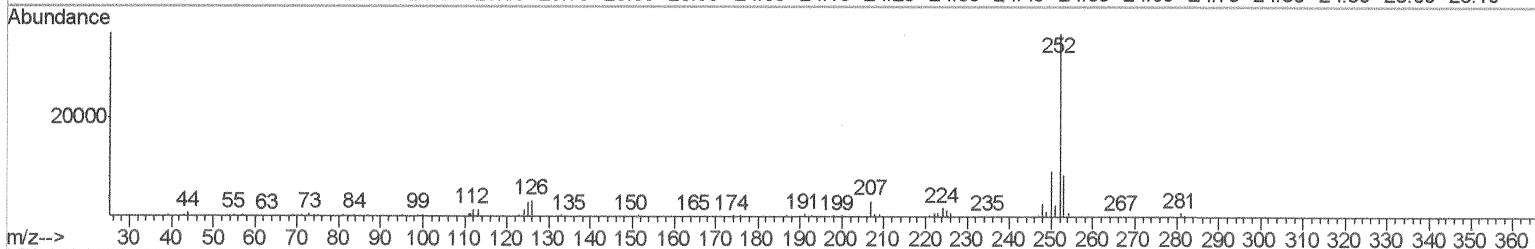
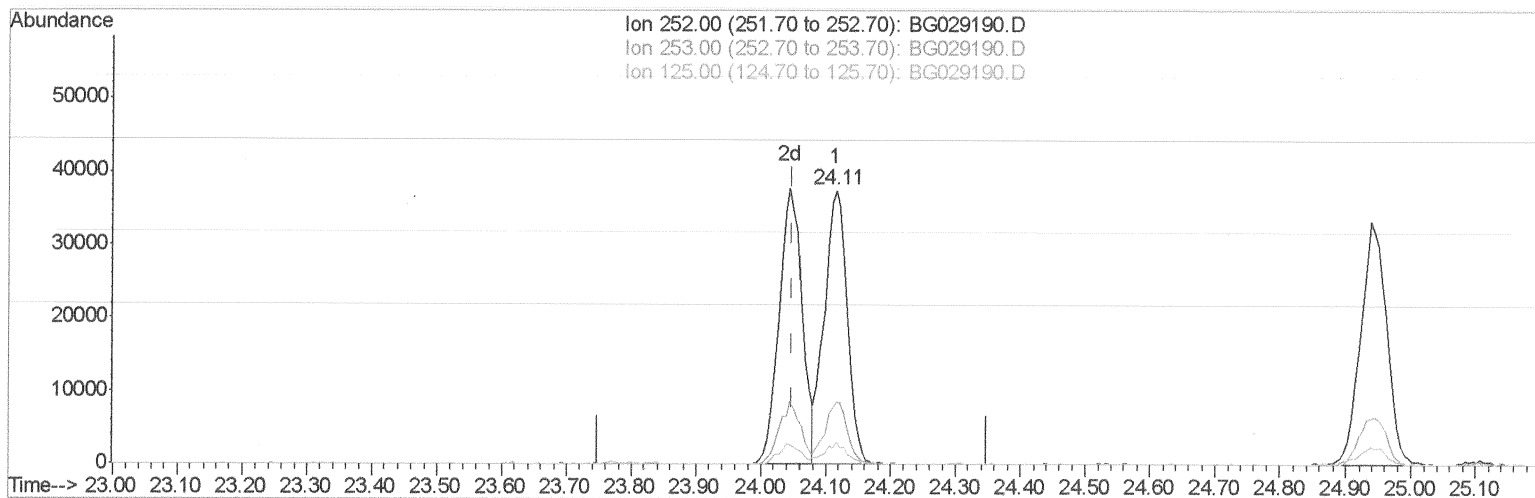
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Instrument :
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ClientSampleId :
SSTD00557

Manual Integrations
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Sohil
10/15/2017 12:32:12 PM

Quant Time: Oct 13 16:54:58 2017
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TIC: BG029190.D

(85) Benzo(b)fluoranthene

24.115min (+0.066) 4.66ng/ul

response 91245

Ion	Exp%	Act%
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252.00	100	100
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253.00	21.20	22.72
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125.00	9.40	8.05
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0.00	0.00	0.00
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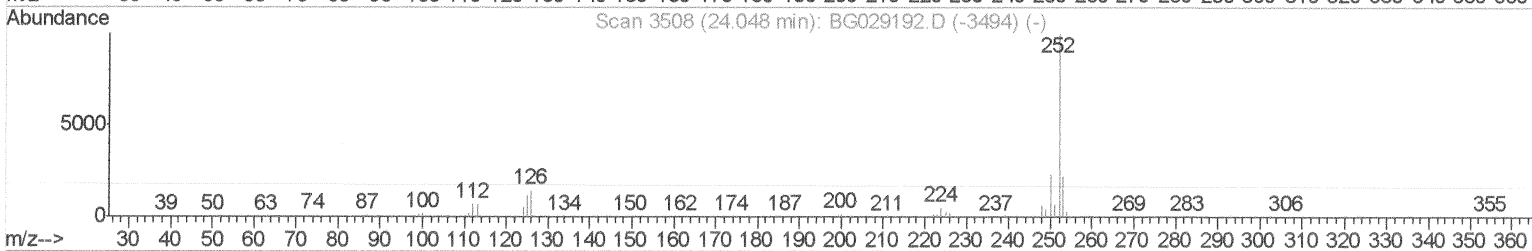
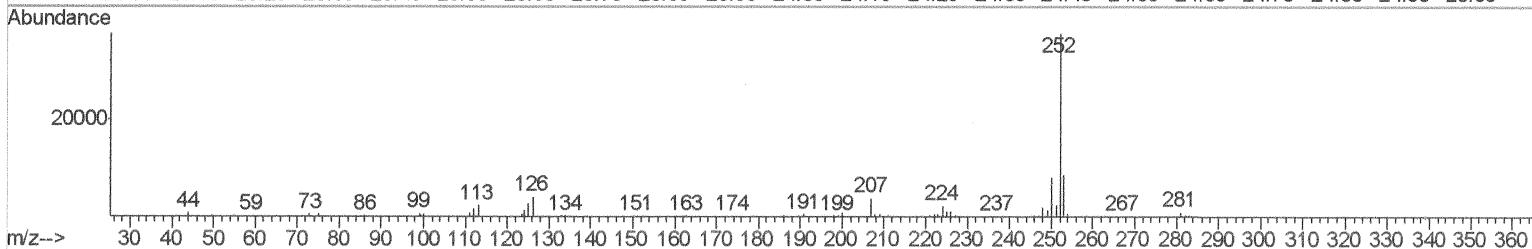
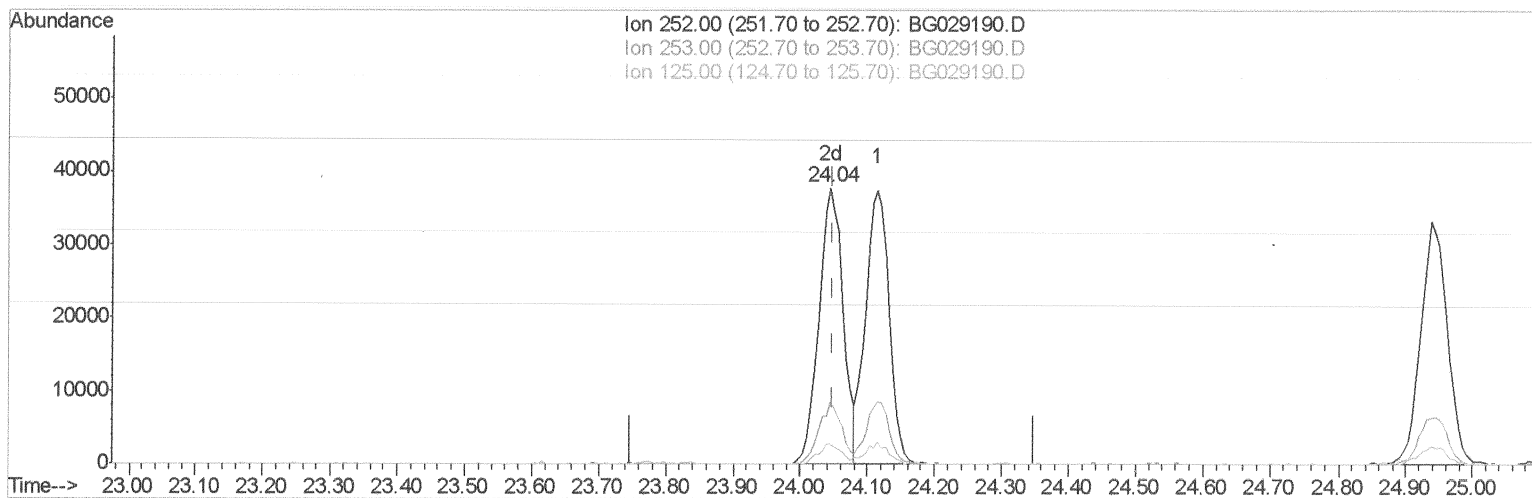
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TIC: BG029190.D

(85) Benzo(b)fluoranthene

24.044min (-0.004) 4.81ng/ul m > SJ 10/15/17

response 94155

Ion	Exp%	Act%
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252.00	100	100
--------	-----	-----

253.00	21.20	22.62
--------	-------	-------

125.00	9.40	7.34#
--------	------	-------

0.00	0.00	0.00
------	------	------

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) 1,4-Dichlorobenzene-d4	8.17	152	35985	20.00	ng/ul	0.00
18) Naphthalene-d8	10.99	136	150179	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.79	164	156812	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.52	188	279533	20.00	ng/ul	0.00
75) Chrysene-d12	21.79	240	304730	20.00	ng/ul	0.00
83) Perylene-d12	25.10	264	299491	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.54	96	1485	4.60	ng/uL	0.00
5) Phenol-d5	0.00	99	0d	0.00	ng/ul	
7) Bis-(2-Chloroethyl) ether-d	0.00	67	0d	0.00	ng/ul	
9) 2-Chlorophenol-d4	7.70	132	11371	4.67	ng/ul	0.00
13) 4-Methylphenol-d8	0.00	113	0d	0.00	ng/ul	
19) Nitrobenzene-d5	9.33	128	4780	4.10	ng/ul	0.00
22) 2-Nitrophenol-d4	10.05	143	4713	4.34	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.60	165	12909	4.80	ng/ul	0.00
29) 4-Chloroaniline-d4	0.00	131	0d	0.00	ng/ul	
43) Dimethylphthalate-d6	14.19	166	70155	5.36	ng/ul	0.00
46) Acenaphthylene-d8	14.48	160	84450	5.15	ng/ul	0.00
51) 4-Nitrophenol-d4	0.00	143	0d	0.00	ng/ul	
57) Fluorene-d10	15.77	176	48673	4.21	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	0.00	200	0d	0.00	ng/ul	
70) Anthracene-d10	17.62	188	67315	5.02	ng/ul	0.00
76) Pyrene-d10	19.89	212	78865	4.79	ng/ul	0.00
87) Benzo(a)pyrene-d12	24.87	264	74337	4.84	ng/ul	-0.01

Target Compounds

					Qvalue	
2) 1,4-Dioxane	3.59	88	2304	5.536	ng/uL#	83
10) 2-Chlorophenol	7.73	128	11680	4.720	ng/ul	95
15) N-Nitroso-di-n-propylamine	8.97	70	12179	4.188	ng/ul#	96
17) Hexachloroethane	9.26	117	4952	4.139	ng/ul#	78
20) Nitrobenzene	9.38	77	13727	4.083	ng/ul#	97
21) Isophorone	9.90	82	32859	4.821	ng/ul	95
23) 2-Nitrophenol	10.09	139	5015	4.088	ng/ul#	89
24) 2,4-Dimethylphenol	10.14	107	15653	4.681	ng/ul	92
25) Bis(2-Chloroethoxy)methane	10.38	93	18895	4.877	ng/ul	97
27) 2,4-Dichlorophenol	10.62	162	11995	4.672	ng/ul	93
28) Naphthalene	11.04	128	39991	4.913	ng/ul	99
31) Hexachlorobutadiene	11.32	225	9895	4.829	ng/ul	91
33) 4-Chloro-3-methylphenol	12.25	107	14736	4.677	ng/ul	89
34) 2-Methylnaphthalene	12.63	142	47930	6.538	ng/ul	90
36) 1,2,4,5-Tetrachlorobenzene	13.00	216	21842	5.063	ng/ul#	94
38) 2,4,6-Trichlorophenol	13.23	196	15223	5.201	ng/ul	95
39) 2,4,5-Trichlorophenol	13.30	196	16169	5.200	ng/ul	97
40) 1,1'-Biphenyl	13.63	154	68210	6.171	ng/ul	99
41) 2-Chloronaphthalene	13.67	162	50434	5.938	ng/ul	94
42) 2-Nitroaniline	13.86	65	13198	4.939	ng/ul	96
44) Dimethylphthalate	14.23	163	69893	5.281	ng/ul	98
45) 2,6-Dinitrotoluene	14.34	165	9005	4.200	ng/ul#	91
47) Acenaphthylene	14.51	152	87793	5.327	ng/ul	99

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49) Acenaphthene	14.85	153	57414	4.975	ng/ul	100
53) Dibenzofuran	15.18	168	82578	4.942	ng/ul	98
54) 2,4-Dinitrotoluene	15.14	165	12968	3.757	ng/ul#	98
55) 2,3,4,6-Tetrachlorophenol	15.40	232	13682	4.198	ng/ul#	99
56) Diethylphthalate	15.59	149	70784	4.575	ng/ul	99
58) Fluorene	15.83	166	51159	4.046	ng/ul	95
59) 4-Chlorophenyl-phenylether	15.82	204	26472	4.215	ng/ul	96
64) N-Nitrosodiphenylamine	16.02	169	39378	4.704	ng/ul	98
65) 4-Bromophenyl-phenylether	16.71	248	14924	4.931	ng/ul	97
66) Hexachlorobenzene	16.83	284	16241	5.093	ng/ul	95
69) Phenanthrene	17.56	178	74604	4.956	ng/ul	98
71) Anthracene	17.66	178	77858	5.003	ng/ul	98
73) Di-n-butylphthalate	18.47	149	81475	4.798	ng/ul	98
77) Pyrene	19.92	202	104284	4.912	ng/ul	99
78) Butylbenzylphthalate	20.80	149	37395	4.314	ng/ul	98
80) Benzo(a)anthracene	21.78	228	101851	5.094	ng/ul	97
81) Bis(2-ethylhexyl)phthalate	21.67	149	57306	4.576	ng/ul	99
82) Chrysene	21.84	228	91013	5.049	ng/ul	97
85) Benzo(b)fluoranthene	24.04	252	94155m>	4.811	ng/ul	97
86) Benzo(k)fluoranthene	24.11	252	91245	4.831	ng/ul	97
88) Benzo(a)pyrene	24.94	252	92172	4.991	ng/ul#	95
89) Indeno(1,2,3-cd)pyrene	28.87	276	103993	5.477	ng/ul#	96
90) Dibenzo(a,h)anthracene	28.94	278	88718	5.525	ng/ul#	93
91) Benzo(g,h,i)perylene	30.05	276	88498	5.566	ng/ul	94

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