

Quantitation Report (Qedit)

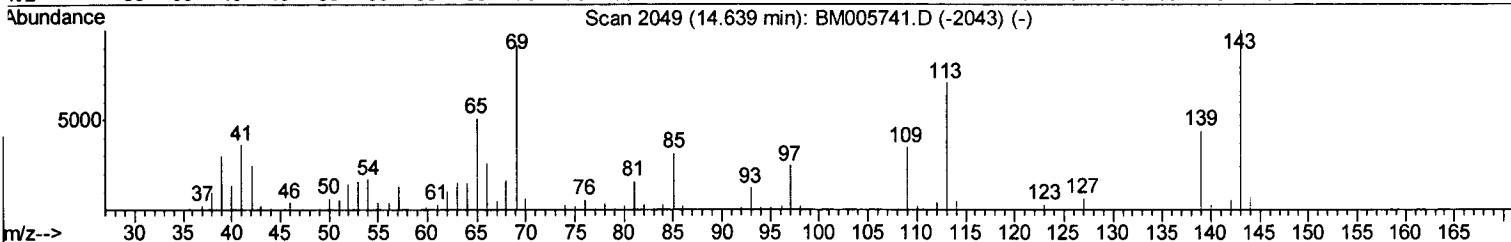
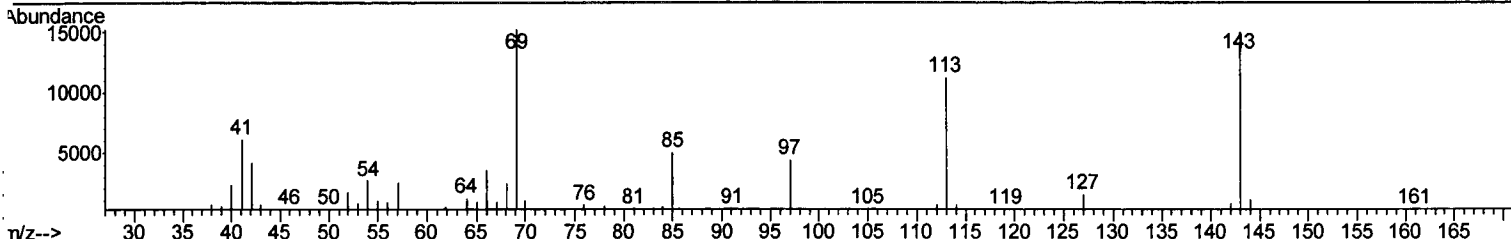
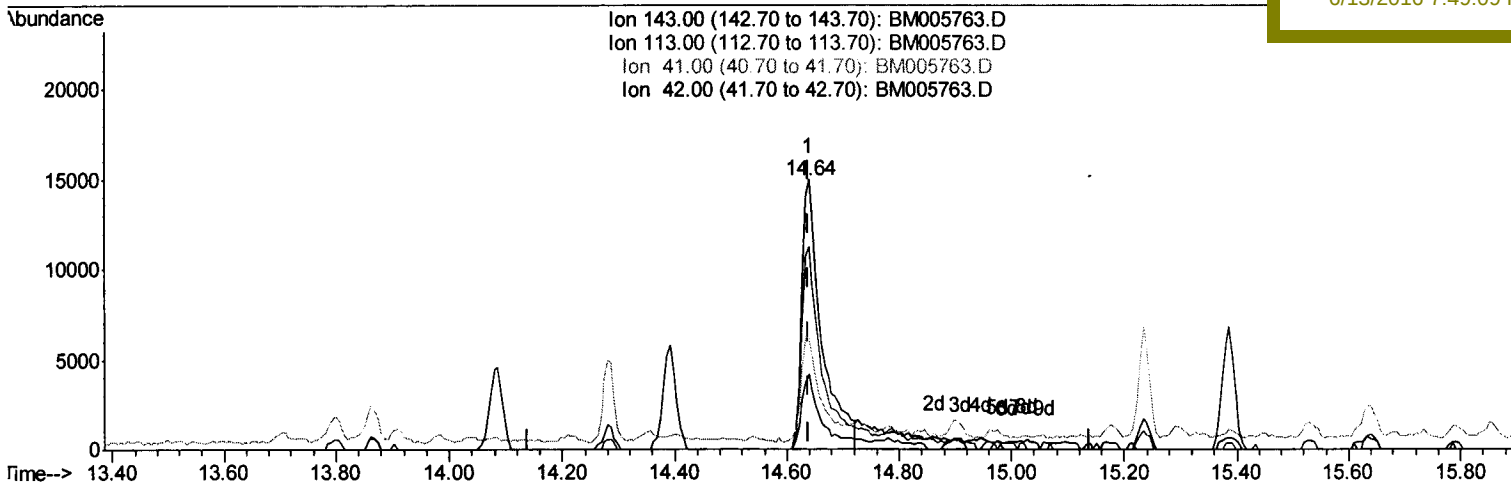
Data Path : Z:\HPCHEM1\BNA_M\DATA\BM061216\
 Data File : BM005763.D
 Acq On : 12 Jun 2016 23:51
 Operator : UM/SJ
 Sample : H3536-11
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 D9Y21

Quant Time: Jun 13 03:32:52 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM061016.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jun 13 03:29:45 2016
 Response via : Initial Calibration

Manual Integrations
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TIC: BM005763.D

(20) 4-Nitrophenol-d4 (S)

14.639min (-0.000) 9.90ng/ul

response 37008

Ion	Exp%	Act%
143.00	100	100
113.00	72.40	75.18
41.00	38.30	40.90
42.00	26.80	27.68

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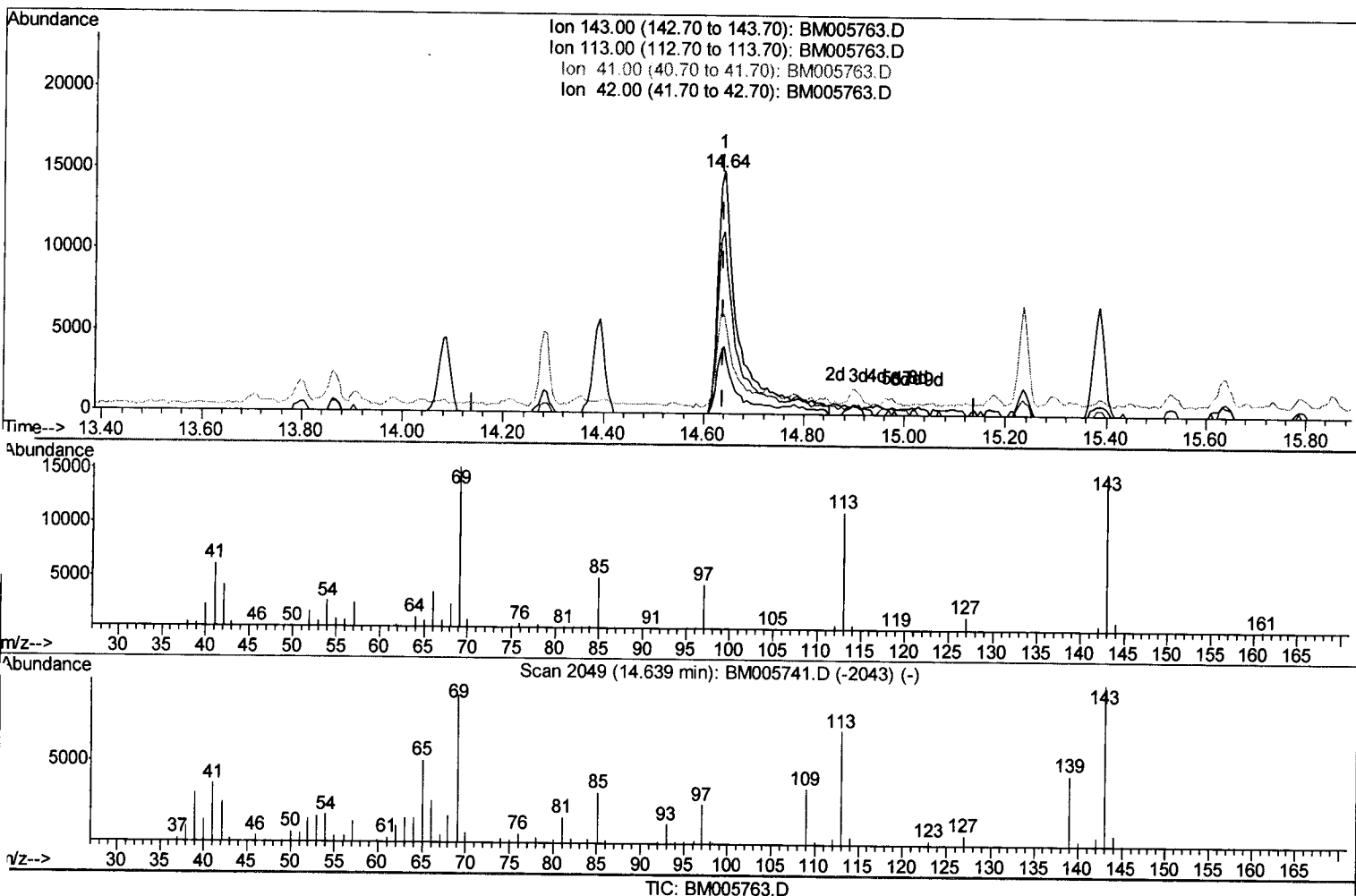
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(20) 4-Nitrophenol-d4 (S)

14.639min (-0.000) 11.88ng/ul m

response 44416

Ion	Exp%	Act%
143.00	100	100
113.00	72.40	75.18
41.00	38.30	40.90
42.00	26.80	27.68

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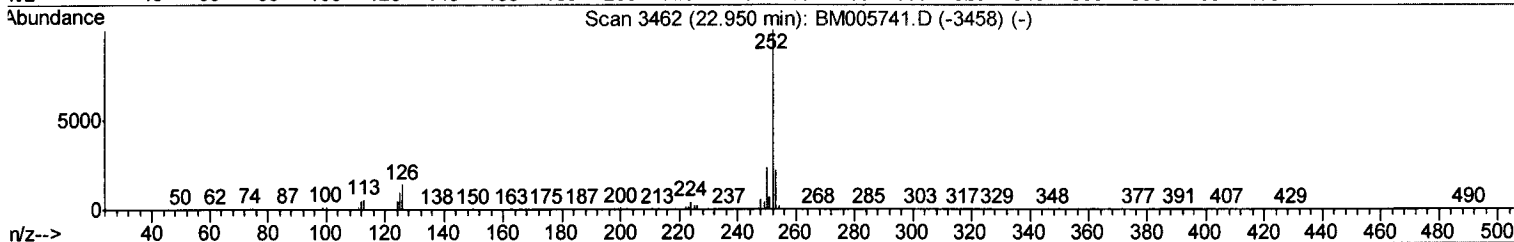
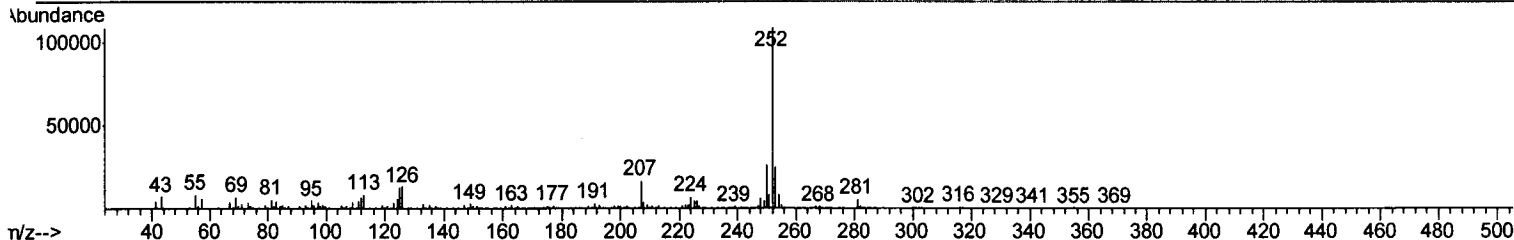
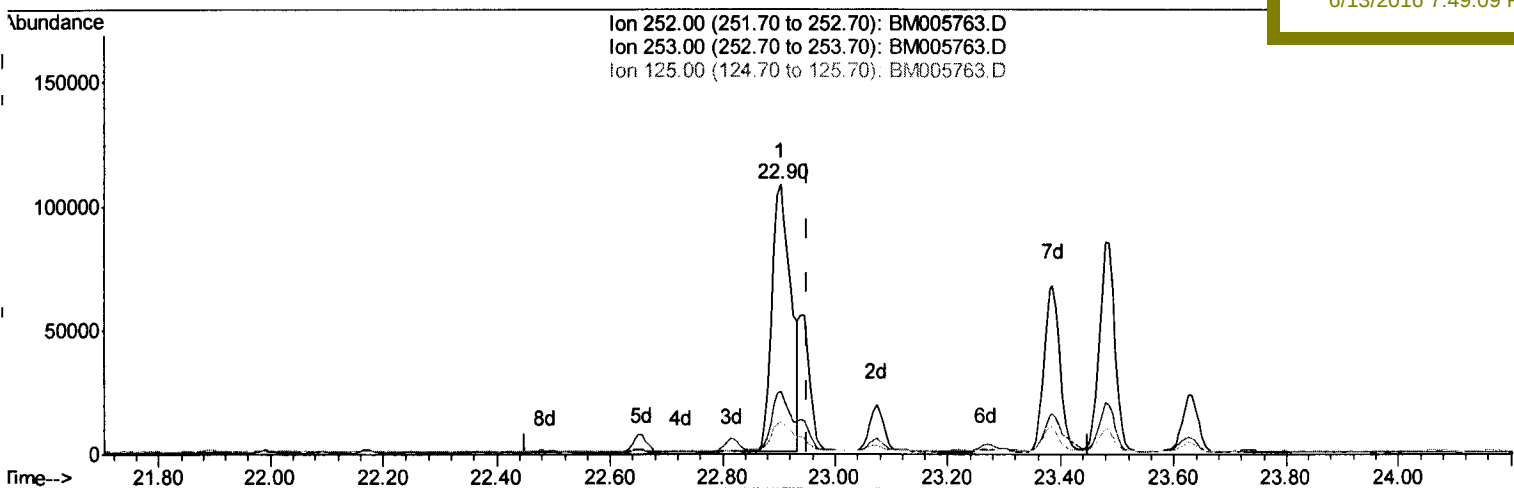
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TIC: BM005763.D

(36) Benzo(k)fluoranthene

22.903min (-0.047) 9.51ng/ul

response 256956

Ion	Exp%	Act%
252.00	100	100
253.00	21.90	23.06
125.00	10.20	11.78
0.00	0.00	0.00

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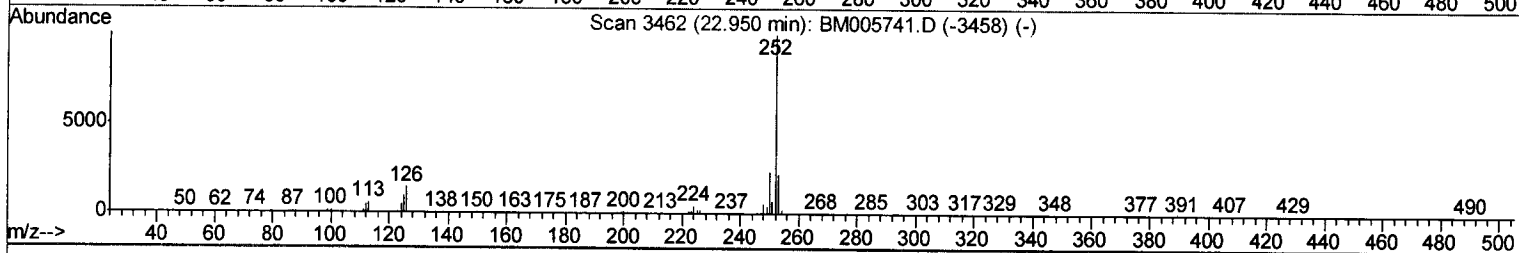
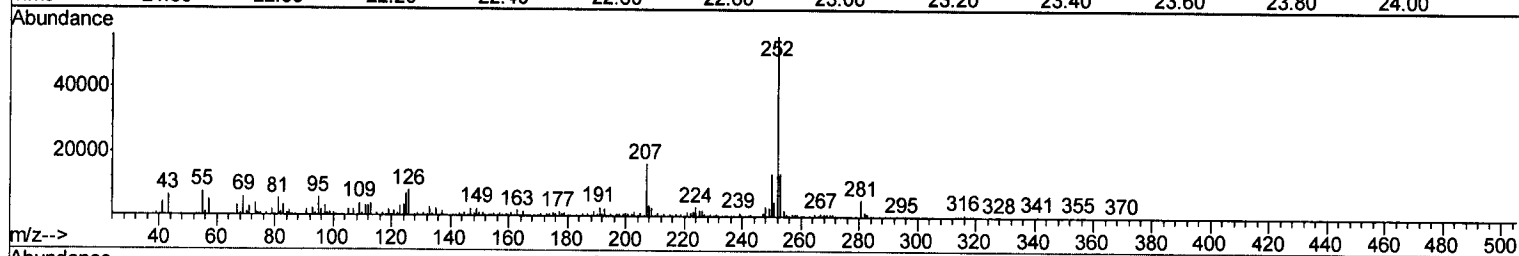
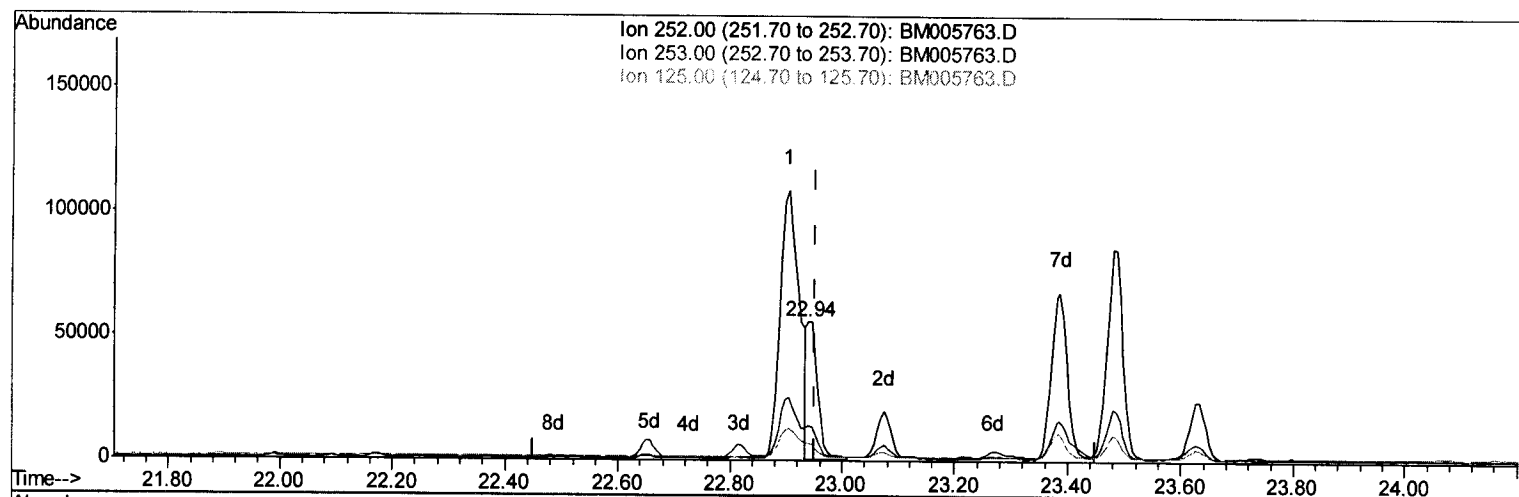
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TIC: BM005763.D

(36) Benzo(k)fluoranthene

22.938min (-0.012) 2.74ng/ul m

response 74050

Ion	Exp%	Act%
252.00	100	100
253.00	21.90	24.58
125.00	10.20	12.77#
0.00	0.00	0.00

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Quant Time: Jun 13 04:09:22 2016
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.75	152	115052	20.00	ng/ul	0.00
7) Naphthalene-d8	10.54	136	567294	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.39	164	309796	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.14	188	539286	20.00	ng/ul	0.00
29) Chrysene-d12	21.32	240	426181	20.00	ng/ul	0.00
34) Perylene-d12	23.58	264	450124	20.00	ng/ul	0.00

System Monitoring Compounds

2) 1,4-Dioxane-d8	3.19	96	3827	1.44	ng/uL	0.00
3) Phenol-d5	6.93	99	167189	16.84	ng/ul	0.00
4) Bis-(2-Chloroethyl) ether-d	7.08	67	112407	18.84	ng/ul	0.00
5) 2-Chlorophenol-d4	7.28	132	155900	19.42	ng/ul	0.00
6) 4-Methylphenol-d8	8.46	113	119879	14.29	ng/ul	0.00
8) Nitrobenzene-d5	8.90	128	79231	19.38	ng/ul	0.00
9) 2-Nitrophenol-d4	9.63	143	89358	19.33	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.17	165	165616	19.75	ng/ul	0.00
12) 4-Chloroaniline-d4	10.68	131	73001	7.22	ng/ul	0.00
16) Dimethylphthalate-d6	13.80	166	540437	21.15	ng/ul	0.00
17) Acenaphthylene-d8	14.09	160	681079	21.79	ng/ul	0.00
20) 4-Nitrophenol-d4	14.64	143	44416m	11.88	ng/ul	0.00
21) Fluorene-d10	15.39	176	447582	20.69	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.53	200	16475	5.91	ng/ul	0.00
26) Anthracene-d10	17.24	188	556524	21.91	ng/ul	0.00
30) Pyrene-d10	19.53	212	490255	23.94	ng/ul	0.00
37) Benzo(a)pyrene-d12	23.44	264	468653	22.83	ng/ul	0.00

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Target Compounds

						Qvalue
25) Phenanthrene	17.18	178	177139	6.00	ng/ul	99
27) Anthracene	17.27	178	40302	1.35	ng/ul	99
28) Fluoranthene	19.20	202	323095	9.32	ng/ul	98
31) Pyrene	19.56	202	267348	10.31	ng/ul	99
32) Benzo(a)anthracene	21.30	228	162509	6.60	ng/ul	98
33) Chrysene	21.36	228	160450	6.61	ng/ul	98
35) Benzo(b)fluoranthene	22.90	252	256956	9.61	ng/ul	98
36) Benzo(k)fluoranthene	22.94	252	74050m	2.74	ng/ul	
38) Benzo(a)pyrene	23.48	252	160875	6.34	ng/ul	96
39) Indeno(1,2,3-cd)pyrene	25.87	276	126751	5.11	ng/ul#	94
40) Dibenzo(a,h)anthracene	25.86	278	33423	1.61	ng/ul#	80
41) Benzo(g,h,i)perylene	26.57	276	120617	5.82	ng/ul	100

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

(QT Reviewed)

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