

Quantitation Report (Qedit)

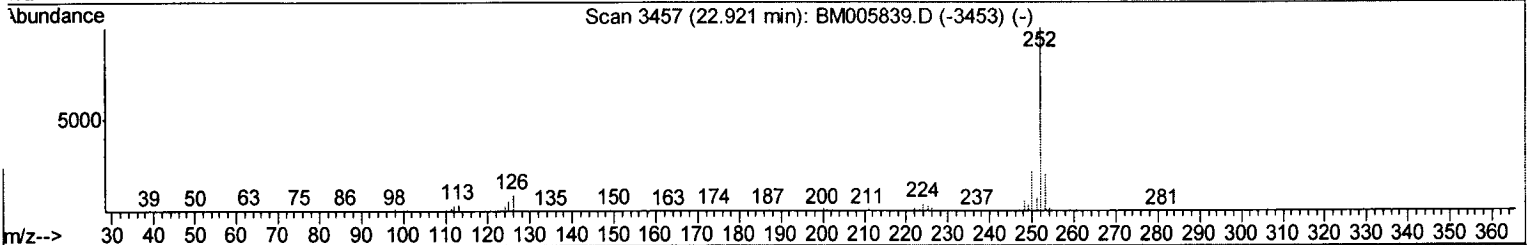
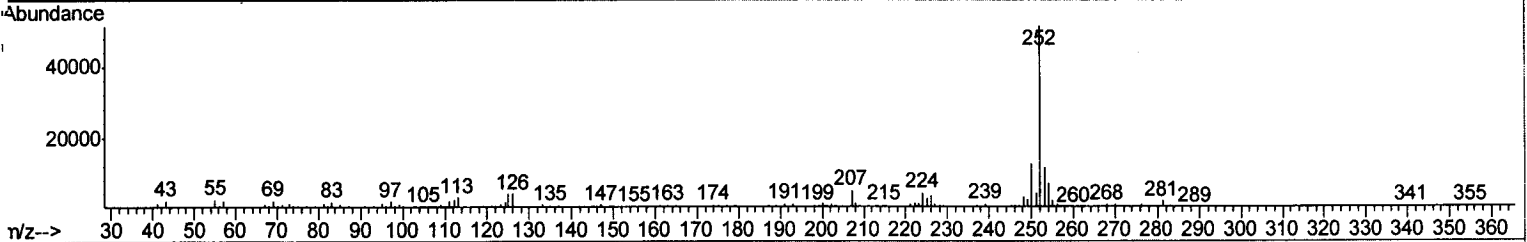
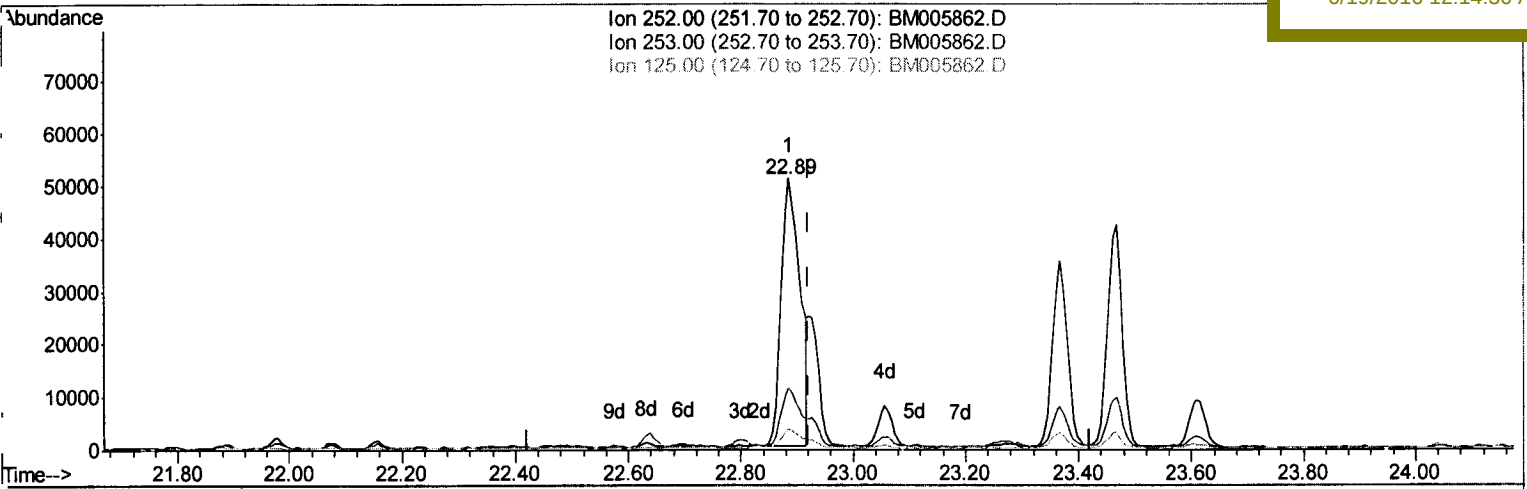
Data Path : Z:\HPCHEM1\BNA_M\Data\BM061816\
 Data File : BM005862.D
 Acq On : 17 Jun 2016 15:22
 Operator : UM/SJ
 Sample : H3537-18
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 D9Y48

Quant Time: Jun 17 15:54:04 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM061716.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jun 17 10:48:37 2016
 Response via : Initial Calibration

Manual Integrations
 APPROVED

umangi
 6/19/2016 12:14:36 AM



TIC: BM005862.D

(36) Benzo(k)fluoranthene

22.886min (-0.035) 11.08ng/ul

response 115474

Ion	Exp%	Act%
252.00	100	100
253.00	21.90	22.54
125.00	6.00	7.85#
0.00	0.00	0.00

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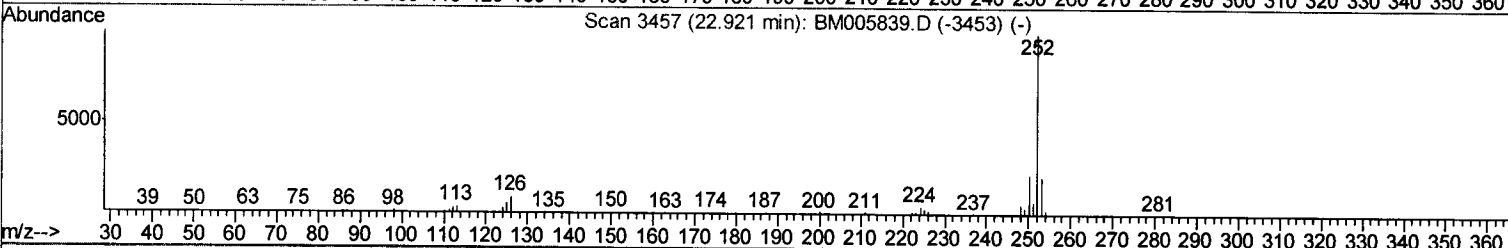
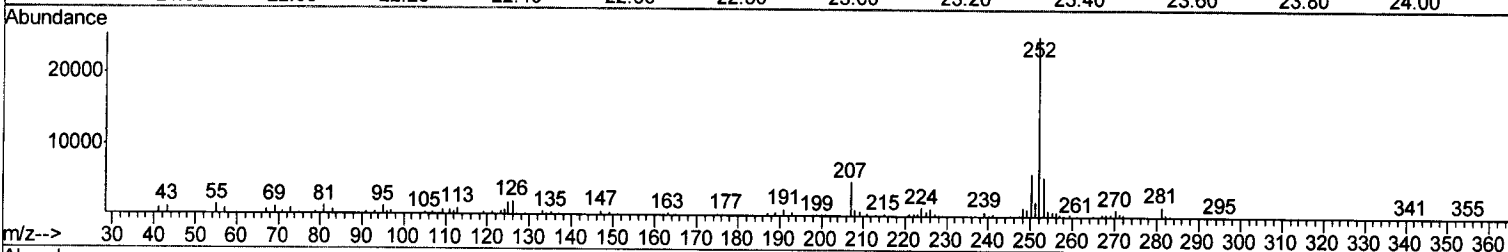
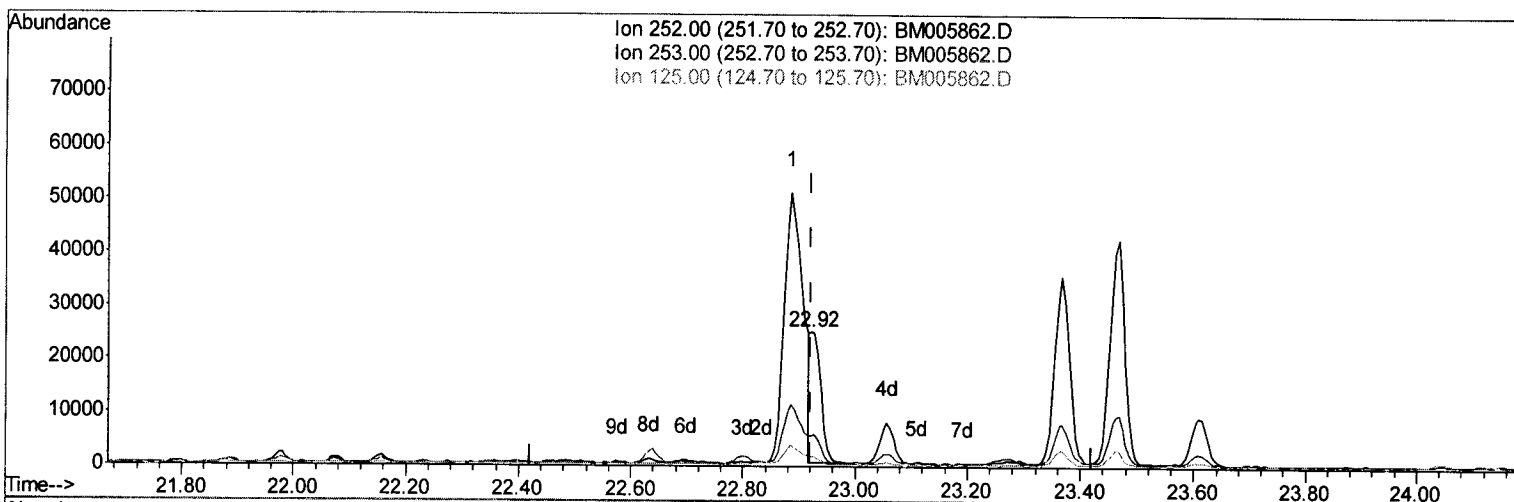
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TIC: BM005862.D

(36) Benzo(k)fluoranthene

22.921min (+0.000) 3.16ng/ul m

response 32984

Ion Exp% Act%

252.00 100 100

253.00 21.90 23.17

125.00 6.00 7.87#

0.00 0.00 0.00

> U.M
 06/20/16

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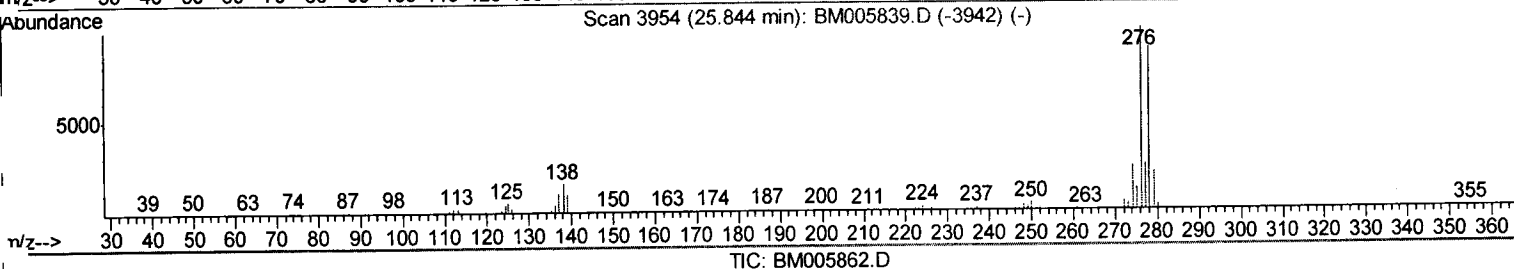
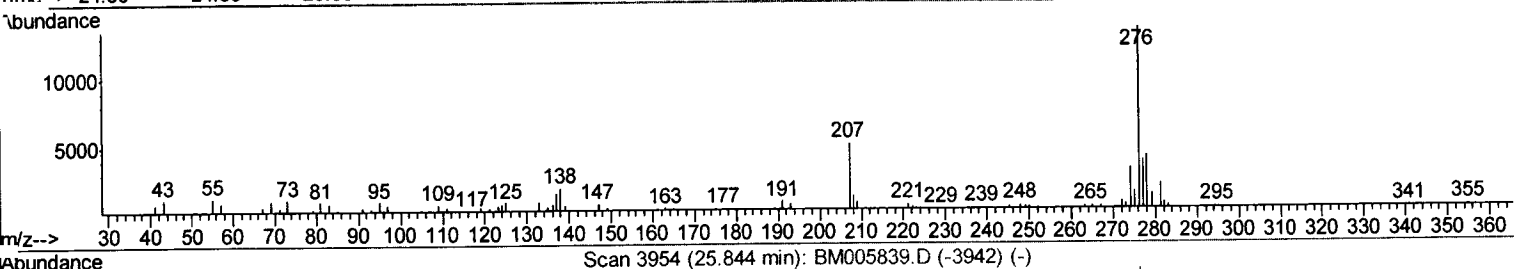
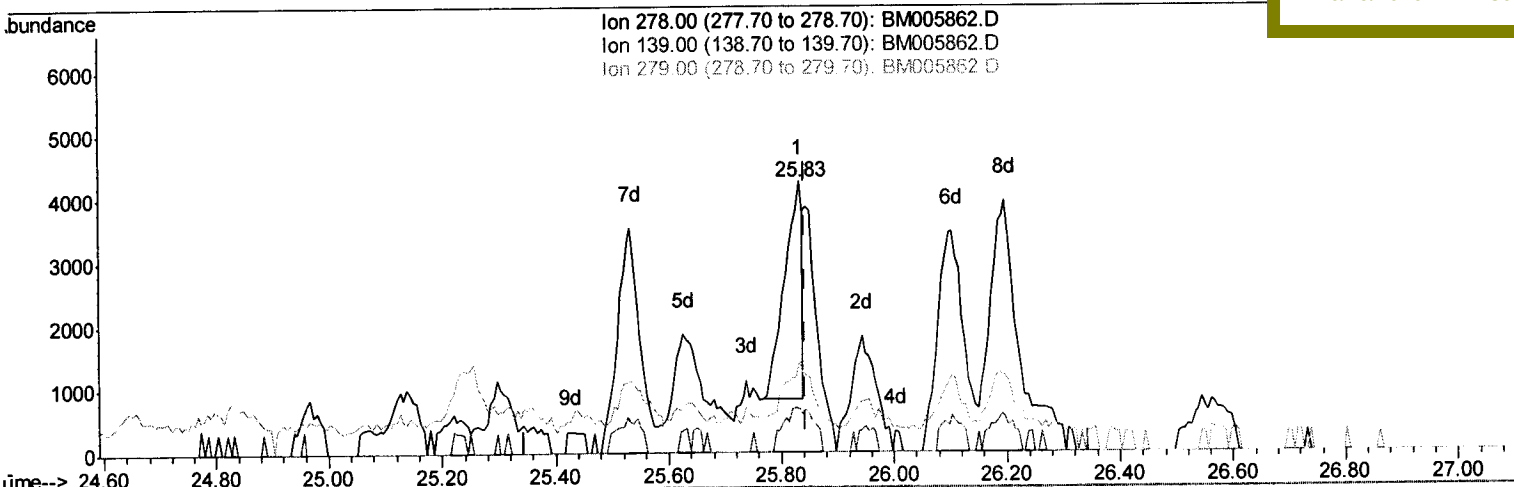
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Quant Time: Jun 17 15:56:44 2016
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(40) Dibenzo(a,h)anthracene

25.833min (-0.012) 0.70ng/ul

response 7414

Ion	Exp%	Act%
278.00	100	100
139.00	10.30	16.08#
279.00	23.20	33.12#
0.00	0.00	0.00

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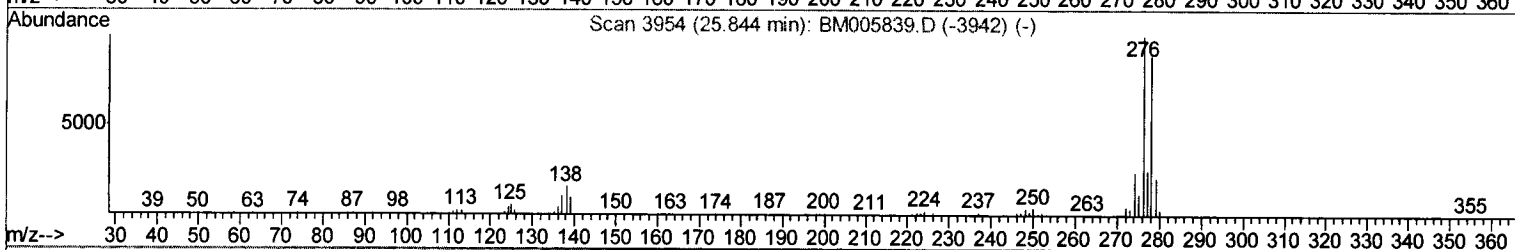
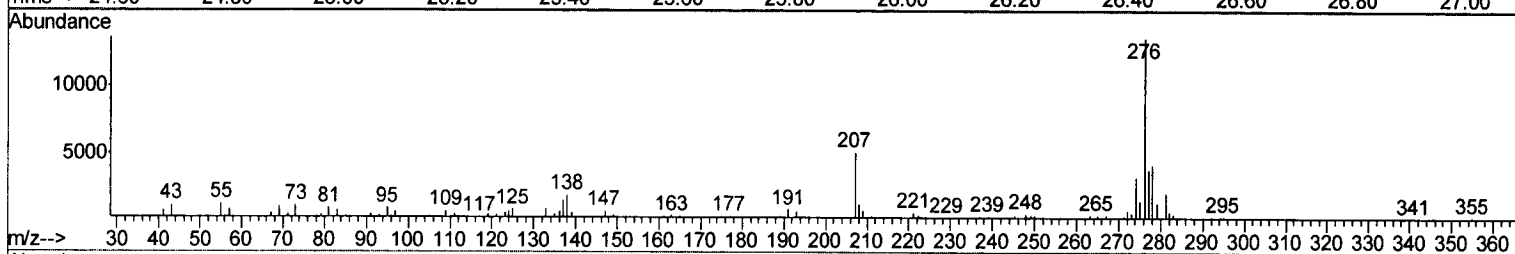
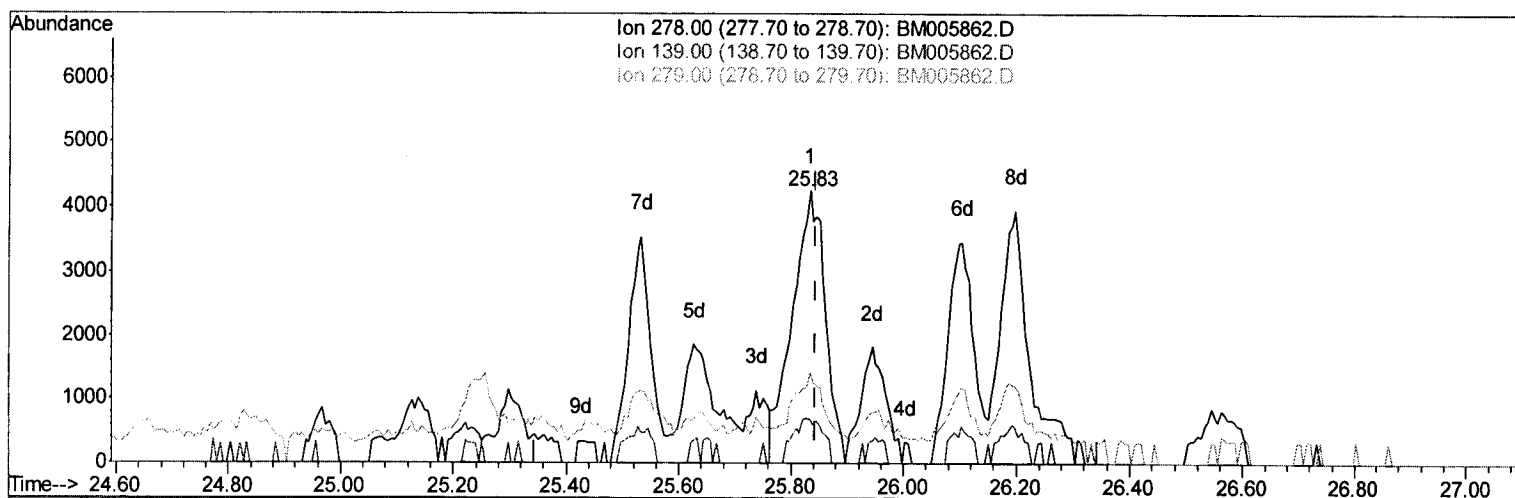
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TIC: BM005862.D

(40) Dibenzo(a,h)anthracene

25.833min (-0.012) 1.63ng/ul m *U.M*
 response 17322
06/20/16

Ion	Exp%	Act%
278.00	100	100
139.00	10.30	16.08#
279.00	23.20	33.12#
0.00	0.00	0.00

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Quant Time: Jun 17 16:00:12 2016
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 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	32262	20.00	ng/ul	0.00
7) Naphthalene-d8	10.52	136	144313	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.37	164	88718	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.12	188	205106	20.00	ng/ul	0.00
29) Chrysene-d12	21.31	240	196720	20.00	ng/ul	0.00
34) Perylene-d12	23.56	264	173405	20.00	ng/ul	0.00

System Monitoring Compounds

2) 1,4-Dioxane-d8	3.18	96	685	2.75	ng/uL	0.00
3) Phenol-d5	6.92	99	40687	15.06	ng/ul	0.00
4) Bis-(2-Chloroethyl) ether-d	7.06	67	22136	15.81	ng/ul	0.00
5) 2-Chlorophenol-d4	7.27	132	38836	17.24	ng/ul	0.00
6) 4-Methylphenol-d8	8.45	113	33492	13.87	ng/ul	0.00
8) Nitrobenzene-d5	8.89	128	18509	18.17	ng/ul	0.00
9) 2-Nitrophenol-d4	9.61	143	22848	18.36	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.16	165	42434	17.26	ng/ul	0.00
12) 4-Chloroaniline-d4	10.67	131	13285	7.99	ng/ul	0.00
16) Dimethylphthalate-d6	13.79	166	142658	17.00	ng/ul	0.00
17) Acenaphthylene-d8	14.07	160	176244	18.23	ng/ul	0.00
20) 4-Nitrophenol-d4	14.63	143	8719	8.62	ng/ul	0.02
21) Fluorene-d10	15.37	176	127544	17.83	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.52	200	10803	7.81	ng/ul	0.00
26) Anthracene-d10	17.22	188	195720	18.28	ng/ul	0.00
30) Pyrene-d10	19.52	212	219458	23.54	ng/ul	0.00
37) Benzo(a)pyrene-d12	23.42	264	163990	18.92	ng/ul	0.00

Target Compounds

						Qvalue
25) Phenanthrene	17.16	178	62494	5.12	ng/ul	99
27) Anthracene	17.26	178	38475	3.08	ng/ul	98
28) Fluoranthene	19.18	202	201223	13.11	ng/ul	99
31) Pyrene	19.54	202	286212	24.62	ng/ul	99
32) Benzo(a)anthracene	21.29	228	125851	10.24	ng/ul	97
33) Chrysene	21.34	228	151785	12.83	ng/ul	96
35) Benzo(b)fluoranthene	22.89	252	115474	10.70	ng/ul#	97
36) Benzo(k)fluoranthene	22.92	252	32984m	3.16	ng/ul	
38) Benzo(a)pyrene	23.47	252	75987	7.17	ng/ul	99
39) Indeno(1,2,3-cd)pyrene	25.84	276	44345	3.42	ng/ul#	96
40) Dibenzo(a,h)anthracene	25.83	278	17322m	1.63	ng/ul	
41) Benzo(g,h,i)perylene	26.54	276	46935	4.26	ng/ul	96

} U.M
 06/20/16

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

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