

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG081519\

Data File : BG042229.D

Acq On : 15 Aug 2019 14:58

Operator : HP/JU

Sample : K4183-17

Misc :

ALS Vial : 12 Sample Multiplier: 1

Quant Time: Aug 15 19:30:54 2019

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080319MA.M

Quant Title : SVOA CALIBRATION

QLast Update : Wed Aug 14 10:42:59 2019

Response via : Initial Calibration

Instrument :

BNA_G

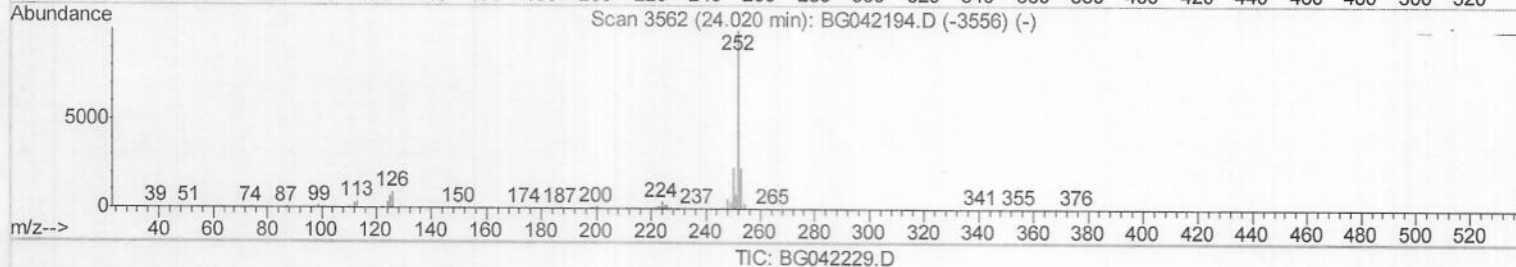
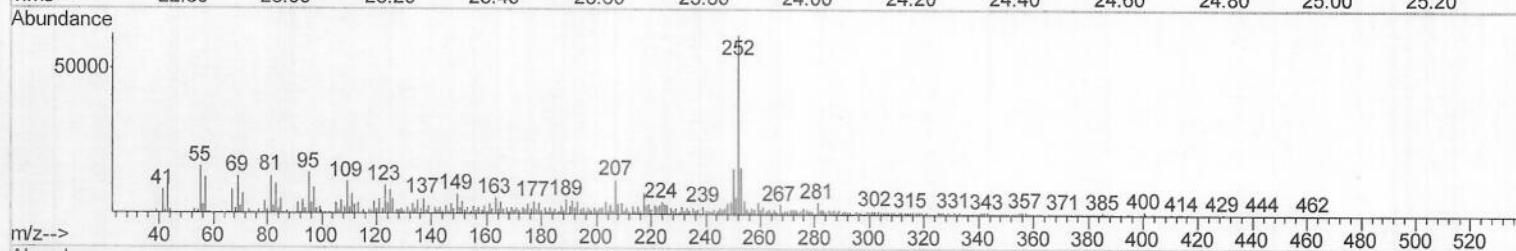
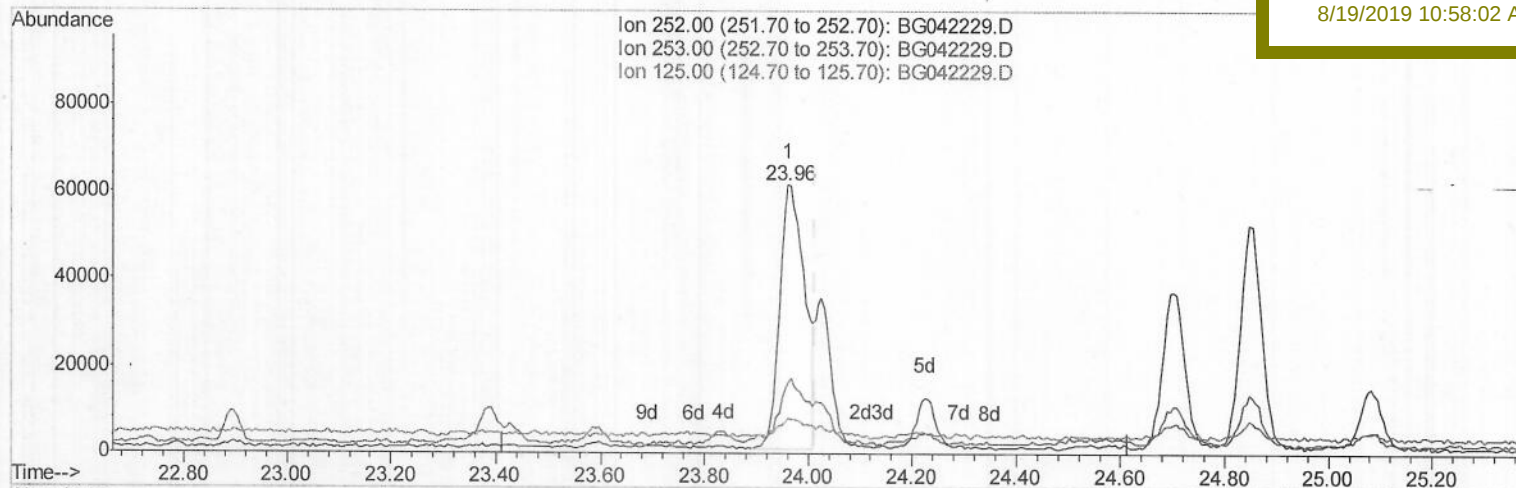
ClientSampleId :

C0AG4

Manual Integrations
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(88) Benzo(k)fluoranthene

23.961min (-0.052) 6.74ng/ul

response 202384

Ion	Exp%	Act%
252.00	100	100
253.00	22.90	26.54
125.00	6.90	13.04#
0.00	0.00	0.00

Quantitation Report (Qedit)

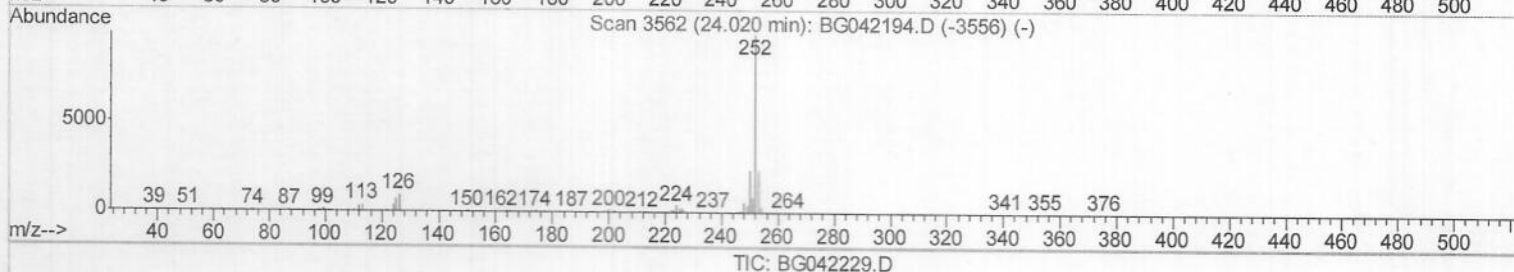
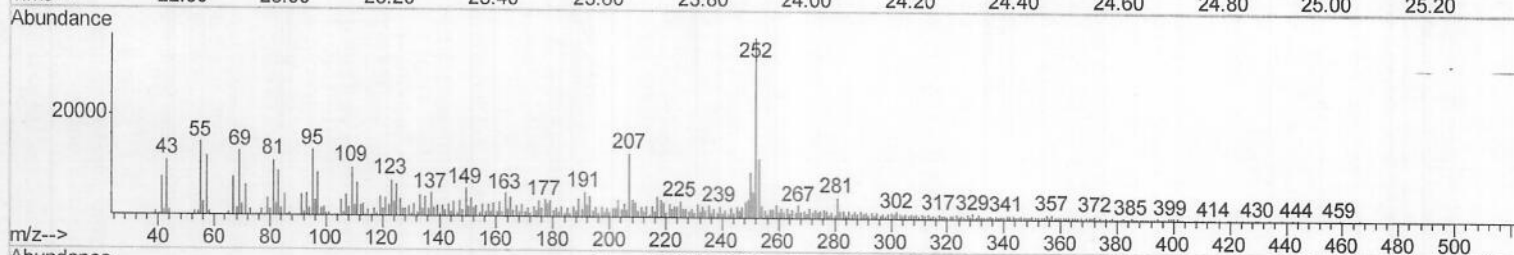
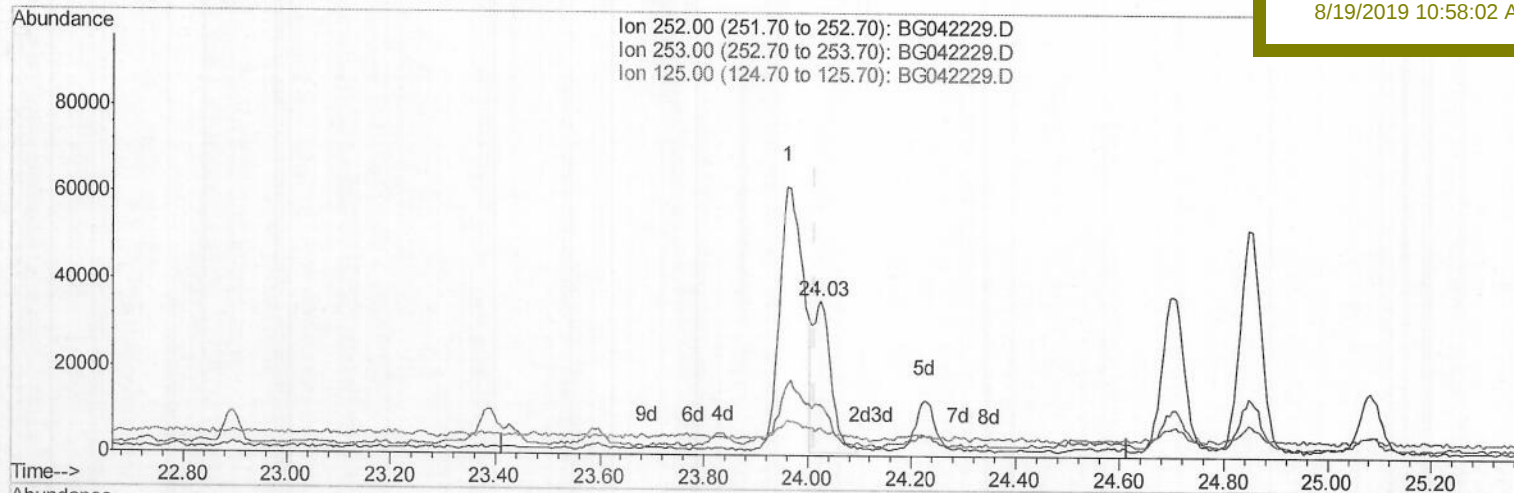
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(88) Benzo(k)fluoranthene

24.025min (+0.012) 2.96ng/ul m

response 88926

Ion	Exp%	Act%
252.00	100	100
253.00	22.90	33.19#
125.00	6.90	17.89#
0.00	0.00	0.00

50
 08/20/19

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Misc :

ALS Vial : 12 Sample Multiplier: 1

Quant Time: Aug 16 12:16:39 2019

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SOM-EPA-BG080319MA.M

Quant Title : SVOA CALIBRATION

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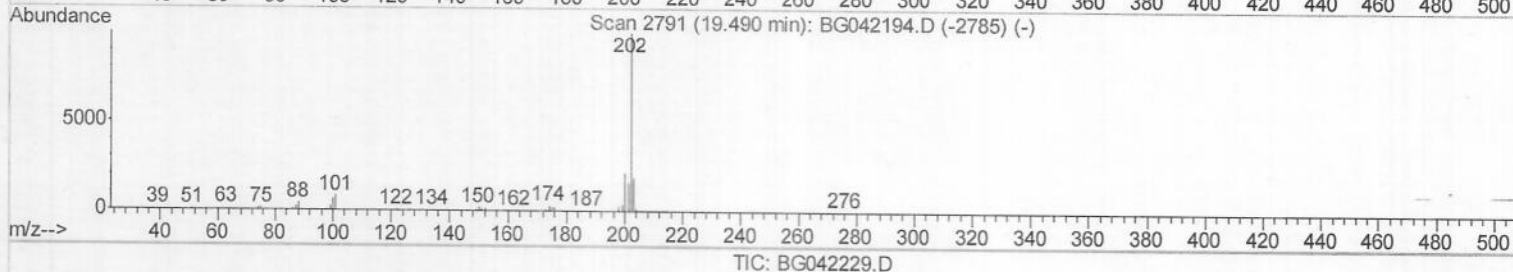
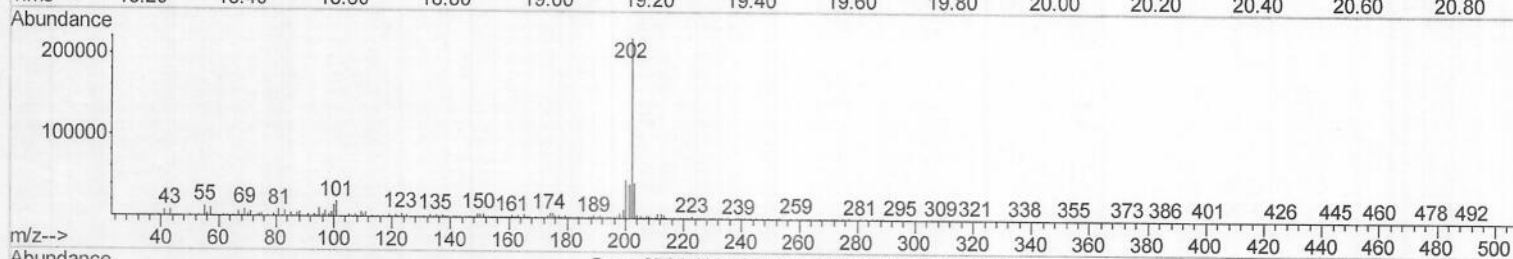
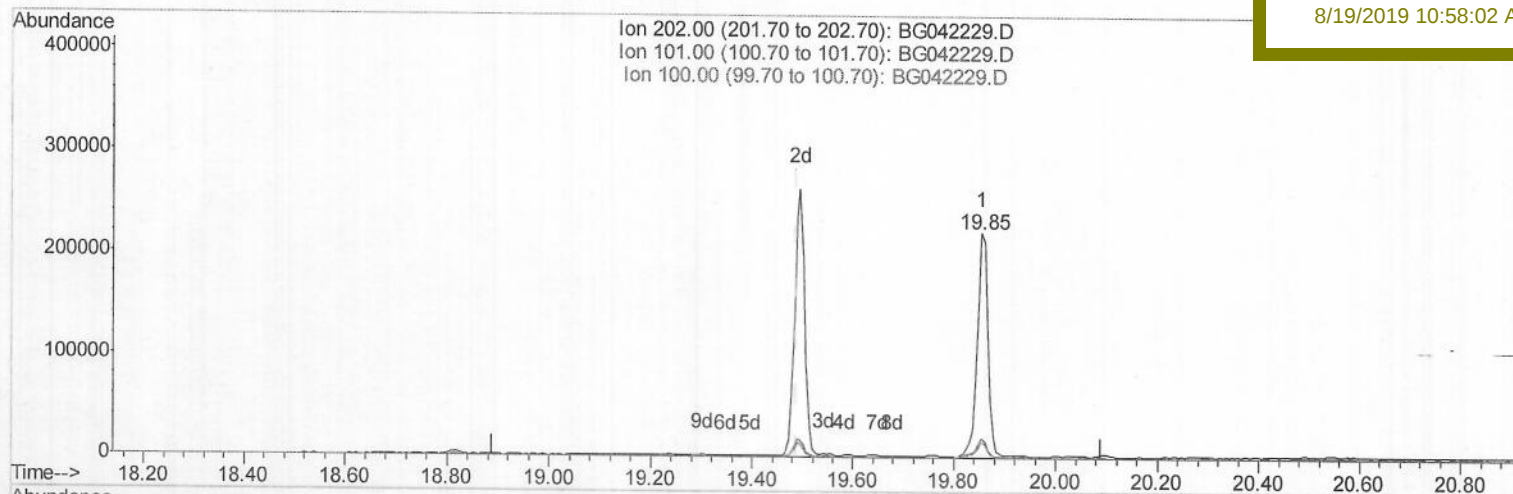
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(76) Fluoranthene (C)

19.854min (+0.365) 12.44ng/ul

response 322337

Ion	Exp%	Act%
202.00	100	100
101.00	8.40	8.66
100.00	6.60	6.97
0.00	0.00	0.00

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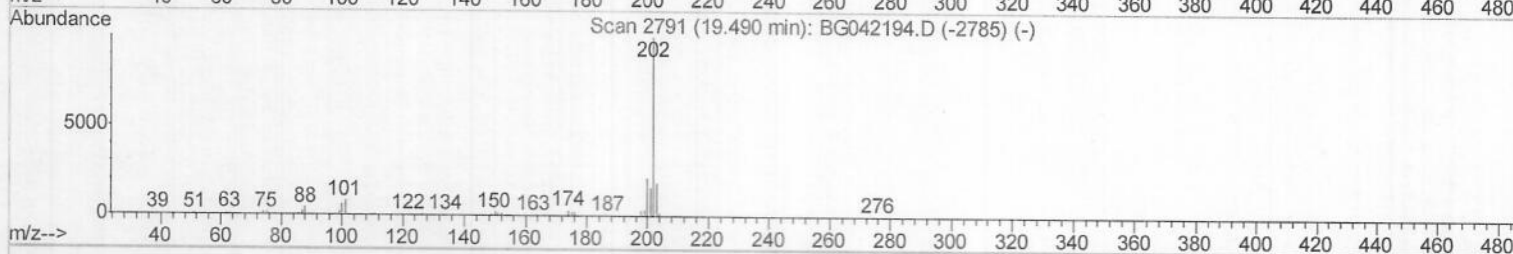
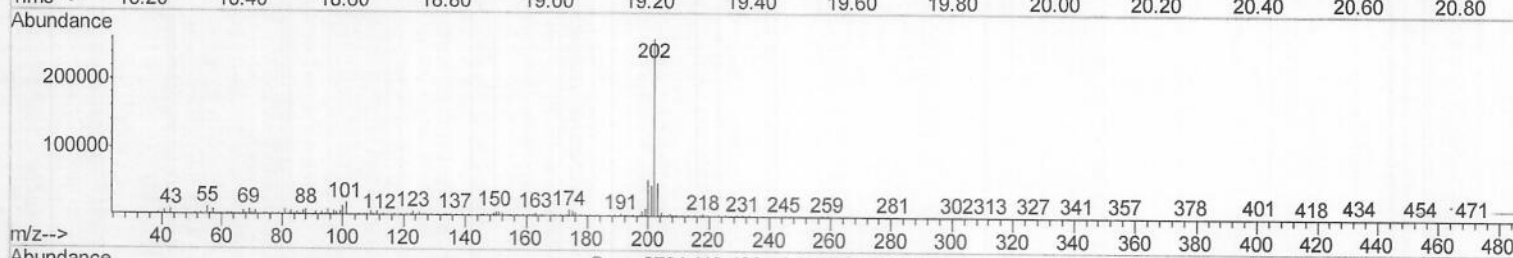
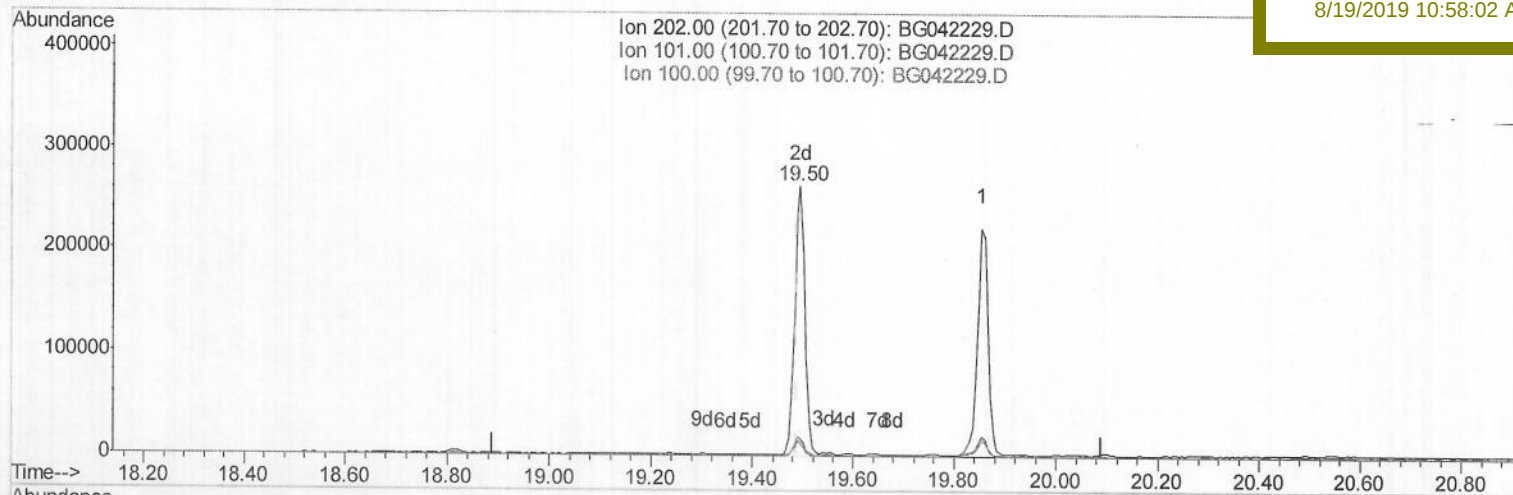
C0AG4

Manual Integrations

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TIC: BG042229.D

(76) Fluoranthene (C)

19.495min (+0.006) 14.33ng/ul m

response 371278

Ion Exp% Act%

202.00 100 100

101.00 8.40 6.59#

100.00 6.60 5.01#

0.00 0.00 0.00

JU
08/20/19

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) 1,4-Dichlorobenzene-d4	8.03	152	75836	20.00	ng/ul	0.00
18) Naphthalene-d8	10.85	136	302833	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.69	164	206363	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.44	188	409548	20.00	ng/ul	0.00
77) Chrysene-d12	21.73	240	413855	20.00	ng/ul	0.00
85) Perylene-d12	25.01	264	497865	20.00	ng/ul	0.02

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.41	96	3195	1.84	ng/uL	0.00
5) Phenol-d5	7.19	99	78425	13.18	ng/ul	0.00
7) Bis-(2-Chloroethyl) ether-d	7.34	67	40427	12.15	ng/ul	0.00
9) 2-Chlorophenol-d4	7.56	132	74712	14.68	ng/ul	0.00
13) 4-Methylphenol-d8	8.74	113	63993	12.81	ng/ul	0.00
19) Nitrobenzene-d5	9.20	128	35686	15.65	ng/ul	0.00
22) 2-Nitrophenol-d4	9.92	143	42658	16.12	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.47	165	81914	15.33	ng/ul	0.00
29) 4-Chloroaniline-d4	10.98	131	55992	9.44	ng/ul	0.00
43) Dimethylphthalate-d6	14.08	166	247665	15.47	ng/ul	0.00
46) Acenaphthylene-d8	14.38	160	291228	15.93	ng/ul	0.00
51) 4-Nitrophenol-d4	14.89	143	27968	10.41	ng/ul	0.01
57) Fluorene-d10	15.68	176	221714	15.57	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.81	200	9291	3.54	ng/ul	0.00
70) Anthracene-d10	17.54	188	322812	16.43	ng/ul	0.00
78) Pyrene-d10	19.82	212	362151	15.68	ng/ul	0.00
89) Benzo(a)pyrene-d12	24.78	264	420838	16.18	ng/ul	0.01

Target Compounds

					Ovalue	
6) Phenol	7.21	94	11347	1.847	ng/ul	96
14) Acetophenone	8.84	105	19420	2.527	ng/ul#	44
34) 2-Methylnaphthalene	12.51	142	20598	1.674	ng/ul	98
44) Dimethylphthalate	14.13	163	68304	4.298	ng/ul	98
69) Phenanthrene	17.48	178	188016	8.386	ng/ul	99
71) Anthracene	17.57	178	54807	2.410	ng/ul	98
76) Fluoranthene	19.50	202	371278m	14.332	ng/ul	
79) Pyrene	19.85	202	322337	11.579	ng/ul	97
80) Butylbenzylphthalate	20.74	149	95759	8.739	ng/ul	96
82) Benzo(a)anthracene	21.70	228	155063	5.404	ng/ul	98
83) Bis(2-ethylhexyl)phthalate	21.61	149	272567	18.231	ng/ul	99
84) Chrysene	21.77	228	154574	5.772	ng/ul	95
87) Benzo(b)fluoranthene	23.96	252	202384	6.478	ng/ul#	90
88) Benzo(k)fluoranthene	24.03	252	88926m	2.960	ng/ul	
90) Benzo(a)pyrene	24.85	252	144647	4.858	ng/ul#	90
91) Indeno(1,2,3-cd)pyrene	28.75	276	82710	2.382	ng/ul#	90
93) Benzo(a,h,i)perylene	29.92	276	75027	2.588	ng/ul	92

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

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