

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM061516\
 Data File : BM005800.D
 Acq On : 15 Jun 2016 00:31
 Operator : UM/SJ
 Sample : H3565-08
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 D9Y01

Manual Integrations
 APPROVED

sohil
 6/15/2016 6:35:54 PM

Quant Time: Jun 15 02:07:33 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM061516.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jun 15 01:42:36 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	90331	20.00	ng/ul	0.00
7) Naphthalene-d8	10.53	136	414457	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.39	164	217595	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.13	188	433696	20.00	ng/ul	0.00
29) Chrysene-d12	21.32	240	465481	20.00	ng/ul	0.00
34) Perylene-d12	23.57	264	502020	20.00	ng/ul	0.00

System Monitoring Compounds

2) 1,4-Dioxane-d8	3.19	96	957	0.49	ng/uL	0.00
3) Phenol-d5	6.93	99	20299	2.74	ng/ul	0.00
4) Bis-(2-Chloroethyl)ether-d	7.07	67	24184	5.48	ng/ul	0.00
5) 2-Chlorophenol-d4	7.28	132	27432	4.44	ng/ul	0.00
6) 4-Methylphenol-d8	8.47	113	7114	1.13	ng/ul	0.01
8) Nitrobenzene-d5	8.90	128	12321	4.20	ng/ul	0.00
9) 2-Nitrophenol-d4	9.63	143	13719	4.21	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.17	165	23918m	3.83	ng/ul	0.01
12) 4-Chloroaniline-d4	10.69	131	23096m	3.37	ng/ul	0.01
16) Dimethylphthalate-d6	13.79	166	133998	7.76	ng/ul	0.00
17) Acenaphthylene-d8	14.08	160	152417	6.95	ng/ul	0.00
20) 4-Nitrophenol-d4	14.64	143	18592m	9.36	ng/ul	0.02
21) Fluorene-d10	15.38	176	117346	8.16	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.53	200	12257	6.32	ng/ul	0.00
26) Anthracene-d10	17.23	188	233140	11.59	ng/ul	0.00
30) Pyrene-d10	19.52	212	381088	16.82	ng/ul	0.00
37) Benzo(a)pyrene-d12	23.43	264	434240	18.73	ng/ul	0.00

Target Compounds

						Qvalue
18) Acenaphthylene	14.11	152	53116	2.40	ng/ul	99
25) Phenanthrene	17.17	178	179314	7.54	ng/ul	100
27) Anthracene	17.26	178	55346	2.35	ng/ul	99
28) Fluoranthene	19.19	202	664583	26.25	ng/ul	100
31) Pyrene	19.56	202	561122	19.63	ng/ul	97
32) Benzo(a)anthracene	21.30	228	356161	13.23	ng/ul	97
33) Chrysene	21.35	228	344628	13.06	ng/ul	97
35) Benzo(b)fluoranthene	22.89	252	527638	17.59	ng/ul	98
36) Benzo(k)fluoranthene	22.93	252	176775m	6.37	ng/ul	
38) Benzo(a)pyrene	23.47	252	339431	11.85	ng/ul	99
39) Indeno(1,2,3-cd)pyrene	25.85	276	261565	7.96	ng/ul	94
40) Dibenzo(a,h)anthracene	25.84	278	68479	2.49	ng/ul#	76
41) Benzo(g,h,i)perylene	26.56	276	222373	7.82	ng/ul	99

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

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