

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM061516\
 Data File : BM005786.D
 Acq On : 14 Jun 2016 15:50
 Operator : UM/SJ
 Sample : SSTD01035
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD01035

Quant Time: Jun 15 01:19:11 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM061516.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jun 15 01:18:34 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	61563	20.00	ng/ul	0.00
7) Naphthalene-d8	10.53	136	311436	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.39	164	197646	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.13	188	449303	20.00	ng/ul	0.00
29) Chrysene-d12	21.31	240	431804	20.00	ng/ul	0.00
34) Perylene-d12	23.57	264	422102	20.00	ng/ul	0.00

System Monitoring Compounds

2) 1,4-Dioxane-d8	3.19	96	5523	3.92	ng/uL	0.00
3) Phenol-d5	6.92	99	50039	9.50	ng/ul	0.00
4) Bis-(2-Chloroethyl)ether-d	7.07	67	32450	10.26	ng/ul	0.00
5) 2-Chlorophenol-d4	7.27	132	43494	10.18	ng/ul	0.00
6) 4-Methylphenol-d8	8.46	113	43235	9.69	ng/ul	0.00
8) Nitrobenzene-d5	8.90	128	21623	9.65	ng/ul	0.00
9) 2-Nitrophenol-d4	9.62	143	23741	9.44	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.17	165	48905	10.59	ng/ul	0.00
12) 4-Chloroaniline-d4	10.67	131	62016	11.53	ng/ul	0.00
16) Dimethylphthalate-d6	13.79	166	163377	10.07	ng/ul	0.00
17) Acenaphthylene-d8	14.08	160	214238	10.76	ng/ul	0.00
20) 4-Nitrophenol-d4	14.63	143	15714	7.12	ng/ul	0.00
21) Fluorene-d10	15.38	176	149234	11.04	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.53	200	13919	6.23	ng/ul	0.00
26) Anthracene-d10	17.23	188	226988	10.80	ng/ul	0.00
30) Pyrene-d10	19.53	212	232340	11.22	ng/ul	0.00
37) Benzo(a)pyrene-d12	23.43	264	207441	10.78	ng/ul	0.00

Target Compounds

						Qvalue
11) Naphthalene	10.58	128	167014	10.74	ng/ul	100
13) 2-Methylnaphthalene	12.20	142	125000	10.80	ng/ul	97
14) 1-Methylnaphthalene	12.42	142	128762	11.26	ng/ul	99
18) Acenaphthylene	14.11	152	218626	10.77	ng/ul	99
19) Acenaphthene	14.45	153	143183	10.83	ng/ul	100
22) Fluorene	15.43	166	166383	11.25	ng/ul	98
25) Phenanthrene	17.17	178	267925	10.94	ng/ul	100
27) Anthracene	17.26	178	266412	10.77	ng/ul	99
28) Fluoranthene	19.19	202	291361	10.27	ng/ul	99
31) Pyrene	19.56	202	296820	11.34	ng/ul	98
32) Benzo(a)anthracene	21.30	228	264349	10.59	ng/ul	99
33) Chrysene	21.35	228	263452	10.78	ng/ul	100
35) Benzo(b)fluoranthene	22.89	252	270351	10.72	ng/ul	98
36) Benzo(k)fluoranthene	22.93	252	257740	10.41	ng/ul	99
38) Benzo(a)pyrene	23.47	252	258375	10.86	ng/ul	99
39) Indeno(1,2,3-cd)pyrene	25.86	276	285647	11.85	ng/ul	99
40) Dibenzo(a,h)anthracene	25.86	278	242405	12.00	ng/ul	99
41) Benzo(g,h,i)perylene	26.56	276	243754	11.94	ng/ul	98

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

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