

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM061216\
 Data File : BM005748.D
 Acq On : 12 Jun 2016 14:14
 Operator : UM/SJ
 Sample : H3536-13
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 D9Y23

Quant Time: Jun 13 03:30:41 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

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.75	152	80252	20.00	ng/ul	0.00
7) Naphthalene-d8	10.54	136	430847	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.39	164	272856	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.14	188	607038	20.00	ng/ul	0.00
29) Chrysene-d12	21.32	240	492992	20.00	ng/ul	0.00
34) Perylene-d12	23.58	264	429176	20.00	ng/ul	0.00

System Monitoring Compounds

2) 1,4-Dioxane-d8	3.19	96	3043	1.64	ng/uL	0.00
3) Phenol-d5	6.93	99	117983	17.04	ng/ul	0.00
4) Bis-(2-Chloroethyl)ether-d	7.07	67	83960	20.18	ng/ul	0.00
5) 2-Chlorophenol-d4	7.28	132	107282	19.16	ng/ul	0.00
6) 4-Methylphenol-d8	8.46	113	94505	16.15	ng/ul	0.00
8) Nitrobenzene-d5	8.91	128	59867	19.28	ng/ul	0.00
9) 2-Nitrophenol-d4	9.63	143	67327	19.18	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.17	165	116665	18.32	ng/ul	0.00
12) 4-Chloroaniline-d4	10.68	131	116287	15.13	ng/ul	0.00
16) Dimethylphthalate-d6	13.80	166	464458	20.64	ng/ul	0.00
17) Acenaphthylene-d8	14.09	160	583066	21.18	ng/ul	0.00
20) 4-Nitrophenol-d4	14.64	143	44846	13.62	ng/ul	0.00
21) Fluorene-d10	15.39	176	401257	21.06	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.53	200	25124	8.00	ng/ul	-0.01
26) Anthracene-d10	17.24	188	592916	20.73	ng/ul	0.00
30) Pyrene-d10	19.53	212	590365	24.92	ng/ul	0.00
37) Benzo(a)pyrene-d12	23.44	264	422517	21.59	ng/ul	0.00

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

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