

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM061716\
 Data File : BM005845.D
 Acq On : 17 Jun 2016 01:38
 Operator : UM/SJ
 Sample : H3537-09
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 D9Y39

Quant Time: Jun 17 02:14:00 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM061716.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jun 17 01:56:56 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	26686	20.00	ng/ul	0.00
7) Naphthalene-d8	10.52	136	127052	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.37	164	85834	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.12	188	207103	20.00	ng/ul	0.00
29) Chrysene-d12	21.30	240	255367	20.00	ng/ul	0.00
34) Perylene-d12	23.55	264	239281	20.00	ng/ul	0.00

System Monitoring Compounds

2) 1,4-Dioxane-d8	3.18	96	297	1.44	ng/uL	0.00
3) Phenol-d5	6.92	99	24536	10.98	ng/ul	0.00
4) Bis-(2-Chloroethyl)ether-d	7.06	67	14361	12.40	ng/ul	0.00
5) 2-Chlorophenol-d4	7.26	132	22685	11.93	ng/ul	0.00
6) 4-Methylphenol-d8	8.45	113	19200	9.61	ng/ul	0.00
8) Nitrobenzene-d5	8.89	128	11288	12.59	ng/ul	0.00
9) 2-Nitrophenol-d4	9.62	143	14370	13.11	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.16	165	25169	11.63	ng/ul	0.00
12) 4-Chloroaniline-d4	10.67	131	17306	11.82	ng/ul	0.00
16) Dimethylphthalate-d6	13.78	166	105103	12.94	ng/ul	0.00
17) Acenaphthylene-d8	14.07	160	124496	13.31	ng/ul	0.00
20) 4-Nitrophenol-d4	14.63	143	7178	7.33	ng/ul	0.02
21) Fluorene-d10	15.37	176	92045	13.30	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.52	200	11835	8.47	ng/ul	0.00
26) Anthracene-d10	17.22	188	146655	13.56	ng/ul	0.00
30) Pyrene-d10	19.52	212	174371	14.41	ng/ul	0.00
37) Benzo(a)pyrene-d12	23.41	264	167127	13.97	ng/ul	0.00

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

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