

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM060916\
 Data File : BM005711.D
 Acq On : 09 Jun 2016 13:09
 Operator : UM/SJ
 Sample : H3411-11
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 D9XQ9

Quant Time: Jun 10 01:25:00 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM060216.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jun 10 01:23:18 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.75	152	100262	20.00	ng/ul	0.00
7) Naphthalene-d8	10.54	136	511984	20.00	ng/ul	0.00
15) Acenaphthene-d10	14.40	164	319735	20.00	ng/ul	0.00
23) Phenanthrene-d10	17.14	188	703358	20.00	ng/ul	0.00
29) Chrysene-d12	21.33	240	649892	20.00	ng/ul	0.00
34) Perylene-d12	23.59	264	515583	20.00	ng/ul	0.00

System Monitoring Compounds

2) 1,4-Dioxane-d8	3.20	96	3479	1.46	ng/uL	0.00
3) Phenol-d5	6.94	99	127060	14.42	ng/ul	0.00
4) Bis-(2-Chloroethyl)ether-d	7.08	67	88367	16.47	ng/ul	0.00
5) 2-Chlorophenol-d4	7.29	132	110986	15.26	ng/ul	0.00
6) 4-Methylphenol-d8	8.47	113	92911	13.33	ng/ul	0.00
8) Nitrobenzene-d5	8.91	128	61522	15.65	ng/ul	0.00
9) 2-Nitrophenol-d4	9.63	143	68661	16.32	ng/ul	0.00
10) 2,4-Dichlorophenol-d3	10.18	165	118662	15.87	ng/ul	0.00
12) 4-Chloroaniline-d4	10.69	131	152486	17.16	ng/ul	0.00
16) Dimethylphthalate-d6	13.80	166	452522	18.41	ng/ul	0.00
17) Acenaphthylene-d8	14.09	160	562014	17.58	ng/ul	0.00
20) 4-Nitrophenol-d4	14.65	143	46792	13.29	ng/ul	0.00
21) Fluorene-d10	15.39	176	396476	18.78	ng/ul	0.00
24) 4,6-Dinitro-2-methylphenol	15.53	200	34922	10.54	ng/ul	0.00
26) Anthracene-d10	17.24	188	592359	17.73	ng/ul	0.00
30) Pyrene-d10	19.54	212	660824	20.23	ng/ul	0.00
37) Benzo(a)pyrene-d12	23.44	264	451900	19.05	ng/ul	0.00

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

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