

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM082916\
 Data File : BM007317.D
 Acq On : 29 Aug 2016 20:09
 Operator : UM/SJ
 Sample : H4575-09
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DA282

Quant Time: Aug 30 02:26:57 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM082316.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Aug 30 02:26:11 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.80	152	145288	20.00	ng/ul	0.00
18) Naphthalene-d8	10.58	136	700275	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.43	164	513229	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.17	188	1452209	20.00	ng/ul	0.00
75) Chrysene-d12	21.35	240	2238031	20.00	ng/ul	0.00
83) Perylene-d12	23.62	264	2527485	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.31	96	9474	3.33	ng/uL	0.00
5) Phenol-d5	6.96	99	214495	15.63	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.13	67	131151	18.19	ng/ul	0.00
9) 2-Chlorophenol-d4	7.33	132	170277	16.79	ng/ul	0.00
13) 4-Methylphenol-d8	8.49	113	193140	16.14	ng/ul	0.00
19) Nitrobenzene-d5	8.95	128	90934	17.69	ng/ul	0.00
22) 2-Nitrophenol-d4	9.67	143	104443	17.40	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.20	165	198201	16.58	ng/ul	0.00
29) 4-Chloroaniline-d4	10.72	131	190590	12.78	ng/ul	0.00
43) Dimethylphthalate-d6	13.83	166	876957	19.81	ng/ul	0.00
46) Acenaphthylene-d8	14.12	160	956670	18.64	ng/ul	0.00
51) 4-Nitrophenol-d4	14.62	143	146656	17.55	ng/ul	0.00
57) Fluorene-d10	15.42	176	765255	19.98	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.54	200	147981	16.21	ng/ul	0.00
70) Anthracene-d10	17.27	188	1371584	20.22	ng/ul	0.00
76) Pyrene-d10	19.56	212	1899874	20.30	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.47	264	2415035	20.75	ng/ul	0.00

Target Compounds						Qvalue
6) Phenol	6.99	94	15306	1.08	ng/ul#	71
44) Dimethylphthalate	13.88	163	673444	15.27	ng/ul	98

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

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