

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM011216\
 Data File : BM003948.D
 Acq On : 12 Jan 2016 10:54
 Operator : SJ/UM
 Sample : H1057-06
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 G9EL3

Quant Time: Jan 12 13:48:39 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM010116.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jan 12 00:35:37 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	41011	20.00	ng/ul	0.00
18) Naphthalene-d8	10.46	136	193018	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.32	164	115098	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.07	188	257193	20.00	ng/ul	0.00
78) Chrysene-d12	21.27	240	272273	20.00	ng/ul	-0.01
86) Perylene-d12	23.51	264	236959	20.00	ng/ul	-0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.13	96	4085	4.46	ng/uL	0.00
5) Phenol-d5	6.85	99	21966	6.45	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.00	67	62209	32.32	ng/ul	0.00
9) 2-Chlorophenol-d4	7.20	132	69478	24.83	ng/ul	0.00
13) 4-Methylphenol-d8	8.38	113	45005	16.31	ng/ul	0.00
19) Nitrobenzene-d5	8.83	128	41325	28.42	ng/ul	0.00
22) 2-Nitrophenol-d4	9.55	143	44940	28.20	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.08	165	76949	27.17	ng/ul	0.00
29) 4-Chloroaniline-d4	10.61	131	1507	0.40	ng/ul	0.01
44) Dimethylphthalate-d6	13.73	166	298961	33.09	ng/ul	0.00
47) Acenaphthylene-d8	14.02	160	360807	32.78	ng/ul	0.00
52) 4-Nitrophenol-d4	14.55	143	6210	3.83	ng/ul	0.00
58) Fluorene-d10	15.32	176	265541	34.65	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.45	200	36031	24.84	ng/ul	0.00
71) Anthracene-d10	17.17	188	413836	34.79	ng/ul	0.00
79) Pyrene-d10	19.47	212	457996	32.48	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.37	264	417524	35.28	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	3.16	88	6529	6.14	ng/uL#	96

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

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