

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG081215\
 Data File : BG018408.D
 Acq On : 12 Aug 2015 22:22
 Operator : UM/MA
 Sample : MDL-01-S-ML
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 MDL-01-S-ML

Quant Time: Aug 13 04:24:51 2015
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG080915.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Aug 13 04:07:48 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.02	152	89125	20.00	ng/ul	0.00
18) Naphthalene-d8	10.83	136	394911	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.65	164	260334	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.39	188	647374	20.00	ng/ul	0.00
78) Chrysene-d12	21.65	240	621657	20.00	ng/ul	0.00
86) Perylene-d12	24.86	264	614834	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.45	96	9957	4.94	ng/uL	0.00
5) Phenol-d5	7.18	99	193727	23.95	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.34	67	121356	24.21	ng/ul	0.00
9) 2-Chlorophenol-d4	7.55	132	149968	24.21	ng/ul	0.00
13) 4-Methylphenol-d8	8.72	113	166031	24.43	ng/ul	0.00
19) Nitrobenzene-d5	9.18	128	81558	26.49	ng/ul	0.00
22) 2-Nitrophenol-d4	9.90	143	91133	29.07	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.45	165	170945	25.69	ng/ul	0.00
29) 4-Chloroaniline-d4	10.96	131	242220	32.21	ng/ul	0.00
44) Dimethylphthalate-d6	14.05	166	589699	26.92	ng/ul	0.00
47) Acenaphthylene-d8	14.34	160	720685	26.13	ng/ul	0.00
52) 4-Nitrophenol-d4	14.84	143	105427	26.76	ng/ul	0.00
58) Fluorene-d10	15.64	176	519839	27.25	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.75	200	68047	21.37	ng/ul	0.00
71) Anthracene-d10	17.49	188	842046	27.20	ng/ul	0.00
79) Pyrene-d10	19.77	212	885905	27.47	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.64	264	900316	27.58	ng/ul	0.00

Target Compounds					Qvalue
35) 1-Methylnaphthalene	12.70	142	34744	2.35 ng/ul#	92

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

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