

Data Path : Z:\HPCHEM1\BNA G\DATA\BG100615\
 Data File : BG019058.D
 Acq On : 6 Oct 2015 23:07
 Operator : UM/NP
 Sample : G3915-03
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C03G5

Quant Time: Oct 07 01:09:08 2015
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG100415.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 07 01:07:34 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.04	152	17644	20.00	ng/ul	0.00
18) Naphthalene-d8	10.85	136	78788	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.67	164	53505	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.42	188	142230	20.00	ng/ul	0.00
78) Chrysene-d12	21.69	240	194869	20.00	ng/ul	0.00
86) Perylene-d12	24.92	264	193281	20.00	ng/ul	-0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.49	96	579	1.48	ng/uL	0.00
5) Phenol-d5	7.20	99	13201	8.68	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.37	67	32879	33.42	ng/ul	0.00
9) 2-Chlorophenol-d4	7.57	132	32445	28.43	ng/ul	0.00
13) 4-Methylphenol-d8	8.74	113	24316	20.34	ng/ul	0.00
19) Nitrobenzene-d5	9.20	128	20718	34.78	ng/ul	0.00
22) 2-Nitrophenol-d4	9.92	143	23285	38.00	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.47	165	41500	33.84	ng/ul	0.00
29) 4-Chloroaniline-d4	10.98	131	797	0.53	ng/ul	0.00
44) Dimethylphthalate-d6	14.07	166	161755	38.34	ng/ul	0.00
47) Acenaphthylene-d8	14.37	160	181365	34.76	ng/ul	0.00
52) 4-Nitrophenol-d4	14.86	143	8531	10.01	ng/ul	0.00
58) Fluorene-d10	15.66	176	138864	36.36	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.77	200	28085	36.40	ng/ul	0.00
71) Anthracene-d10	17.52	188	239240	35.72	ng/ul	0.00
79) Pyrene-d10	19.80	212	312152	34.58	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.70	264	352866	35.68	ng/ul	0.00

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

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