

Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA G\DATA\BG110517\
 Data File : BG030739.D
 Acq On : 5 Nov 2017 15:36
 Operator : SJ/JU
 Sample : I6107-06
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 B0E86

Quant Time: Nov 06 00:53:46 2017
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM-EPA-BG101417.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Nov 06 00:53:03 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.15	152	62407	20.00	ng/ul	0.00
18) Naphthalene-d8	10.97	136	271314	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.77	164	178383	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.51	188	476735	20.00	ng/ul	0.00
75) Chrysene-d12	21.77	240	650557	20.00	ng/ul	-0.02
83) Perylene-d12	25.08	264	642628	20.00	ng/ul	-0.03

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.53	96	6659	4.34	ng/uL	0.00
5) Phenol-d5	7.31	99	113154	18.89	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.47	67	68860	18.37	ng/ul	0.00
9) 2-Chlorophenol-d4	7.68	132	91956	21.76	ng/ul	0.00
13) 4-Methylphenol-d8	8.85	113	92111	19.04	ng/ul	0.00
19) Nitrobenzene-d5	9.32	128	43530	22.35	ng/ul	0.00
22) 2-Nitrophenol-d4	10.04	143	49797	24.95	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.59	165	93348	20.53	ng/ul	0.00
29) 4-Chloroaniline-d4	11.10	131	128684	24.36	ng/ul	0.00
43) Dimethylphthalate-d6	14.17	166	371457	24.36	ng/ul	0.00
46) Acenaphthylene-d8	14.47	160	420918	22.70	ng/ul	0.00
51) 4-Nitrophenol-d4	14.96	143	60842	21.61	ng/ul	0.00
57) Fluorene-d10	15.76	176	328732	26.32	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.87	200	59969	25.77	ng/ul	0.00
70) Anthracene-d10	17.60	188	579983	25.72	ng/ul	0.00
76) Pyrene-d10	19.88	212	784139	23.29	ng/ul	-0.01
87) Benzo(a)pyrene-d12	24.85	264	813527	26.73	ng/ul	-0.03

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
6) Phenol	7.33	94	7411	1.196	ng/ul#	87
44) Dimethylphthalate	14.21	163	188106	12.646	ng/ul	99

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

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