

Data Path : Z:\HPCHEM1\BNA G\DATA\BG101716\
 Data File : BG024379.D
 Acq On : 17 Oct 2016 12:35
 Operator : UM/SJ
 Sample : H5240-10
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BDPG8

Quant Time: Oct 17 13:32:24 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG0100716.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Oct 15 05:09:35 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.28	152	137350	20.00	ng/ul	0.00
18) Naphthalene-d8	11.11	136	667657	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.90	164	548902	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.64	188	1217949	20.00	ng/ul	0.00
75) Chrysene-d12	21.94	240	1389050	20.00	ng/ul	0.00
83) Perylene-d12	25.38	264	1389646	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.65	96	3391	1.17	ng/uL	0.02
5) Phenol-d5	7.41	99	75009	5.85	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.60	67	218074	25.17	ng/ul	0.00
9) 2-Chlorophenol-d4	7.80	132	177684	20.14	ng/ul	0.00
13) 4-Methylphenol-d8	8.97	113	156757	15.13	ng/ul	0.00
19) Nitrobenzene-d5	9.46	128	120213	25.37	ng/ul	0.00
22) 2-Nitrophenol-d4	10.18	143	139073	26.24	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.72	165	247674	21.45	ng/ul	0.00
29) 4-Chloroaniline-d4	11.25	131	4216	0.35	ng/ul	0.00
43) Dimethylphthalate-d6	14.29	166	1023989	26.72	ng/ul	0.00
46) Acenaphthylene-d8	14.60	160	1152069	26.24	ng/ul	0.00
51) 4-Nitrophenol-d4	15.06	143	39229	5.77	ng/ul	0.00
57) Fluorene-d10	15.89	176	973026	26.85	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.99	200	183958	25.13	ng/ul	0.00
70) Anthracene-d10	17.74	188	1488485	27.06	ng/ul	0.00
76) Pyrene-d10	20.01	212	1709244	26.36	ng/ul	0.00
87) Benzo(a)pyrene-d12	25.14	264	1748262	27.96	ng/ul	0.00

Target Compounds					Ovalue
33) 4-Chloro-3-methylphenol	12.36	107	15576	1.15	ng/ul 96

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

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