

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG120616\
 Data File : BG024977.D
 Acq On : 6 Dec 2016 13:10
 Operator : UM/SJ
 Sample : H5733-11
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 YA8Z3

Quant Time: Dec 07 02:38:26 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG113016.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Dec 07 02:35:14 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.24	152	80213	20.00	ng/ul	0.00
18) Naphthalene-d8	11.07	136	337895	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.86	164	288366	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.70	188	1066225	20.00	ng/ul	0.10
75) Chrysene-d12	21.89	240	1083931	20.00	ng/ul	0.00
83) Perylene-d12	25.31	264	1123796	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.60	96	2317	1.49	ng/uL	0.00
5) Phenol-d5	7.38	99	42767	6.18	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.56	67	127861	29.87	ng/ul	0.00
9) 2-Chlorophenol-d4	7.77	132	113539	23.46	ng/ul	0.00
13) 4-Methylphenol-d8	8.94	113	84013	14.53	ng/ul	0.00
19) Nitrobenzene-d5	9.42	128	74824	31.27	ng/ul	0.00
22) 2-Nitrophenol-d4	10.15	143	93126	31.22	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.68	165	146745	25.58	ng/ul	0.00
29) 4-Chloroaniline-d4	11.21	131	9334	1.66	ng/ul	0.00
43) Dimethylphthalate-d6	14.25	166	659515	33.25	ng/ul	0.00
46) Acenaphthylene-d8	14.56	160	696563	32.06	ng/ul	0.00
51) 4-Nitrophenol-d4	15.05	143	19250	5.65	ng/ul	0.00
57) Fluorene-d10	15.85	176	599463	32.11	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	16.08	200	183	0.03	ng/ul	0.12
70) Anthracene-d10	17.70	188	1063793	23.64	ng/ul	0.00
76) Pyrene-d10	19.98	212	1444577	32.36	ng/ul	0.00
87) Benzo(a)pyrene-d12	25.06	264	1534332	33.70	ng/ul	0.00

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

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