

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG111416\
 Data File : BG024771.D
 Acq On : 14 Nov 2016 20:07
 Operator : UM/SJ
 Sample : H5634-02
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BD4R1

Quant Time: Nov 15 01:02:18 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG111416.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 15 00:49:36 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.26	152	98473	20.00	ng/ul	0.00
18) Naphthalene-d8	11.09	136	401826	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.88	164	325901	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.62	188	805959	20.00	ng/ul	0.00
75) Chrysene-d12	21.92	240	1060276	20.00	ng/ul	0.00
83) Perylene-d12	25.34	264	1114212	20.00	ng/ul	-0.02

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.62	96	2625	1.38	ng/uL	0.00
5) Phenol-d5	7.40	99	51601	6.12	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.58	67	153622	28.95	ng/ul	0.00
9) 2-Chlorophenol-d4	7.79	132	138720	22.93	ng/ul	0.00
13) 4-Methylphenol-d8	8.95	113	102343	14.45	ng/ul	-0.01
19) Nitrobenzene-d5	9.43	128	88322	29.03	ng/ul	0.00
22) 2-Nitrophenol-d4	10.16	143	109938	30.27	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.70	165	191517	26.49	ng/ul	0.00
29) 4-Chloroaniline-d4	0.00	131	0d	0.00	ng/ul	
43) Dimethylphthalate-d6	14.27	166	760992	32.02	ng/ul	0.00
46) Acenaphthylene-d8	14.58	160	802982	31.01	ng/ul	0.00
51) 4-Nitrophenol-d4	15.06	143	28555	6.75	ng/ul	0.00
57) Fluorene-d10	15.87	176	688527	30.78	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.97	200	154207	27.83	ng/ul	-0.01
70) Anthracene-d10	17.72	188	1145655	32.18	ng/ul	0.00
76) Pyrene-d10	19.99	212	1495043	32.52	ng/ul	0.00
87) Benzo(a)pyrene-d12	25.10	264	1539807	31.57	ng/ul	-0.01

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

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