

Data Path : Z:\HPCHEM1\BNA G\DATA\BG101916\
 Data File : BG024441.D
 Acq On : 19 Oct 2016 14:07
 Operator : UM/SJ
 Sample : H5231-11
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BDNN9

Quant Time: Oct 19 17:29:24 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG0100716.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 19 10:58:30 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.28	152	180847	20.00	ng/ul	0.00
18) Naphthalene-d8	11.11	136	863088	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.90	164	726527	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.64	188	1439991	20.00	ng/ul	0.00
75) Chrysene-d12	21.93	240	1592850	20.00	ng/ul	0.00
83) Perylene-d12	25.38	264	1656456	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.64	96	11452	3.01	ng/uL	0.00
5) Phenol-d5	7.41	99	349952	20.74	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.60	67	237199	20.80	ng/ul	0.00
9) 2-Chlorophenol-d4	7.81	132	251015	21.60	ng/ul	0.00
13) 4-Methylphenol-d8	8.97	113	289594	21.23	ng/ul	0.00
19) Nitrobenzene-d5	9.46	128	139375	22.75	ng/ul	0.00
22) 2-Nitrophenol-d4	10.18	143	165004	24.08	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.71	165	316501	21.21	ng/ul	0.00
29) 4-Chloroaniline-d4	11.24	131	373655	23.83	ng/ul	0.00
43) Dimethylphthalate-d6	14.29	166	1106079	21.80	ng/ul	0.00
46) Acenaphthylene-d8	14.59	160	1280255	22.03	ng/ul	0.00
51) 4-Nitrophenol-d4	15.06	143	181197	20.14	ng/ul	0.00
57) Fluorene-d10	15.88	176	1035857	21.60	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.99	200	191239	22.10	ng/ul	0.00
70) Anthracene-d10	17.74	188	1525148	23.45	ng/ul	0.00
76) Pyrene-d10	20.01	212	1791162	24.09	ng/ul	0.00
87) Benzo(a)pyrene-d12	25.13	264	1847298	24.79	ng/ul	0.00

Target Compounds					Ovalue
6) Phenol	7.44	94	18816	1.10 ng/ul	89
44) Dimethylphthalate	14.34	163	491851	9.89 ng/ul	98

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

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