

Data Path : Z:\HPCHEM1\BNA G\DATA\BG101116\
 Data File : BG024282.D
 Acq On : 12 Oct 2016 2:01
 Operator : UM/SJ
 Sample : H5113-07
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BDP64

Quant Time: Oct 12 04:52:32 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG0100716.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 12 04:50:43 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.29	152	146630	20.00	ng/ul	0.00
18) Naphthalene-d8	11.12	136	634211	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.91	164	508532	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.64	188	1110097	20.00	ng/ul	0.00
75) Chrysene-d12	21.94	240	1261985	20.00	ng/ul	0.00
83) Perylene-d12	25.39	264	1282662	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.65	96	12008	3.89	ng/uL	0.00
5) Phenol-d5	7.42	99	307854	22.50	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.60	67	216712	23.43	ng/ul	0.00
9) 2-Chlorophenol-d4	7.81	132	224980	23.88	ng/ul	0.00
13) 4-Methylphenol-d8	8.97	113	252466	22.83	ng/ul	0.00
19) Nitrobenzene-d5	9.46	128	115441	25.65	ng/ul	0.00
22) 2-Nitrophenol-d4	10.19	143	130206	25.86	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.72	165	246856	22.51	ng/ul	0.00
29) 4-Chloroaniline-d4	11.24	131	186385	16.18	ng/ul	0.00
43) Dimethylphthalate-d6	14.30	166	941218	26.51	ng/ul	0.00
46) Acenaphthylene-d8	14.61	160	1076647	26.47	ng/ul	0.00
51) 4-Nitrophenol-d4	15.06	143	170603	27.10	ng/ul	0.00
57) Fluorene-d10	15.89	176	891632	26.56	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.99	200	173348	25.98	ng/ul	0.00
70) Anthracene-d10	17.74	188	1417843	28.28	ng/ul	0.00
76) Pyrene-d10	20.01	212	1677814	28.48	ng/ul	0.00
87) Benzo(a)pyrene-d12	25.15	264	1655090	28.68	ng/ul	0.00

Target Compounds

					Ovalue	
6) Phenol	7.45	94	15983	1.15	ng/ul#	69
44) Dimethylphthalate	14.35	163	763485	21.93	ng/ul	98

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

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