

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG091515\
 Data File : BG018687.D
 Acq On : 15 Sep 2015 17:47
 Operator : UM/IZ
 Sample : G3641-06
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 B0AM4

Quant Time: Sep 15 18:21:36 2015
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG091015.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 15 13:56:56 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.01	152	68818	20.00	ng/ul	0.00
18) Naphthalene-d8	10.81	136	305475	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.63	164	211006	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.37	188	537435	20.00	ng/ul	0.00
78) Chrysene-d12	21.62	240	554741	20.00	ng/ul	0.00
86) Perylene-d12	24.81	264	561108	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.44	96	1857	1.30	ng/uL	0.00
5) Phenol-d5	7.16	99	46414	8.21	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.32	67	108170	33.03	ng/ul	0.00
9) 2-Chlorophenol-d4	7.54	132	122087	28.53	ng/ul	0.00
13) 4-Methylphenol-d8	8.70	113	86079	18.77	ng/ul	0.00
19) Nitrobenzene-d5	9.16	128	76368	33.00	ng/ul	0.00
22) 2-Nitrophenol-d4	9.89	143	81345	34.95	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.42	165	162815	33.23	ng/ul	0.00
29) 4-Chloroaniline-d4	10.94	131	8229	1.44	ng/ul	0.00
44) Dimethylphthalate-d6	14.03	166	549994	32.39	ng/ul	0.00
47) Acenaphthylene-d8	14.32	160	679378	33.76	ng/ul	0.00
52) 4-Nitrophenol-d4	14.82	143	20951	6.43	ng/ul	0.00
58) Fluorene-d10	15.62	176	508557	34.25	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.73	200	75822	29.80	ng/ul	0.00
71) Anthracene-d10	17.46	188	816737	34.52	ng/ul	0.00
79) Pyrene-d10	19.75	212	880228	34.49	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.59	264	952123	34.87	ng/ul	-0.02

Target Compounds						Qvalue
2) 1,4-Dioxane	3.48	88	22635	14.62	ng/uL#	77
8) Bis(2-Chloroethyl)ether	7.42	93	890605	204.00	ng/ul	96

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

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