

Data Path : Z:\HPCHEM1\BNA G\DATA\BG101915\
 Data File : BG019251.D
 Acq On : 19 Oct 2015 22:20
 Operator : UM/NP
 Sample : G4075-17
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 JG9M9

Quant Time: Oct 20 05:51:36 2015
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG101515.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Oct 20 05:49:53 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.01	152	11461	20.00	ng/ul	0.00
18) Naphthalene-d8	10.81	136	55105	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.63	164	43010	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.38	188	109838	20.00	ng/ul	0.00
78) Chrysene-d12	21.64	240	131686	20.00	ng/ul	0.00
86) Perylene-d12	24.85	264	122133	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.44	96	208	1.01	ng/uL	0.00
5) Phenol-d5	7.18	99	7677	7.68	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.33	67	20444	31.76	ng/ul	0.00
9) 2-Chlorophenol-d4	7.54	132	20343	27.91	ng/ul	0.00
13) 4-Methylphenol-d8	8.72	113	14406	17.00	ng/ul	0.00
19) Nitrobenzene-d5	9.16	128	13852	32.63	ng/ul	0.00
22) 2-Nitrophenol-d4	9.89	143	15828	35.13	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.43	165	28747	32.21	ng/ul	0.00
29) 4-Chloroaniline-d4	10.95	131	3008	2.62	ng/ul	0.00
44) Dimethylphthalate-d6	14.04	166	130320	34.64	ng/ul	0.00
47) Acenaphthylene-d8	14.33	160	136706	33.62	ng/ul	0.00
52) 4-Nitrophenol-d4	14.85	143	4996	7.39	ng/ul	0.00
58) Fluorene-d10	15.63	176	108687	33.80	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.75	200	23145	34.16	ng/ul	0.00
71) Anthracene-d10	17.48	188	182422	35.15	ng/ul	0.00
79) Pyrene-d10	19.76	212	205591	34.42	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.64	264	230194	37.31	ng/ul	0.00

Target Compounds Ovalue

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

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