

Data Path : Z:\HPCHEM1\BNA G\DATA\BG101915\
 Data File : BG019247.D
 Acq On : 19 Oct 2015 19:49
 Operator : UM/NP
 Sample : G4075-16
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 JG9M8

Quant Time: Oct 20 05:51:12 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	8.00	152	12915	20.00	ng/ul	0.00
18) Naphthalene-d8	10.81	136	60502	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.64	164	43688	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.38	188	116764	20.00	ng/ul	0.00
78) Chrysene-d12	21.65	240	135389	20.00	ng/ul	0.00
86) Perylene-d12	24.86	264	126895	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.43	96	68	0.29	ng/uL	0.00
5) Phenol-d5	7.18	99	6944	6.16	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.33	67	23910	32.97	ng/ul	0.00
9) 2-Chlorophenol-d4	7.54	132	21766	26.50	ng/ul	0.00
13) 4-Methylphenol-d8	8.72	113	15168	15.88	ng/ul	0.00
19) Nitrobenzene-d5	9.17	128	16586	35.58	ng/ul	0.00
22) 2-Nitrophenol-d4	9.89	143	19130	38.67	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.44	165	31624	32.27	ng/ul	0.00
29) 4-Chloroaniline-d4	0.00	131	0	0.00	ng/ul	
44) Dimethylphthalate-d6	14.04	166	141798	37.11	ng/ul	0.00
47) Acenaphthylene-d8	14.33	160	152855	37.01	ng/ul	0.00
52) 4-Nitrophenol-d4	14.86	143	4590	6.68	ng/ul	0.00
58) Fluorene-d10	15.63	176	121272	37.13	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.75	200	25717	35.70	ng/ul	0.00
71) Anthracene-d10	17.48	188	202015	36.62	ng/ul	0.00
79) Pyrene-d10	19.77	212	223145	36.34	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.64	264	242895	37.89	ng/ul	0.00

Target Compounds Ovalue

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

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