

Data Path : Z:\HPCHEM1\BNA G\DATA\BG101515\
 Data File : BG019213.D
 Acq On : 15 Oct 2015 19:51
 Operator : UM/NP
 Sample : PB86122BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK22

Quant Time: Oct 16 04:08:53 2015
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG101515.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Oct 16 04:07:16 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.01	152	13530	20.00	ng/ul	0.00
18) Naphthalene-d8	10.81	136	61335	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.64	164	44348	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.38	188	123250	20.00	ng/ul	0.00
78) Chrysene-d12	21.65	240	173308	20.00	ng/ul	0.00
86) Perylene-d12	24.85	264	161523	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.44	96	1580	6.50	ng/uL	0.00
5) Phenol-d5	7.17	99	35347	29.94	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.33	67	24775	32.61	ng/ul	0.00
9) 2-Chlorophenol-d4	7.54	132	27182	31.59	ng/ul	0.00
13) 4-Methylphenol-d8	8.71	113	30059	30.05	ng/ul	0.00
19) Nitrobenzene-d5	9.17	128	15999	33.86	ng/ul	0.00
22) 2-Nitrophenol-d4	9.89	143	17422	34.74	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.43	165	29534	29.73	ng/ul	0.00
29) 4-Chloroaniline-d4	10.95	131	46319	36.21	ng/ul	0.00
44) Dimethylphthalate-d6	14.04	166	144835	37.34	ng/ul	0.00
47) Acenaphthylene-d8	14.33	160	141796	33.82	ng/ul	0.00
52) 4-Nitrophenol-d4	14.84	143	23945	34.34	ng/ul	0.00
58) Fluorene-d10	15.63	176	112849	34.04	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.74	200	23263	30.59	ng/ul	0.00
71) Anthracene-d10	17.48	188	211014	36.24	ng/ul	0.00
79) Pyrene-d10	19.77	212	274996	34.98	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.64	264	304797	37.35	ng/ul	0.00

Target Compounds Ovalue

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

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