

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG032816\
 Data File : BG021539.D
 Acq On : 28 Mar 2016 23:33
 Operator : UM/SJ
 Sample : H1910-13
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0AC3

Quant Time: Mar 29 02:05:12 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG031616.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Mar 29 02:01:55 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.06	152	7516	20.00	ng/ul	0.00
18) Naphthalene-d8	10.86	136	37726	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.67	164	26465	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.41	188	72111	20.00	ng/ul	0.00
78) Chrysene-d12	21.67	240	86807	20.00	ng/ul	0.00
86) Perylene-d12	24.88	264	76163	20.00	ng/ul	-0.02

System Monitoring Compounds

3) 1,4-Dioxane-d8	0.00	96	0	0.00	ng/uL	
5) Phenol-d5	7.34	99	3000	4.22	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.38	67	13481	32.03	ng/ul	0.00
9) 2-Chlorophenol-d4	7.63	132	13272	24.59	ng/ul	0.00
13) 4-Methylphenol-d8	8.84	113	8524	14.01	ng/ul	0.00
19) Nitrobenzene-d5	9.22	128	9222	32.58	ng/ul	0.00
22) 2-Nitrophenol-d4	9.94	143	10170	31.43	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.56	165	15928	27.54	ng/ul	0.00
29) 4-Chloroaniline-d4	11.02	131	365	0.48	ng/ul	0.02
44) Dimethylphthalate-d6	14.07	166	72626	32.96	ng/ul	0.00
47) Acenaphthylene-d8	14.37	160	85559	31.50	ng/ul	0.00
52) 4-Nitrophenol-d4	15.11	143	1365	3.88	ng/ul	0.02
58) Fluorene-d10	15.66	176	65530	32.98	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.79	200	12794	26.86	ng/ul	0.00
71) Anthracene-d10	17.51	188	113280	32.06	ng/ul	0.00
79) Pyrene-d10	19.79	212	141137	34.87	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.66	264	124797	35.50	ng/ul	-0.01

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

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