

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG032216\
 Data File : BG021478.D
 Acq On : 22 Mar 2016 17:37
 Operator : UM/SJ
 Sample : H1835-25
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 A4T36

Quant Time: Mar 23 03:33:50 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG031616.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Mar 23 03:32:13 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.07	152	9499	20.00	ng/ul	0.00
18) Naphthalene-d8	10.88	136	45130	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.68	164	28960	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.42	188	66465	20.00	ng/ul	0.00
78) Chrysene-d12	21.68	240	79745	20.00	ng/ul	0.00
86) Perylene-d12	24.89	264	75550	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	0.00	96	0	0.00	ng/uL	
5) Phenol-d5	7.35	99	5834	6.50	ng/ul	0.02
7) Bis-(2-Chloroethyl)ether-d	7.40	67	15510	29.15	ng/ul	0.00
9) 2-Chlorophenol-d4	7.64	132	16508	24.21	ng/ul	0.00
13) 4-Methylphenol-d8	8.85	113	12675	16.48	ng/ul	0.00
19) Nitrobenzene-d5	9.24	128	10760	31.78	ng/ul	0.00
22) 2-Nitrophenol-d4	9.96	143	12111	31.29	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.57	165	22461	32.47	ng/ul	0.00
29) 4-Chloroaniline-d4	0.00	131	0	0.00	ng/ul	
44) Dimethylphthalate-d6	14.08	166	83162	34.49	ng/ul	0.00
47) Acenaphthylene-d8	14.38	160	100358	33.76	ng/ul	0.00
52) 4-Nitrophenol-d4	15.09	143	1796	4.66	ng/ul	0.02
58) Fluorene-d10	15.67	176	72252	33.23	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.79	200	13134	29.91	ng/ul	0.00
71) Anthracene-d10	17.52	188	115809	35.56	ng/ul	0.00
79) Pyrene-d10	19.79	212	139625	37.55	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.67	264	141032	40.45	ng/ul	0.00

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

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