

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG032816\
 Data File : BG021537.D
 Acq On : 28 Mar 2016 22:14
 Operator : UM/SJ
 Sample : H1910-08
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0AE4

Quant Time: Mar 29 02:05:01 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	8.05	152	9607	20.00	ng/ul	0.00
18) Naphthalene-d8	10.86	136	47533	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.67	164	32546	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.41	188	89180	20.00	ng/ul	0.00
78) Chrysene-d12	21.67	240	106336	20.00	ng/ul	-0.01
86) Perylene-d12	24.89	264	95455	20.00	ng/ul	-0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	0.00	96	0	0.00	ng/uL	
5) Phenol-d5	7.35	99	4403	4.85	ng/ul	0.01
7) Bis-(2-Chloroethyl)ether-d	7.38	67	12627	23.47	ng/ul	0.00
9) 2-Chlorophenol-d4	7.64	132	13490	19.56	ng/ul	0.00
13) 4-Methylphenol-d8	8.85	113	9052	11.64	ng/ul	0.00
19) Nitrobenzene-d5	9.22	128	9282	26.03	ng/ul	0.00
22) 2-Nitrophenol-d4	9.95	143	9664	23.71	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.57	165	15510	21.28	ng/ul	0.00
29) 4-Chloroaniline-d4	11.01	131	1074	1.13	ng/ul	0.00
44) Dimethylphthalate-d6	14.07	166	70730	26.10	ng/ul	0.00
47) Acenaphthylene-d8	14.37	160	82916	24.82	ng/ul	0.00
52) 4-Nitrophenol-d4	15.11	143	1689	3.90	ng/ul	0.02
58) Fluorene-d10	15.66	176	62140	25.43	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.78	200	12288	20.86	ng/ul	0.00
71) Anthracene-d10	17.51	188	106993	24.48	ng/ul	0.00
79) Pyrene-d10	19.79	212	134068	27.04	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.67	264	121404	27.56	ng/ul	0.00

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

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