

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG032816\
 Data File : BG021536.D
 Acq On : 28 Mar 2016 21:34
 Operator : UM/SJ
 Sample : H1910-07
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0AE2

Quant Time: Mar 29 02:04:56 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	7928	20.00	ng/ul	0.00
18) Naphthalene-d8	10.86	136	38241	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.67	164	26351	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.41	188	73852	20.00	ng/ul	0.00
78) Chrysene-d12	21.67	240	89636	20.00	ng/ul	0.00
86) Perylene-d12	24.89	264	80051	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	3444	4.59	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.38	67	12003	27.03	ng/ul	0.00
9) 2-Chlorophenol-d4	7.63	132	12538	22.03	ng/ul	0.00
13) 4-Methylphenol-d8	8.84	113	8522	13.27	ng/ul	0.00
19) Nitrobenzene-d5	9.23	128	8467	29.51	ng/ul	0.00
22) 2-Nitrophenol-d4	9.95	143	9268	28.26	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.57	165	15912	27.14	ng/ul	0.00
29) 4-Chloroaniline-d4	11.02	131	1068	1.39	ng/ul	0.01
44) Dimethylphthalate-d6	14.07	166	72651	33.11	ng/ul	0.00
47) Acenaphthylene-d8	14.37	160	81840	30.26	ng/ul	0.00
52) 4-Nitrophenol-d4	15.10	143	1491	4.25	ng/ul	0.01
58) Fluorene-d10	15.66	176	63740	32.22	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.78	200	13776	28.24	ng/ul	0.00
71) Anthracene-d10	17.51	188	111637	30.85	ng/ul	0.00
79) Pyrene-d10	19.78	212	142042	33.98	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.66	264	130245	35.25	ng/ul	-0.01

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

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