

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG032816\
 Data File : BG021545.D
 Acq On : 29 Mar 2016 3:31
 Operator : UM/SJ
 Sample : H1910-19
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0AG6

Quant Time: Mar 29 06:10:29 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	8038	20.00	ng/ul	0.00
18) Naphthalene-d8	10.86	136	37629	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.67	164	26656	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.41	188	73168	20.00	ng/ul	0.00
78) Chrysene-d12	21.66	240	87369	20.00	ng/ul	-0.01
86) Perylene-d12	24.88	264	76647	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	3789	4.99	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.38	67	13175	29.27	ng/ul	0.00
9) 2-Chlorophenol-d4	7.63	132	14011	24.28	ng/ul	0.00
13) 4-Methylphenol-d8	8.84	113	8607	13.22	ng/ul	0.00
19) Nitrobenzene-d5	9.23	128	9297	32.93	ng/ul	0.00
22) 2-Nitrophenol-d4	9.95	143	10260	31.79	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.56	165	16213	28.11	ng/ul	0.00
29) 4-Chloroaniline-d4	11.02	131	186	0.25	ng/ul	0.02
44) Dimethylphthalate-d6	14.07	166	76770	34.59	ng/ul	0.00
47) Acenaphthylene-d8	14.37	160	88560	32.37	ng/ul	0.00
52) 4-Nitrophenol-d4	15.11	143	1427	4.02	ng/ul	0.02
58) Fluorene-d10	15.66	176	67001	33.48	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.78	200	13729	28.40	ng/ul	0.00
71) Anthracene-d10	17.51	188	118549	33.06	ng/ul	0.00
79) Pyrene-d10	19.78	212	147147	36.12	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.66	264	130542	36.90	ng/ul	-0.01

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

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