

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG032216\
 Data File : BG021468.D
 Acq On : 22 Mar 2016 10:31
 Operator : UM/SJ
 Sample : PB89055BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK55

Quant Time: Mar 23 01:32:22 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG031616.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Mar 23 01:30:27 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.07	152	9888	20.00	ng/ul	0.00
18) Naphthalene-d8	10.87	136	48670	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.68	164	31375	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.42	188	74953	20.00	ng/ul	0.00
78) Chrysene-d12	21.67	240	87315	20.00	ng/ul	0.00
86) Perylene-d12	24.89	264	83445	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.47	96	1086	4.93	ng/uL	0.00
5) Phenol-d5	7.34	99	26020	27.83	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.40	67	17275	31.19	ng/ul	0.00
9) 2-Chlorophenol-d4	7.64	132	20645	29.08	ng/ul	0.00
13) 4-Methylphenol-d8	8.85	113	23964	29.93	ng/ul	0.00
19) Nitrobenzene-d5	9.24	128	11364	31.12	ng/ul	0.00
22) 2-Nitrophenol-d4	9.96	143	12856	30.80	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.57	165	22787	30.54	ng/ul	0.00
29) 4-Chloroaniline-d4	11.02	131	32988	33.82	ng/ul	0.00
44) Dimethylphthalate-d6	14.09	166	87045	33.32	ng/ul	0.00
47) Acenaphthylene-d8	14.38	160	101363	31.48	ng/ul	0.00
52) 4-Nitrophenol-d4	15.08	143	10564	25.30	ng/ul	0.01
58) Fluorene-d10	15.67	176	73363	31.15	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.79	200	13066	26.39	ng/ul	0.00
71) Anthracene-d10	17.52	188	119232	32.46	ng/ul	0.00
79) Pyrene-d10	19.79	212	140485	34.50	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.67	264	144896	37.62	ng/ul	0.00

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

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