

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG033016\
 Data File : BG021584.D
 Acq On : 30 Mar 2016 16:48
 Operator : UM/SJ
 Sample : PB89255BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK55

Quant Time: Mar 31 10:56:27 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG031616.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Mar 31 02:07:33 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.05	152	7678	20.00	ng/ul	0.00
18) Naphthalene-d8	10.86	136	34502	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.67	164	22209	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.41	188	58574	20.00	ng/ul	0.00
78) Chrysene-d12	21.67	240	66589	20.00	ng/ul	0.00
86) Perylene-d12	24.88	264	66202	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.45	96	1246	7.28	ng/uL	0.00
5) Phenol-d5	7.34	99	24043	33.12	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.37	67	15077	35.06	ng/ul	0.00
9) 2-Chlorophenol-d4	7.63	132	20211	36.66	ng/ul	0.00
13) 4-Methylphenol-d8	8.85	113	20868	33.56	ng/ul	0.00
19) Nitrobenzene-d5	9.22	128	10472	40.46	ng/ul	0.00
22) 2-Nitrophenol-d4	9.94	143	11426	38.62	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.56	165	20438	38.64	ng/ul	0.00
29) 4-Chloroaniline-d4	11.01	131	9015	13.04	ng/ul	0.01
44) Dimethylphthalate-d6	14.07	166	76330	41.28	ng/ul	0.00
47) Acenaphthylene-d8	14.36	160	91000	39.92	ng/ul	0.00
52) 4-Nitrophenol-d4	15.10	143	9042	30.59	ng/ul	0.00
58) Fluorene-d10	15.66	176	68621	41.16	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.79	200	13846	35.78	ng/ul	0.00
71) Anthracene-d10	17.51	188	115264	40.16	ng/ul	0.00
79) Pyrene-d10	19.79	212	134132	43.20	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.66	264	136709	44.74	ng/ul	0.00

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

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