

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG032216\
 Data File : BG021469.D
 Acq On : 22 Mar 2016 11:13
 Operator : UM/SJ
 Sample : PB89055TB
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SLEB16

Quant Time: Mar 23 01:32:26 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	10849	20.00	ng/ul	0.00
18) Naphthalene-d8	10.88	136	54259	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.69	164	33878	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.42	188	71843	20.00	ng/ul	0.00
78) Chrysene-d12	21.67	240	83382	20.00	ng/ul	0.00
86) Perylene-d12	24.89	264	80189	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.47	96	1512	6.25	ng/uL	0.00
5) Phenol-d5	7.34	99	31134	30.35	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.39	67	21071	34.68	ng/ul	0.00
9) 2-Chlorophenol-d4	7.64	132	25599	32.86	ng/ul	0.00
13) 4-Methylphenol-d8	8.85	113	28318	32.23	ng/ul	0.00
19) Nitrobenzene-d5	9.24	128	13803	33.91	ng/ul	0.00
22) 2-Nitrophenol-d4	9.96	143	15565	33.45	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.57	165	27361	32.89	ng/ul	0.00
29) 4-Chloroaniline-d4	11.02	131	39341	36.18	ng/ul	0.00
44) Dimethylphthalate-d6	14.09	166	96155	34.09	ng/ul	0.00
47) Acenaphthylene-d8	14.38	160	119499	34.37	ng/ul	0.00
52) 4-Nitrophenol-d4	15.07	143	10227	22.68	ng/ul	0.00
58) Fluorene-d10	15.67	176	83219	32.72	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.80	200	13010	27.41	ng/ul	0.00
71) Anthracene-d10	17.52	188	120524	34.23	ng/ul	0.00
79) Pyrene-d10	19.80	212	141499	36.39	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.67	264	147453	39.84	ng/ul	0.00

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

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