

Data Path : Z:\HPCHEM1\BNA G\DATA\BG020916\
 Data File : BG020813.D
 Acq On : 9 Feb 2016 12:13
 Operator : SJ/UM
 Sample : PB88249BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK49

Quant Time: Feb 10 00:17:22 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG020816.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Feb 10 00:13:30 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.27	152	16013	20.00	ng/ul	0.00
18) Naphthalene-d8	11.09	136	80769	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.88	164	51625	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.62	188	121858	20.00	ng/ul	0.00
78) Chrysene-d12	21.90	240	112894	20.00	ng/ul	0.00
86) Perylene-d12	25.29	264	104339	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.60	96	2227	6.75	ng/uL	0.00
5) Phenol-d5	7.40	99	51960	32.04	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.58	67	30146	34.83	ng/ul	0.00
9) 2-Chlorophenol-d4	7.80	132	43058	33.92	ng/ul	0.00
13) 4-Methylphenol-d8	8.96	113	45651	32.37	ng/ul	0.00
19) Nitrobenzene-d5	9.44	128	23494	36.25	ng/ul	0.00
22) 2-Nitrophenol-d4	10.17	143	23695	36.76	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.71	165	40064	32.29	ng/ul	0.00
29) 4-Chloroaniline-d4	11.22	131	71177	43.08	ng/ul	0.00
44) Dimethylphthalate-d6	14.28	166	169464	38.15	ng/ul	0.00
47) Acenaphthylene-d8	14.58	160	194044	35.76	ng/ul	0.00
52) 4-Nitrophenol-d4	15.05	143	32741	36.81	ng/ul	0.00
58) Fluorene-d10	15.87	176	139160	36.81	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.97	200	15683	26.56	ng/ul	0.00
71) Anthracene-d10	17.72	188	224841	37.34	ng/ul	0.00
79) Pyrene-d10	19.99	212	219485	37.82	ng/ul	0.00
90) Benzo(a)pyrene-d12	25.06	264	213593	36.94	ng/ul	0.00

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

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