

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG080816\
 Data File : BG023539.D
 Acq On : 8 Aug 2016 18:13
 Operator : UM/SJ
 Sample : PB92714BL
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK14

Quant Time: Aug 09 04:43:58 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG080416.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Aug 09 02:15:40 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.12	152	109581	20.00	ng/ul	0.00
18) Naphthalene-d8	10.95	136	548095	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.78	164	399522	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.54	188	1005389	20.00	ng/ul	0.00
75) Chrysene-d12	21.86	240	996776	20.00	ng/ul	0.00
83) Perylene-d12	25.23	264	975316	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.49	96	16809	6.52	ng/uL	0.00
5) Phenol-d5	7.27	99	411478	37.28	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.43	67	293559	41.50	ng/ul	0.00
9) 2-Chlorophenol-d4	7.64	132	298841	39.68	ng/ul	0.00
13) 4-Methylphenol-d8	8.82	113	326578	37.89	ng/ul	0.00
19) Nitrobenzene-d5	9.29	128	153200	35.46	ng/ul	0.00
22) 2-Nitrophenol-d4	10.02	143	189257	38.14	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.57	165	319630	34.15	ng/ul	0.00
29) 4-Chloroaniline-d4	11.08	131	466289	42.47	ng/ul	0.00
43) Dimethylphthalate-d6	14.17	166	1209429	36.97	ng/ul	0.00
46) Acenaphthylene-d8	14.47	160	1379725	37.48	ng/ul	0.00
51) 4-Nitrophenol-d4	14.99	143	191237	34.25	ng/ul	0.00
57) Fluorene-d10	15.77	176	1114694	36.88	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.90	200	210408	32.62	ng/ul	0.00
70) Anthracene-d10	17.64	188	1695575	37.43	ng/ul	0.00
76) Pyrene-d10	19.93	212	1858290	36.83	ng/ul	0.00
87) Benzo(a)pyrene-d12	25.00	264	1719265	38.94	ng/ul	0.00

Target Compounds	Qvalue

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

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