

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG081415\
 Data File : BG018434.D
 Acq On : 14 Aug 2015 18:38
 Operator : UM/MA
 Sample : MDL-BLK-W
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Aug 17 03:17:12 2015
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG080915.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Aug 17 03:14:49 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.02	152	89774	20.00	ng/ul	0.00
18) Naphthalene-d8	10.82	136	417288	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.65	164	271074	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.39	188	669402	20.00	ng/ul	0.00
78) Chrysene-d12	21.65	240	670735	20.00	ng/ul	0.00
86) Perylene-d12	24.85	264	647547	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.45	96	13625	6.71	ng/uL	0.00
5) Phenol-d5	7.17	99	274740	33.72	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.34	67	176176	34.89	ng/ul	0.00
9) 2-Chlorophenol-d4	7.55	132	207765	33.29	ng/ul	0.00
13) 4-Methylphenol-d8	8.72	113	235789	34.44	ng/ul	0.00
19) Nitrobenzene-d5	9.17	128	117012	35.97	ng/ul	0.00
22) 2-Nitrophenol-d4	9.90	143	122466	36.97	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.44	165	230691	32.81	ng/ul	0.00
29) 4-Chloroaniline-d4	10.95	131	325754	41.00	ng/ul	0.00
44) Dimethylphthalate-d6	14.05	166	781141	34.25	ng/ul	0.00
47) Acenaphthylene-d8	14.34	160	975910	33.98	ng/ul	0.00
52) 4-Nitrophenol-d4	14.83	143	151897	37.03	ng/ul	0.00
58) Fluorene-d10	15.63	176	687916	34.63	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.75	200	117901	35.81	ng/ul	0.00
71) Anthracene-d10	17.49	188	1123781	35.10	ng/ul	0.00
79) Pyrene-d10	19.77	212	1205449	34.64	ng/ul	0.00
90) Benzo(a)pyrene-d12	24.64	264	1237411	36.00	ng/ul	0.00

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

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