

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM110115\
 Data File : BM003031.D
 Acq On : 01 Nov 2015 16:37
 Operator : UM/IZ
 Sample : G4210-16
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BC7D6

Quant Time: Nov 02 02:38:19 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM103015.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Nov 02 01:03:38 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.77	152	127351	20.00	ng/ul	0.00
18) Naphthalene-d8	10.56	136	548129	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.41	164	293508	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.15	188	627075	20.00	ng/ul	0.00
78) Chrysene-d12	21.34	240	654257	20.00	ng/ul	0.00
86) Perylene-d12	23.61	264	635548	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.23	96	16882	6.25	ng/uL	0.00
5) Phenol-d5	6.93	99	334176	33.19	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.10	67	201678	34.26	ng/ul	0.00
9) 2-Chlorophenol-d4	7.30	132	286467	33.57	ng/ul	0.00
13) 4-Methylphenol-d8	8.46	113	276082	34.01	ng/ul	0.00
19) Nitrobenzene-d5	8.92	128	123597	36.98	ng/ul	0.00
22) 2-Nitrophenol-d4	9.65	143	129054	38.11	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.17	165	257361	33.88	ng/ul	0.00
29) 4-Chloroaniline-d4	10.69	131	304687	31.00	ng/ul	0.00
44) Dimethylphthalate-d6	13.82	166	842687	38.16	ng/ul	0.00
47) Acenaphthylene-d8	14.10	160	1020706	36.48	ng/ul	0.00
52) 4-Nitrophenol-d4	14.59	143	119337	32.07	ng/ul	0.00
58) Fluorene-d10	15.40	176	730043	38.24	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.52	200	84215	29.00	ng/ul	0.00
71) Anthracene-d10	17.25	188	1084154	39.00	ng/ul	0.00
79) Pyrene-d10	19.55	212	1183943	36.76	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.47	264	1220646	40.42	ng/ul	0.00

Target Compounds						Qvalue
6) Phenol	6.95	94	18452	1.78	ng/ul#	86
45) Dimethylphthalate	13.87	163	502741	22.34	ng/ul	99

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

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