

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM020915\
 Data File : BM000472.D
 Acq On : 10 Feb 2015 23:53
 Operator : TP/IZ
 Sample : G1249-24
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4FN7

Quant Time: Feb 11 04:13:03 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM01.2-EPA-BM020615.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Feb 11 04:03:52 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.72	152	108342	20.00	ng/ul	0.00
18) Naphthalene-d8	10.51	136	444865	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.37	164	274259	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.12	188	596247	20.00	ng/ul	0.00
75) Chrysene-d12	21.31	240	599761	20.00	ng/ul	0.00
83) Perylene-d12	23.55	264	541290	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	0.00	96	0	0.00	ng/uL	
5) Phenol-d5	6.89	99	39802	5.53	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.06	67	94927	21.44	ng/ul	0.00
9) 2-Chlorophenol-d4	7.26	132	113608	18.15	ng/ul	0.00
13) 4-Methylphenol-d8	8.43	113	72429	12.15	ng/ul	0.00
19) Nitrobenzene-d5	8.88	128	62893	22.57	ng/ul	0.00
22) 2-Nitrophenol-d4	9.60	143	73346	23.28	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.13	165	138391	20.39	ng/ul	0.00
29) 4-Chloroaniline-d4	10.65	131	7651	1.23	ng/ul	0.00
43) Dimethylphthalate-d6	13.78	166	468148	24.77	ng/ul	0.00
46) Acenaphthylene-d8	14.06	160	573242	22.20	ng/ul	0.00
51) 4-Nitrophenol-d4	14.57	143	9790	3.17	ng/ul	0.00
57) Fluorene-d10	15.37	176	419799	22.07	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.49	200	59907	18.99	ng/ul	0.00
70) Anthracene-d10	17.22	188	614543	22.66	ng/ul	0.00
76) Pyrene-d10	19.52	212	667811	23.62	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.41	264	522440	21.61	ng/ul	0.00

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

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