

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG022515\
 Data File : BG016038.D
 Acq On : 25 Feb 2015 16:24
 Operator : TP/IZ
 Sample : G1354-04
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0AD9

Quant Time: Feb 26 01:14:51 2015
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM01.2-EPA-MA-BG021915.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Feb 26 00:35:32 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.81	152	93151	20.00	ng/ul	0.00
16) Naphthalene-d8	10.60	136	432171	20.00	ng/ul	0.00
33) Acenaphthene-d10	14.44	164	274533	20.00	ng/ul	0.00
59) Phenanthrene-d10	17.17	188	573663	20.00	ng/ul	0.00
75) Chrysene-d12	21.35	240	571335	20.00	ng/ul	0.00
83) Perylene-d12	23.65	264	568618	20.00	ng/ul	0.00

System Monitoring Compounds

3) Phenol-d5	6.97	99	98661	11.13	ng/ul	0.01
5) Bis-(2-Chloroethyl)ether-d	7.14	67	59286	11.33	ng/ul	0.00
7) 2-Chlorophenol-d4	7.34	132	75965	11.59	ng/ul	0.00
11) 4-Methylphenol-d8	8.50	113	73389	11.19	ng/ul	0.00
17) Nitrobenzene-d5	8.96	128	34887	11.42	ng/ul	0.00
20) 2-Nitrophenol-d4	9.68	143	26230	9.31	ng/ul	0.00
24) 2,4-Dichlorophenol-d3	10.21	165	62049	10.15	ng/ul	0.00
27) 4-Chloroaniline-d4	10.73	131	95711	14.22	ng/ul	0.00
41) Dimethylphthalate-d6	13.84	166	223352	12.17	ng/ul	0.00
44) Acenaphthylene-d8	14.13	160	293651	11.30	ng/ul	0.00
49) 4-Nitrophenol-d4	14.62	143	21455	5.81	ng/ul	0.00
55) Fluorene-d10	15.42	176	207671	11.26	ng/ul	0.00
60) 4,6-Dinitro-2-methylphenol	15.54	200	9328	4.22	ng/ul	0.00
68) Anthracene-d10	17.27	188	304958	11.21	ng/ul	0.00
76) Pyrene-d10	19.56	212	309077	11.52	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.50	264	264142	10.24	ng/ul	0.00

Target Compounds

						Qvalue
2) Benzaldehyde	6.95	77	6905	1.23	ng/ul	82
4) Phenol	6.99	94	12373	1.32	ng/ul	94
12) Acetophenone	8.63	105	14338	1.44	ng/ul	96
42) Dimethylphthalate	13.89	163	94361	4.93	ng/ul	97

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

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