

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM052815\
 Data File : BM001440.D
 Acq On : 28 May 2015 23:54
 Operator : TP/IZ
 Sample : G2406-02MSD
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 KM-SP-002-150518MSD

Quant Time: May 29 04:14:05 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIM-BM052815.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 29 03:45:53 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.38	152	12625	5.00	ng	0.00
6) Naphthalene-d8	10.16	136	43817	5.00	ng	0.00
9) Acenaphthene-d10	14.07	164	26371	5.00	ng	0.00
15) Phenanthrene-d10	16.81	188	61379	5.00	ng	0.00
17) Chrysene-d12	21.03	240	54868	5.00	ng	0.00
21) Perylene-d12	23.11	264	46541	5.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
2) 2-Fluorophenol	4.97	112	20057	10.38	ng	0.00
3) Phenol-d6	6.57	99	26201	12.84	ng	-0.03
7) Nitrobenzene-d5	8.53	82	18184	7.78	ng	0.00
10) 2,4,6-Tribromophenol	15.58	330	11843	10.77	ng	0.00
11) 2-Fluorobiphenyl	12.68	172	51462	7.03	ng	0.00
18) Terphenyl-d14	19.47	244	48084	7.05	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
4) bis(2-Chloroethyl)ether	6.82	93	12580	5.96	ng	75
5) n-Nitroso-di-n-propylamine	8.18	70	11717	8.44	ng	# 73
8) bis(2-Chloroethoxy)methane	9.60	93	19774	6.61	ng	98
12) 2,6-Dinitrotoluene	13.66	165	10213	6.29	ng	99
13) 2,4-Dinitrotoluene	14.46	165	14866	7.79	ng	# 51
14) 4-Chlorophenyl-phenylether	15.11	204	25974	6.40	ng	# 76
16) 4-Bromophenyl-phenylether	16.02	248	15416	6.37	ng	# 68
19) Benzo(a)anthracene	21.00	228	87619	6.40	ng	# 88
20) 3,3'-Dichlorobenzidine	20.95	252	34884	7.51	ng	# 83

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

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