

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM052815\
 Data File : BM001441.D
 Acq On : 29 May 2015 00:30
 Operator : TP/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDCCC001EC

Quant Time: May 29 04:15:06 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	12554	5.00	ng	0.00
6) Naphthalene-d8	10.16	136	41139	5.00	ng	0.00
9) Acenaphthene-d10	14.07	164	24168	5.00	ng	0.00
15) Phenanthrene-d10	16.82	188	53949	5.00	ng	0.01
17) Chrysene-d12	21.03	240	49590	5.00	ng	0.00
21) Perylene-d12	23.11	264	43504	5.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
2) 2-Fluorophenol	4.97	112	1989	1.03	ng	0.00
3) Phenol-d6	6.57	99	2210	1.09	ng	-0.03
7) Nitrobenzene-d5	8.53	82	2212	1.01	ng	0.00
10) 2,4,6-Tribromophenol	15.56	330	791	1.00	ng	0.00
11) 2-Fluorobiphenyl	12.68	172	6932	1.03	ng	0.00
18) Terphenyl-d14	19.47	244	6356	1.03	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
4) bis(2-Chloroethyl)ether	6.82	93	2078	0.99	ng	92
5) n-Nitroso-di-n-propylamine	8.18	70	1485	1.08	ng	83
8) bis(2-Chloroethoxy)methane	9.60	93	2485	1.01	ng	98
12) 2,6-Dinitrotoluene	13.66	165	1033	0.94	ng	# 87
13) 2,4-Dinitrotoluene	14.46	165	1189	0.91	ng	# 79
14) 4-Chlorophenyl-phenylether	15.12	204	3823	1.03	ng	# 89
16) 4-Bromophenyl-phenylether	16.01	248	2204	1.04	ng	# 80
19) Benzo(a)anthracene	21.00	228	12854	1.04	ng	# 85
20) 3,3'-Dichlorobenzidine	20.95	252	3485	1.08	ng	# 75

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

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