

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM052815\
 Data File : BM001434.D
 Acq On : 28 May 2015 19:04
 Operator : TP/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDCCC001EC

Quant Time: May 29 03:40:43 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIM-BM052815.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 28 15:41:25 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.38	152	11559	5.00	ng	0.00
6) Naphthalene-d8	10.16	136	38307	5.00	ng	0.00
9) Acenaphthene-d10	14.07	164	23376	5.00	ng	0.00
15) Phenanthrene-d10	16.82	188	54668	5.00	ng	0.01
17) Chrysene-d12	21.03	240	50347	5.00	ng	0.00
21) Perylene-d12	23.12	264	43087	5.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
2) 2-Fluorophenol	4.98	112	1817	1.02	ng	0.00
3) Phenol-d6	6.59	99	1896	1.02	ng	0.00
7) Nitrobenzene-d5	8.53	82	2046	1.00	ng	0.00
10) 2,4,6-Tribromophenol	15.58	330	699	0.93	ng	0.01
11) 2-Fluorobiphenyl	12.68	172	6643	1.02	ng	0.00
18) Terphenyl-d14	19.47	244	6480	1.04	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
4) bis(2-Chloroethyl)ether	6.82	93	2123	1.09	ng	96
5) n-Nitroso-di-n-propylamine	8.18	70	1246	0.99	ng	94
8) bis(2-Chloroethoxy)methane	9.61	93	2137	0.94	ng	# 89
12) 2,6-Dinitrotoluene	13.66	165	965	0.92	ng	# 54
13) 2,4-Dinitrotoluene	14.46	165	1103	0.88	ng	87
14) 4-Chlorophenyl-phenylether	15.12	204	3679	1.03	ng	# 78
16) 4-Bromophenyl-phenylether	16.03	248	2104	0.97	ng	# 71
19) Benzo(a)anthracene	21.01	228	12286	0.98	ng	100
20) 3,3'-Dichlorobenzidine	20.96	252	2853	0.93	ng	100

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

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