

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM052815\
 Data File : BM001424.D
 Acq On : 28 May 2015 10:44
 Operator : TP/IZ
 Sample : SSTDICC005
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC005

Quant Time: May 28 15:26:35 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 15:25:05 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.38	152	11731	5.00	ng	0.00
6) Naphthalene-d8	10.16	136	39229	5.00	ng	0.00
9) Acenaphthene-d10	14.07	164	23248	5.00	ng	0.00
15) Phenanthrene-d10	16.82	188	53618	5.00	ng	0.01
17) Chrysene-d12	21.03	240	52944	5.00	ng	0.00
21) Perylene-d12	23.12	264	46780	5.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
2) 2-Fluorophenol	4.98	112	9239	4.91	ng	0.00
3) Phenol-d6	6.57	99	10145	5.30	ng	-0.03
7) Nitrobenzene-d5	8.53	82	11530	5.40	ng	0.00
10) 2,4,6-Tribromophenol	15.56	330	4384	6.17	ng	0.00
11) 2-Fluorobiphenyl	12.68	172	32356	4.94	ng	0.00
18) Terphenyl-d14	19.47	244	32064	4.89	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
4) bis(2-Chloroethyl)ether	6.82	93	9389	4.43	ng	85
5) n-Nitroso-di-n-propylamine	8.18	70	7292	5.97	ng	# 87
8) bis(2-Chloroethoxy)methane	9.60	93	13051	6.21	ng	97
12) 2,6-Dinitrotoluene	13.66	165	6698	6.83	ng	# 82
13) 2,4-Dinitrotoluene	14.46	165	8456	7.43	ng	# 61
14) 4-Chlorophenyl-phenylether	15.12	204	18149	5.10	ng	# 85
16) 4-Bromophenyl-phenylether	16.01	248	10470	4.97	ng	# 67
19) Benzo(a)anthracene	21.01	228	64340	19.98	ng	98
20) 3,3'-Dichlorobenzidine	20.96	252	19413	5.81	ng	# 95

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

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