

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM052815\
 Data File : BM001425.D
 Acq On : 28 May 2015 11:20
 Operator : TP/IZ
 Sample : SSTDICC002
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC002

Quant Time: May 28 15:26:46 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	11564	5.00	ng	0.00
6) Naphthalene-d8	10.16	136	38089	5.00	ng	0.00
9) Acenaphthene-d10	14.07	164	22556	5.00	ng	0.00
15) Phenanthrene-d10	16.81	188	51706	5.00	ng	0.00
17) Chrysene-d12	21.03	240	49930	5.00	ng	0.00
21) Perylene-d12	23.12	264	42932	5.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
2) 2-Fluorophenol	4.98	112	3714	2.00	ng	0.00
3) Phenol-d6	6.57	99	3959	2.10	ng	-0.03
7) Nitrobenzene-d5	8.53	82	4351	2.10	ng	0.00
10) 2,4,6-Tribromophenol	15.57	330	1555	2.26	ng	0.00
11) 2-Fluorobiphenyl	12.68	172	13109	2.06	ng	0.00
18) Terphenyl-d14	19.47	244	12927	2.09	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
4) bis(2-Chloroethyl)ether	6.82	93	3875	1.86	ng	90
5) n-Nitroso-di-n-propylamine	8.18	70	2802	2.33	ng	# 87
8) bis(2-Chloroethoxy)methane	9.60	93	4952	2.43	ng	98
12) 2,6-Dinitrotoluene	13.65	165	2194	2.31	ng	# 29
13) 2,4-Dinitrotoluene	14.45	165	2609	2.36	ng	# 84
14) 4-Chlorophenyl-phenylether	15.12	204	7368	2.13	ng	97
16) 4-Bromophenyl-phenylether	16.02	248	4448	2.19	ng	94
19) Benzo(a)anthracene	21.01	228	25070	8.26	ng	99
20) 3,3'-Dichlorobenzidine	20.97	252	6891	2.19	ng	# 96

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

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