

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM052815\
 Data File : BM001428.D
 Acq On : 28 May 2015 13:45
 Operator : TP/IZ
 Sample : SSTDICC0.5
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC0.5

Quant Time: May 28 15:27:18 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	10845	5.00	ng	0.00
6) Naphthalene-d8	10.16	136	34296	5.00	ng	0.00
9) Acenaphthene-d10	14.07	164	20685	5.00	ng	0.00
15) Phenanthrene-d10	16.81	188	46031	5.00	ng	0.00
17) Chrysene-d12	21.02	240	45866	5.00	ng	0.00
21) Perylene-d12	23.12	264	39853	5.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
2) 2-Fluorophenol	4.98	112	836	0.48	ng	0.00
3) Phenol-d6	6.59	99	843	0.48	ng	0.00
7) Nitrobenzene-d5	8.53	82	895	0.48	ng	0.00
10) 2,4,6-Tribromophenol	15.57	330	277	0.44	ng	0.00
11) 2-Fluorobiphenyl	12.68	172	2902	0.50	ng	0.00
18) Terphenyl-d14	19.48	244	2711	0.48	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
4) bis(2-Chloroethyl)ether	6.82	93	966	0.49	ng	90
5) n-Nitroso-di-n-propylamine	8.18	70	558	0.49	ng	# 92
8) bis(2-Chloroethoxy)methane	9.61	93	831	0.45	ng	# 81
12) 2,6-Dinitrotoluene	13.67	165	370	0.42	ng	# 83
13) 2,4-Dinitrotoluene	14.46	165	404	0.40	ng	# 85
14) 4-Chlorophenyl-phenylether	15.12	204	1609	0.51	ng	95
16) 4-Bromophenyl-phenylether	16.02	248	913	0.51	ng	97
19) Benzo(a)anthracene	21.01	228	5683	2.04	ng	94
20) 3,3'-Dichlorobenzidine	20.96	252	1138	0.39	ng	# 87

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

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