

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM052815\
 Data File : BM001423.D
 Acq On : 28 May 2015 10:08
 Operator : TP/IZ
 Sample : SSTDICC010
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC010

Quant Time: May 28 15:39:29 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	12467	5.00	ng	0.00
6) Naphthalene-d8	10.16	136	43697	5.00	ng	0.00
9) Acenaphthene-d10	14.07	164	26490	5.00	ng	0.00
15) Phenanthrene-d10	16.81	188	61486	5.00	ng	0.00
17) Chrysene-d12	21.03	240	54241	5.00	ng	0.00
21) Perylene-d12	23.12	264	47721	5.00	ng	0.00
System Monitoring Compounds						
2) 2-Fluorophenol	4.97	112	19871	9.93	ng	0.00
3) Phenol-d6	6.57	99	22348	10.98	ng	-0.03
7) Nitrobenzene-d5	8.53	82	27117	11.41	ng	0.00
10) 2,4,6-Tribromophenol	15.58	330	11252	13.90	ng	0.00
11) 2-Fluorobiphenyl	12.68	172	71813	9.62	ng	0.00
18) Terphenyl-d14	19.48	244	71854	10.70	ng	0.00
Target Compounds						
4) bis(2-Chloroethyl)ether	6.82	93	20041	8.91	ng	Qvalue 83
5) n-Nitroso-di-n-propylamine	8.18	70	16562	12.77	ng	# 94
8) bis(2-Chloroethoxy)methane	9.60	93	30211	12.92	ng	97
12) 2,6-Dinitrotoluene	13.66	165	16880	15.11	ng	# 83
14) 4-Chlorophenyl-phenylether	15.11	204	40253	9.93	ng	# 65
16) 4-Bromophenyl-phenylether	16.02	248	24928	10.32	ng	# 75
19) Benzo(a)anthracene	21.01	228	141104	42.77	ng	98
20) 3,3'-Dichlorobenzidine	20.95	252	47780	13.95	ng	# 72

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

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