

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM053115\
 Data File : BM001504.D
 Acq On : 31 May 2015 09:55
 Operator : TP/IZ
 Sample : G2438-03MSD
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 OM-SP-002-150528MSD

Manual Integrations
 APPROVED

apatel
 6/1/2015 8:45:07 PM

Quant Time: Jun 01 05:55:49 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIM-BM053115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun May 31 07:43:08 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.38	152	13284	5.00	ng	0.00
6) Naphthalene-d8	10.15	136	51021	5.00	ng	0.00
9) Acenaphthene-d10	14.07	164	27602	5.00	ng	0.00
15) Phenanthrene-d10	16.81	188	56533	5.00	ng	0.00
17) Chrysene-d12	21.01	240	43484	5.00	ng	-0.01
21) Perylene-d12	23.11	264	37970	5.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
2) 2-Fluorophenol	4.97	112	33502	12.82	ng	0.00
3) Phenol-d6	6.57	99	50273	14.74	ng	0.00
7) Nitrobenzene-d5	8.53	82	25496	9.25	ng	0.00
10) 2,4,6-Tribromophenol	15.57	330	16112	16.47	ng	0.00
11) 2-Fluorobiphenyl	12.67	172	72472	9.06	ng	0.00
18) Terphenyl-d14	19.47	244	53569	9.21	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
4) bis(2-Chloroethyl)ether	6.80	93	24958	9.49	ng	99
5) n-Nitroso-di-n-propylamine	8.18	70	20737	9.65	ng	# 92
8) bis(2-Chloroethoxy)methane	9.58	93	33010	10.03	ng	97
12) 2,6-Dinitrotoluene	13.65	165	16541	9.69	ng	# 86
13) 2,4-Dinitrotoluene	14.45	165	22759	10.38	ng	# 65
14) 4-Chlorophenyl-phenylether	15.11	204	38261	9.55	ng	95
16) 4-Bromophenyl-phenylether	16.02	248	22401	9.51	ng	# 93
19) Benzo(a)anthracene	21.00	228	109807	9.26	ng	# 89
20) 3,3'-Dichlorobenzidine	20.95	252	44672m	11.54	ng	

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

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