

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM052815\
 Data File : BM001433.D
 Acq On : 28 May 2015 18:27
 Operator : TP/IZ
 Sample : PB83634BS
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB83634BS

Quant Time: May 29 03:38:41 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 15:41:25 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.38	152	13063	5.00	ng	0.00
6) Naphthalene-d8	10.16	136	46286	5.00	ng	0.00
9) Acenaphthene-d10	14.07	164	29206	5.00	ng	0.00
15) Phenanthrene-d10	16.81	188	67776	5.00	ng	0.00
17) Chrysene-d12	21.03	240	59283	5.00	ng	0.00
21) Perylene-d12	23.11	264	50697	5.00	ng	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
2) 2-Fluorophenol	4.97	112	23360	11.65	ng	0.00
3) Phenol-d6	6.57	99	29015	13.79	ng	-0.03
7) Nitrobenzene-d5	8.53	82	21345	8.62	ng	0.00
10) 2,4,6-Tribromophenol	15.57	330	15644	12.80	ng	0.00
11) 2-Fluorobiphenyl	12.68	172	62925	7.76	ng	0.00
18) Terphenyl-d14	19.47	244	62636	8.50	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
4) bis(2-Chloroethyl)ether	6.82	93	16364	7.43	ng	78
5) n-Nitroso-di-n-propylamine	8.18	70	15757	11.02	ng	# 79
8) bis(2-Chloroethoxy)methane	9.60	93	26056	8.21	ng	99
12) 2,6-Dinitrotoluene	13.65	165	15051	8.28	ng	# 44
13) 2,4-Dinitrotoluene	14.45	165	22334	10.48	ng	# 68
14) 4-Chlorophenyl-phenylether	15.12	204	37716	8.42	ng	97
16) 4-Bromophenyl-phenylether	16.02	248	23279	8.70	ng	92
19) Benzo(a)anthracene	21.00	228	124055	8.37	ng	# 85
20) 3,3'-Dichlorobenzidine	20.95	252	43602	8.65	ng	# 74

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

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