

Data Path : Z:\HPCHEM1\BNA M\DATA\BM030116\
 Data File : BM004522.D
 Acq On : 02 Mar 2016 16:21
 Operator : SJ/UM
 Sample : H1509-06DL 5X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MW-3DL

Quant Time: Mar 02 18:02:37 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-SIM-BM030116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 02 14:46:57 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	17393	5.00	ng	0.00
7) Naphthalene-d8	10.39	136	68991	5.00	ng	0.00
13) Acenaphthene-d10	14.26	164	37018	5.00	ng	0.00
19) Phenanthrene-d10	17.01	188	84484	5.00	ng	0.00
26) Chrysene-d12	21.23	240	65430	5.00	ng	0.00
35) Perylene-d12	23.45	264	61038	5.00	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.24	112	1202	0.33	ng	0.00
5) Phenol-d6	6.87	99	982	0.33	ng	0.03
8) Nitrobenzene-d5	8.81	82	1720	0.60	ng	0.01
14) 2,4,6-Tribromophenol	15.77	330	2236	1.82	ng	-0.02
15) 2-Fluorobiphenyl	12.88	172	8544	0.75	ng	0.00
29) Terphenyl-d14	19.67	244	8884	0.82	ng	0.00
Target Compounds						
17) Acenaphthene	14.33	154	744	0.04	ng	# 45
18) Fluorene	15.33	166	1072	0.06	ng	# 93
22) Pentachlorophenol	16.71	266	2998	4.53	ng	# 87
33) Bis(2-ethylhexyl)phthalate	21.12	149	941	0.18	ng	# 98

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

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