

Data Path : S:\HPCHEM1\BNA_M\DATA\BM022916\
 Data File : BM004489.D
 Acq On : 01 Mar 2016 03:12
 Operator : SJ/UM
 Sample : H1509-04
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MW-4

Quant Time: Mar 01 07:45:33 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-SIM-BM022916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 01 00:52:28 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	15960	5.00	ng	0.00
7) Naphthalene-d8	10.39	136	64985	5.00	ng	0.00
13) Acenaphthene-d10	14.26	164	35569	5.00	ng	0.00
19) Phenanthrene-d10	17.01	188	75927	5.00	ng	0.00
26) Chrysene-d12	21.23	240	65313	5.00	ng	-0.01
35) Perylene-d12	23.45	264	62833	5.00	ng	-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.22	112	10015	2.70	ng	0.00
5) Phenol-d6	6.81	99	7261	1.73	ng	0.00
8) Nitrobenzene-d5	8.78	82	14510	4.39	ng	0.00
14) 2,4,6-Tribromophenol	15.77	330	11848	8.67	ng	-0.03
15) 2-Fluorobiphenyl	12.88	172	53787	4.94	ng	0.00
29) Terphenyl-d14	19.67	244	52281	6.22	ng	0.00
Target Compounds						
23) Phenanthrene	17.06	178	1656	0.08	ng	# 79
24) Anthracene	17.16	178	629	0.03	ng	# 93
25) Fluoranthene	19.11	202	3479	0.12	ng	# 93
28) Pyrene	19.47	202	3320	0.13	ng	# 98
33) Bis(2-ethylhexyl)phthalate	21.12	149	6014	0.32	ng	# 91
34) Indeno(1,2,3-cd)pyrene	25.70	276	886	0.04	ng	# 95
40) Benzo(g,h,i)perylene	26.39	276	1131	0.05	ng	# 53

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

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