

Data Path : Z:\HPCHEM1\BNA F\DATA\BF102616\
 Data File : BF090847.D
 Acq On : 26 Oct 2016 18:34
 Operator : UM/SJ
 Sample : H5308-17
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-105D

Quant Time: Oct 27 02:15:16 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF102616.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 26 16:35:31 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.78	152	207441	20.00	ng	-0.01
21) Naphthalene-d8	8.08	136	887811	20.00	ng	-0.01
38) Acenaphthene-d10	9.82	164	432753	20.00	ng	-0.02
63) Phenanthrene-d10	11.31	188	761502	20.00	ng	-0.01
75) Chrysene-d12	13.95	240	691809	20.00	ng	-0.01
86) Perylene-d12	15.37	264	517104	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.38	112	814213	68.63	ng	-0.01
7) Phenol-d6	6.42	99	651972	40.73	ng	-0.02
23) Nitrobenzene-d5	7.36	82	903180	66.60	ng	-0.01
41) 2,4,6-Tribromophenol	10.62	330	631254	141.94	ng	-0.01
44) 2-Fluorobiphenyl	9.15	172	1391261	56.45	ng	-0.02
78) Terphenyl-d14	12.90	244	1590976	57.94	ng	-0.02
Target Compounds						
10) Phenol	6.43	94	106462	6.04	ng	# 53
17) 2-Methylphenol	7.05	107	132073	11.86	ng	# 84
20) 3+4-Methylphenols	7.21	107	203427	15.78	ng	# 86
32) Benzoic acid	7.90	122	116799	12.93	ng	# 65
49) Dimethylphthalate	9.55	163	119226	4.12	ng	# 97

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

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