

Data Path : Z:\HPCHEM1\BNA F\DATA\BF102116\
 Data File : BF090741.D
 Acq On : 22 Oct 2016 5:34
 Operator : UM/SJ
 Sample : H5308-17
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-105D

Quant Time: Oct 22 06:37:23 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF102116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 21 17:36:31 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	263604	20.00	ng	-0.01
21) Naphthalene-d8	8.14	136	1133992	20.00	ng	0.00
38) Acenaphthene-d10	9.89	164	536663	20.00	ng	-0.01
63) Phenanthrene-d10	11.38	188	831097	20.00	ng	-0.01
75) Chrysene-d12	14.02	240	558919	20.00	ng	0.00
86) Perylene-d12	15.45	264	388780	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	1089154	66.63	ng	0.01
7) Phenol-d6	6.49	99	804492	36.34	ng	-0.01
23) Nitrobenzene-d5	7.42	82	1198329	57.96	ng	0.00
41) 2,4,6-Tribromophenol	10.69	330	678691	159.78	ng	0.00
44) 2-Fluorobiphenyl	9.22	172	1763856	51.02	ng	-0.01
78) Terphenyl-d14	12.97	244	1658801	68.02	ng	0.00
Target Compounds						
10) Phenol	6.50	94	133444	5.36	ng	# 55
17) 2-Methylphenol	7.11	107	156305	10.57	ng	# 85
20) 3+4-Methylphenols	7.26	107	263453	15.03	ng	# 82
32) Benzoic acid	7.97	122	126523	16.86	ng	# 66
49) Dimethylphthalate	9.62	163	126191	3.74	ng	# 98

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

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