

Data Path : Z:\HPCHEM1\BNA F\DATA\BF011317\
 Data File : BF092259.D
 Acq On : 13 Jan 2017 23:21
 Operator : UM/SJ
 Sample : I1179-01
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-1

Quant Time: Jan 16 02:40:45 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF122116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 16 01:02:36 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.58	152	176114	20.00	ng	-0.01
21) Naphthalene-d8	7.87	136	695781	20.00	ng	-0.01
38) Acenaphthene-d10	9.62	164	348850	20.00	ng	-0.01
63) Phenanthrene-d10	11.09	188	519549	20.00	ng	-0.01
75) Chrysene-d12	13.72	240	437887	20.00	ng	-0.01
86) Perylene-d12	15.07	264	347481	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.17	112	868581	82.35	ng	0.00
7) Phenol-d6	6.24	99	688003	53.43	ng	-0.01
23) Nitrobenzene-d5	7.15	82	1062799	84.94	ng	-0.01
41) 2,4,6-Tribromophenol	10.42	330	385181	133.42	ng	0.00
44) 2-Fluorobiphenyl	8.95	172	1993710	122.84	ng	0.00
78) Terphenyl-d14	12.68	244	1495966	82.86	ng	-0.01
Target Compounds						Qvalue
10) Phenol	6.26	94	48006	3.27	ng	87
31) Naphthalene	7.88	128	404185	12.69	ng	98
37) 2-Methylnaphthalene	8.58	142	314320	12.73	ng	98
49) Dimethylphthalate	9.35	163	73868	3.31	ng	# 87
51) Acenaphthene	9.65	154	47150	2.39	ng	92
57) Fluorene	10.16	166	81972	4.23	ng	# 94
70) Phenanthrene	11.11	178	66118	2.35	ng	# 85

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

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