

Data Path : Z:\HPCHEM1\BNA F\DATA\BF010918\
 Data File : BF101970.D
 Acq On : 9 Jan 2018 23:18
 Operator : SJ/JU
 Sample : J1049-09
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-C

Quant Time: Jan 10 03:37:51 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF010918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 09 15:20:36 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.73	152	218998	20.00	ng	0.00
21) Naphthalene-d8	8.01	136	794564	20.00	ng	0.00
38) Acenaphthene-d10	9.76	164	269601	20.00	ng	0.00
63) Phenanthrene-d10	11.25	188	413158	20.00	ng	0.00
75) Chrysene-d12	13.89	240	384255	20.00	ng	0.01
86) Perylene-d12	15.31	264	294508	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.33	112	595270	38.38	ng	0.00
7) Phenol-d6	6.35	99	407241	22.01	ng	0.00
23) Nitrobenzene-d5	7.29	82	976031	72.18	ng	0.00
41) 2,4,6-Tribromophenol	10.55	330	222110	94.40	ng	0.00
44) 2-Fluorobiphenyl	9.09	172	1367014	79.22	ng	0.00
78) Terphenyl-d14	12.84	244	1092928	59.05	ng	0.00
Target Compounds						
10) Phenol	6.37	94	584658	27.959	ng	98
20) 3+4-Methylphenols	7.13	107	109005	6.470	ng	97
27) 2,4-Dimethylphenol	7.67	122	42275	3.722	ng	96
49) Dimethylphthalate	9.49	163	78100	3.611	ng	99

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

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