

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF041020\
 Data File : BF119900.D
 Acq On : 10 Apr 2020 17:45
 Operator : MA/SJ
 Sample : L2178-01
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-41

Quant Time: Apr 11 02:05:02 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 10 13:48:19 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.75	152	176387	20.00	ng	0.00
21) Naphthalene-d8	8.03	136	662230	20.00	ng	0.00
39) Acenaphthene-d10	9.79	164	371403	20.00	ng	0.00
64) Phenanthrene-d10	11.28	188	724025	20.00	ng	0.00
76) Chrysene-d12	13.92	240	532821	20.00	ng	0.00
87) Perylene-d12	15.36	264	434985	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.36	112	697524	68.34	ng	0.00
7) Phenol-d6	6.38	99	554082	42.13	ng	-0.01
23) Nitrobenzene-d5	7.32	82	1272867	99.44	ng	0.00
42) 2,4,6-Tribromophenol	10.59	330	613362	134.89	ng	0.00
45) 2-Fluorobiphenyl	9.12	172	2377759	90.90	ng	0.00
79) Terphenyl-d14	12.87	244	2693517	85.06	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.51	163	95997	3.342	ng	98
52) Acenaphthene	9.82	154	68798	2.808	ng	98
71) Phenanthrene	11.30	178	239071	6.204	ng	99
75) Fluoranthene	12.49	202	114933	2.764	ng	96

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

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