

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF041020\
 Data File : BF119897.D
 Acq On : 10 Apr 2020 16:25
 Operator : MA/SJ
 Sample : L2178-05
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-1A

Quant Time: Apr 11 01:42:11 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	184617	20.00	ng	0.00
21) Naphthalene-d8	8.04	136	710113	20.00	ng	0.00
39) Acenaphthene-d10	9.79	164	374707	20.00	ng	0.00
64) Phenanthrene-d10	11.28	188	667638	20.00	ng	0.00
76) Chrysene-d12	13.92	240	438737	20.00	ng	0.00
87) Perylene-d12	15.36	264	362992	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.36	112	754572	70.63	ng	0.00
7) Phenol-d6	6.38	99	594107	43.16	ng	-0.01
23) Nitrobenzene-d5	7.32	82	1316721	95.93	ng	0.00
42) 2,4,6-Tribromophenol	10.59	330	606290	132.16	ng	0.00
45) 2-Fluorobiphenyl	9.12	172	2368069	89.73	ng	0.00
79) Terphenyl-d14	12.87	244	2258330	86.61	ng	0.00
Target Compounds						Qvalue
38) 1-Methylnaphthalene	8.85	142	115964	4.857	ng	96
50) Dimethylphthalate	9.51	163	76505	2.640	ng	98
52) Acenaphthene	9.82	154	112867	4.566	ng	96
58) Fluorene	10.34	166	69789	2.573	ng	97
71) Phenanthrene	11.30	178	86134	2.424	ng	96
73) Carbazole	11.51	167	102901	2.998	ng	96

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

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