

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF070121\
 Data File : BF124511.D
 Acq On : 1 Jul 2021 14:54
 Operator : JU/CG
 Sample : M2875-10DL 5X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-14DL

Quant Time: Jul 01 15:24:08 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF062321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 01 14:15:13 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.681	152	29088	20.000	ng	0.00
21) Naphthalene-d8	7.969	136	113301	20.000	ng	# 0.00
39) Acenaphthene-d10	9.716	164	67231	20.000	ng	0.00
64) Phenanthrene-d10	11.204	188	123605	20.000	ng	0.00
76) Chrysene-d12	13.851	240	84522	20.000	ng	0.00
86) Perylene-d12	15.274	264	75595	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.316	112	13629	7.369	ng	0.01
7) Phenol-d6	6.351	99	12200	5.052	ng	0.01
23) Nitrobenzene-d5	7.251	82	39999	14.638	ng	0.00
42) 2,4,6-Tribromophenol	10.510	330	18971	23.618	ng	0.00
45) 2-Fluorobiphenyl	9.039	172	88457	16.117	ng	0.00
79) Terphenyl-d14	12.786	244	108486	19.224	ng	0.00
Target Compounds						
						Qvalue
31) Naphthalene	8.004	128	2768639	415.345	ng	97
37) 2-Methylnaphthalene	8.681	142	434908	92.198	ng	100
38) 1-Methylnaphthalene	8.781	142	401840	91.588	ng	93
46) 1,1'-Biphenyl	9.139	154	33190	5.808	ng	97
52) Acenaphthene	9.751	154	91971	21.438	ng	97
58) Fluorene	10.263	166	33798	6.433	ng	84
71) Phenanthrene	11.227	178	85497	10.973	ng	98
72) Anthracene	11.280	178	19784	2.548	ng	97
78) Pyrene	12.651	202	17670	2.290	ng	# 92

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

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