

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF011022\
 Data File : BF126564.D
 Acq On : 10 Jan 2022 14:22
 Operator : JU/CG
 Sample : M5011-17DL 25X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HSB-8-(10-10.5)DL

Quant Time: Jan 10 17:16:33 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 10 14:50:31 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.769	152	93771	20.000	ng	0.00
21) Naphthalene-d8	8.057	136	388218	20.000	ng	# 0.00
39) Acenaphthene-d10	9.810	164	188106	20.000	ng	0.00
64) Phenanthrene-d10	11.298	188	294248	20.000	ng	# 0.00
76) Chrysene-d12	13.939	240	175404	20.000	ng	0.00
86) Perylene-d12	15.404	264	188829	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.375	112	18182	2.784	ng	0.00
7) Phenol-d6	6.398	99	32196	3.807	ng	-0.01
23) Nitrobenzene-d5	7.328	82	25196	3.332	ng	-0.01
42) 2,4,6-Tribromophenol	10.598	330	2217	1.525	ng	0.00
45) 2-Fluorobiphenyl	9.128	172	36548	3.419	ng	-0.01
79) Terphenyl-d14	12.880	244	28047	3.107	ng	-0.01
Target Compounds						
						Qvalue
31) Naphthalene	8.075	128	167754	9.155	ng	# 97
37) 2-Methylnaphthalene	8.769	142	358442	33.160	ng	100
38) 1-Methylnaphthalene	8.869	142	198562	19.000	ng	100
71) Phenanthrene	11.322	178	44999	2.991	ng	96

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

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