

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF010722\
 Data File : BF126547.D
 Acq On : 7 Jan 2022 17:41
 Operator : JU/CG
 Sample : M5011-11DL 10X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HSB-5-(11.5-12)DL

Quant Time: Jan 08 04:39:04 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 05 01:57:07 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.781	152	117017	20.000	ng	0.00
21) Naphthalene-d8	8.069	136	439197	20.000	ng	0.00
39) Acenaphthene-d10	9.822	164	237723	20.000	ng	0.00
64) Phenanthrene-d10	11.304	188	383007	20.000	ng	0.00
76) Chrysene-d12	13.951	240	226203	20.000	ng	# 0.00
86) Perylene-d12	15.409	264	195385	20.000	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.387	112	58723	7.207	ng	0.00
7) Phenol-d6	6.410	99	109609	10.386	ng	-0.01
23) Nitrobenzene-d5	7.339	82	67989	7.949	ng	-0.01
42) 2,4,6-Tribromophenol	10.604	330	18052	9.827	ng	-0.01
45) 2-Fluorobiphenyl	9.139	172	106749	7.902	ng	-0.01
79) Terphenyl-d14	12.892	244	94702	8.134	ng	0.00
Target Compounds						
						Qvalue
31) Naphthalene	8.086	128	933852	45.048	ng	97
37) 2-Methylnaphthalene	8.780	142	607899	49.710	ng	98
38) 1-Methylnaphthalene	8.875	142	303051	25.632	ng	99
71) Phenanthrene	11.327	178	40709	2.079	ng	96

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

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