

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF021822\
 Data File : BF126960.D
 Acq On : 18 Feb 2022 19:03
 Operator : CG\JU
 Sample : N1600-06
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 C0AA7

Quant Time: Feb 18 23:53:21 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF021622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 16 14:55:48 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.933	152	91120	20.000	ng	0.00
21) Naphthalene-d8	8.222	136	342376	20.000	ng	# 0.00
39) Acenaphthene-d10	9.986	164	176424	20.000	ng	0.01
64) Phenanthrene-d10	11.474	188	275453	20.000	ng	0.01
76) Chrysene-d12	14.104	240	161383	20.000	ng	# 0.00
86) Perylene-d12	15.609	264	147844	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.551	112	548434	93.025	ng	0.00
7) Phenol-d6	6.557	99	744570	93.596	ng	0.00
23) Nitrobenzene-d5	7.498	82	541350	71.369	ng	0.00
42) 2,4,6-Tribromophenol	10.774	330	159779	78.659	ng	0.00
45) 2-Fluorobiphenyl	9.298	172	740921	61.826	ng	0.00
79) Terphenyl-d14	13.051	244	660829	60.195	ng	0.00
Target Compounds						
						Qvalue
31) Naphthalene	8.245	128	113511	6.351	ng	# 78
37) 2-Methylnaphthalene	8.939	142	200983	17.330	ng	98
38) 1-Methylnaphthalene	9.033	142	130465	11.748	ng	95
46) 1,1'-Biphenyl	9.398	154	77537	5.547	ng	95
71) Phenanthrene	11.498	178	38015	2.466	ng	# 85
78) Pyrene	12.915	202	37468	2.672	ng	# 93

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

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