

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111121\
 Data File : BF126134.D
 Acq On : 11 Nov 2021 19:50
 Operator : JU/CG
 Sample : M4630-13
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-13W-16.0-16.5

Quant Time: Nov 12 09:42:40 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF110321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 03 17:21:25 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.851	152	132628	20.000	ng	0.00
21) Naphthalene-d8	8.134	136	470430	20.000	ng	# 0.00
39) Acenaphthene-d10	9.875	164	281362	20.000	ng	-0.02
64) Phenanthrene-d10	11.363	188	526097	20.000	ng	#-0.01
76) Chrysene-d12	14.004	240	352194	20.000	ng	# 0.00
86) Perylene-d12	15.468	264	392330	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.457	112	692279	89.637	ng	0.00
7) Phenol-d6	6.469	99	923183	84.796	ng	-0.02
23) Nitrobenzene-d5	7.551	82	38004	3.810	ng	0.13
42) 2,4,6-Tribromophenol	10.663	330	300797	93.287	ng	-0.02
45) 2-Fluorobiphenyl	9.198	172	1187086	57.015	ng	-0.02
79) Terphenyl-d14	12.951	244	1215809	55.667	ng	-0.01
Target Compounds						Qvalue
27) 2,4-Dimethylphenol	7.798	122	30566	4.668	ng	# 47
31) Naphthalene	8.157	128	1125035	46.172	ng	99
37) 2-Methylnaphthalene	8.839	142	118896	7.065	ng	98
38) 1-Methylnaphthalene	8.933	142	53317	3.271	ng	98
75) Fluoranthene	12.574	202	87306	2.748	ng	95
78) Pyrene	12.804	202	91319	3.233	ng	99

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

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