

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111022\
 Data File : BF131135.D
 Acq On : 10 Nov 2022 21:41
 Operator : CG\JU
 Sample : N5494-11
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-24

Quant Time: Nov 11 00:51:42 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110922.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 10 04:15:17 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.887	152	65417	20.000	ng	0.00
21) Naphthalene-d8	8.175	136	258071	20.000	ng	0.00
39) Acenaphthene-d10	9.933	164	131480	20.000	ng	0.00
64) Phenanthrene-d10	11.422	188	172785	20.000	ng	0.00
76) Chrysene-d12	14.074	240	136716	20.000	ng	0.00
86) Perylene-d12	15.568	264	131885	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.510	112	453442	117.879	ng	0.00
7) Phenol-d6	6.516	99	575402	114.416	ng	0.00
23) Nitrobenzene-d5	7.451	82	362909	63.448	ng	-0.01
42) 2,4,6-Tribromophenol	10.722	330	124584	91.686	ng	-0.01
45) 2-Fluorobiphenyl	9.251	172	656689	69.723	ng	0.00
79) Terphenyl-d14	13.010	244	420285	53.900	ng	0.00
Target Compounds						
87) Indeno(1,2,3-cd)pyrene	17.074	276	4000	4.206	ng	# 94
92) Benzo(g,h,i)perylene	17.539	276	4111	4.769	ng	# 89

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

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