

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF022122\
 Data File : BF126993.D
 Acq On : 22 Feb 2022 00:32
 Operator : CG\JU
 Sample : N1602-06
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 20220082-AMENDED-BERKELEY-2-COMP

Quant Time: Feb 22 03:54:51 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF021622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 21 13:16:57 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.916	152	86543	20.000	ng	0.00
21) Naphthalene-d8	8.198	136	350344	20.000	ng	# 0.00
39) Acenaphthene-d10	9.951	164	184115	20.000	ng	-0.01
64) Phenanthrene-d10	11.445	188	325168	20.000	ng	# 0.00
76) Chrysene-d12	14.086	240	139010	20.000	ng	#-0.01
86) Perylene-d12	15.580	264	151710	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.546	112	759669	135.669	ng	0.00
7) Phenol-d6	6.540	99	1113386	147.359	ng	0.00
23) Nitrobenzene-d5	7.481	82	780357	100.539	ng	-0.01
42) 2,4,6-Tribromophenol	10.745	330	200588	94.624	ng	0.00
45) 2-Fluorobiphenyl	9.275	172	1151657	92.085	ng	0.00
79) Terphenyl-d14	13.033	244	1045146	110.525	ng	0.00
Target Compounds						
10) Phenol	6.551	94	17459	2.287	ng	# 60

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

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