

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF021122\
 Data File : BF126858.D
 Acq On : 11 Feb 2022 15:35
 Operator : CG\JU
 Sample : N1477-03
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 20220087-AMENDED-BERKELEY-7-COMP

Quant Time: Feb 11 22:42:45 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF020922.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 10 00:44:00 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.940	152	115798	20.000	ng	0.00
21) Naphthalene-d8	8.222	136	477938	20.000	ng	0.00
39) Acenaphthene-d10	9.981	164	253866	20.000	ng	0.00
64) Phenanthrene-d10	11.469	188	481908	20.000	ng	0.00
76) Chrysene-d12	14.110	240	314703	20.000	ng	-0.01
86) Perylene-d12	15.621	264	235746	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.563	112	966770	128.542	ng	0.00
7) Phenol-d6	6.563	99	1339122	131.890	ng	0.00
23) Nitrobenzene-d5	7.504	82	901826	87.062	ng	0.00
42) 2,4,6-Tribromophenol	10.775	330	362599	133.229	ng	0.00
45) 2-Fluorobiphenyl	9.298	172	1330117	78.869	ng	0.00
79) Terphenyl-d14	13.063	244	1559965	71.061	ng	0.00
Target Compounds						
10) Phenol	6.569	94	26622	2.601	ng	Qvalue 86

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

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