

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF021422\
 Data File : BF126872.D
 Acq On : 14 Feb 2022 15:46
 Operator : CG\JU
 Sample : N1477-04
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 20220088-AMENDED-BERKELEY-8-COMP

Quant Time: Feb 15 00:59:40 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	6.939	152	116975	20.000	ng	0.00
21) Naphthalene-d8	8.222	136	478977	20.000	ng	0.00
39) Acenaphthene-d10	9.980	164	250553	20.000	ng	0.00
64) Phenanthrene-d10	11.468	188	426018	20.000	ng	0.00
76) Chrysene-d12	14.109	240	203344	20.000	ng	-0.01
86) Perylene-d12	15.621	264	205688	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.563	112	998394	131.411	ng	0.00
7) Phenol-d6	6.563	99	1367820	133.361	ng	0.00
23) Nitrobenzene-d5	7.504	82	975264	93.947	ng	0.00
42) 2,4,6-Tribromophenol	10.769	330	341910	127.288	ng	0.00
45) 2-Fluorobiphenyl	9.298	172	1485796	89.265	ng	0.00
79) Terphenyl-d14	13.057	244	1347713	95.013	ng	0.00
Target Compounds						
10) Phenol	6.569	94	29246	2.829	ng	Qvalue # 42

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

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