

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF022122\
 Data File : BF126989.D
 Acq On : 21 Feb 2022 22:40
 Operator : CG\JU
 Sample : N1602-02
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 20220076-BERKELEY-3-COMP

Quant Time: Feb 22 03:53:11 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	85475	20.000	ng	0.00
21) Naphthalene-d8	8.198	136	351504	20.000	ng	# 0.00
39) Acenaphthene-d10	9.957	164	187137	20.000	ng	0.00
64) Phenanthrene-d10	11.445	188	331398	20.000	ng	# 0.00
76) Chrysene-d12	14.086	240	142212	20.000	ng	#-0.01
86) Perylene-d12	15.580	264	159352	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.540	112	534595	96.666	ng	0.00
7) Phenol-d6	6.534	99	699699	93.764	ng	-0.01
23) Nitrobenzene-d5	7.475	82	486951	62.530	ng	-0.02
42) 2,4,6-Tribromophenol	10.745	330	144637	67.129	ng	0.00
45) 2-Fluorobiphenyl	9.275	172	757538	59.593	ng	0.00
79) Terphenyl-d14	13.027	244	704624	72.837	ng	-0.01

Target Compounds	Qvalue

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

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