

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112222\
 Data File : BF131338.D
 Acq On : 22 Nov 2022 19:34
 Operator : CG\JU
 Sample : N5598-13
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FRAC-TANK-SLURRY-0760533

Quant Time: Nov 23 04:18:40 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF112122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 22 01:17:52 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.963	152	105811	20.000	ng	0.00
21) Naphthalene-d8	8.251	136	384926	20.000	ng	0.00
39) Acenaphthene-d10	10.016	164	217474	20.000	ng	-0.01
64) Phenanthrene-d10	11.510	188	389691	20.000	ng	-0.01
76) Chrysene-d12	14.169	240	194835	20.000	ng	#-0.01
86) Perylene-d12	15.710	264	210443	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.587	112	718342	129.208	ng	0.02
7) Phenol-d6	6.581	99	824892	116.290	ng	0.00
23) Nitrobenzene-d5	7.528	82	670097	87.754	ng	-0.01
42) 2,4,6-Tribromophenol	10.810	330	295334	161.418	ng	0.00
45) 2-Fluorobiphenyl	9.334	172	1242848	83.142	ng	0.00
79) Terphenyl-d14	13.110	244	1237595	103.872	ng	0.00

Target Compounds Qvalue

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

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