

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF093022\
 Data File : BF130489.D
 Acq On : 30 Sep 2022 17:47
 Operator : CG\JU
 Sample : N4898-01RE
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CRUSHED-MASONRY-PILE-11RE

Quant Time: Oct 01 01:36:32 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 28 02:59:31 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.634	152	99081	20.000	ng	-0.01
21) Naphthalene-d8	7.916	136	374641	20.000	ng	-0.01
39) Acenaphthene-d10	9.669	164	182716	20.000	ng	-0.01
64) Phenanthrene-d10	11.145	188	291979	20.000	ng	-0.01
76) Chrysene-d12	13.780	240	187615	20.000	ng	-0.01
86) Perylene-d12	15.168	264	226933	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.245	112	34523	6.043	ng	0.00
7) Phenol-d6	6.275	99	355572	49.747	ng	-0.02
23) Nitrobenzene-d5	7.204	82	438115	59.744	ng	-0.01
42) 2,4,6-Tribromophenol	10.463	330	223	0.127	ng	0.00
45) 2-Fluorobiphenyl	8.998	172	844068	67.270	ng	-0.01
79) Terphenyl-d14	12.733	244	740613	65.390	ng	-0.01
Target Compounds						
75) Fluoranthene	12.357	202	33015	2.084	ng	Qvalue 96

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

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