

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF093022\
 Data File : BF130488.D
 Acq On : 30 Sep 2022 17:15
 Operator : CG\JU
 Sample : N4897-01RE
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CRUSHED-MASONRY-PILE-10RE

Quant Time: Oct 01 01:36:20 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	106458	20.000	ng	-0.01
21) Naphthalene-d8	7.916	136	410682	20.000	ng	-0.01
39) Acenaphthene-d10	9.669	164	204847	20.000	ng	-0.01
64) Phenanthrene-d10	11.151	188	329968	20.000	ng	0.00
76) Chrysene-d12	13.780	240	191260	20.000	ng	-0.01
86) Perylene-d12	15.168	264	222507	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.245	112	19130	3.117	ng	0.00
7) Phenol-d6	6.275	99	252896	32.930	ng	-0.02
23) Nitrobenzene-d5	7.204	82	493075	61.338	ng	-0.01
42) 2,4,6-Tribromophenol	10.463	330	133	0.068	ng	0.00
45) 2-Fluorobiphenyl	8.998	172	940807	66.879	ng	-0.01
79) Terphenyl-d14	12.733	244	815002	70.587	ng	-0.01

Target Compounds	Qvalue

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

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