

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081222\
 Data File : BF129743.D
 Acq On : 12 Aug 2022 14:33
 Operator : CG\JU
 Sample : N4082-01RE
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CRUSHED-MASONARY-PILE-ARE

Quant Time: Aug 12 15:55:20 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 12 15:52:45 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.828	152	131286	20.000	ng	0.00
21) Naphthalene-d8	8.110	136	524596	20.000	ng	0.00
39) Acenaphthene-d10	9.869	164	275397	20.000	ng	0.00
64) Phenanthrene-d10	11.357	188	448584	20.000	ng	0.00
76) Chrysene-d12	14.004	240	272292	20.000	ng	0.00
86) Perylene-d12	15.480	264	259174	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.475	112	199960	24.811	ng	0.00
7) Phenol-d6	6.487	99	619976	61.202	ng	0.00
23) Nitrobenzene-d5	7.398	82	647593	67.387	ng	0.00
42) 2,4,6-Tribromophenol	10.663	330	10876	4.108	ng	0.00
45) 2-Fluorobiphenyl	9.186	172	1233683	69.298	ng	0.00
79) Terphenyl-d14	12.945	244	1172858	81.262	ng	0.00

Target Compounds	Qvalue

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

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