

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080122\
 Data File : BF129470.D
 Acq On : 01 Aug 2022 13:23
 Operator : CG\JU
 Sample : N3969-01RE
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CHAMBERS-STREET-A3-9065X-MHRE

Quant Time: Aug 02 01:15:32 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 02 01:11:25 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.875	152	150591	20.000	ng	0.00
21) Naphthalene-d8	8.157	136	578840	20.000	ng	0.00
39) Acenaphthene-d10	9.916	164	280800	20.000	ng	0.00
64) Phenanthrene-d10	11.404	188	406654	20.000	ng	0.00
76) Chrysene-d12	14.045	240	287492	20.000	ng	0.00
86) Perylene-d12	15.527	264	343285	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.516	112	283850	30.705	ng	0.01
7) Phenol-d6	6.528	99	205512	17.687	ng	0.00
23) Nitrobenzene-d5	7.439	82	774824	73.071	ng	0.00
42) 2,4,6-Tribromophenol	10.710	330	165819	61.422	ng	0.00
45) 2-Fluorobiphenyl	9.233	172	1336361	73.621	ng	0.00
79) Terphenyl-d14	12.986	244	1009298	66.232	ng	0.00

Target Compounds	Qvalue

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

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