

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082322\
 Data File : BF129958.D
 Acq On : 23 Aug 2022 17:48
 Operator : CG\JU
 Sample : N4120-07
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BLDG-19C

Quant Time: Aug 24 01:56:29 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF081422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 23 04:41:40 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.775	152	144663	20.000	ng	0.00
21) Naphthalene-d8	8.057	136	588037	20.000	ng	0.00
39) Acenaphthene-d10	9.810	164	312830	20.000	ng	0.00
64) Phenanthrene-d10	11.298	188	505589	20.000	ng	0.00
76) Chrysene-d12	13.951	240	288027	20.000	ng	0.00
86) Perylene-d12	15.404	264	246742	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.434	112	1086954	124.404	ng	0.02
7) Phenol-d6	6.440	99	1294777	118.336	ng	0.00
23) Nitrobenzene-d5	7.345	82	967099	88.040	ng	0.00
42) 2,4,6-Tribromophenol	10.610	330	453317	144.381	ng	0.00
45) 2-Fluorobiphenyl	9.134	172	1816777	88.205	ng	0.00
79) Terphenyl-d14	12.892	244	1784516	103.092	ng	0.00

Target Compounds	Qvalue

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

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