

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091224\
 Data File : BF139342.D
 Acq On : 12 Sep 2024 16:48
 Operator : RC/JU
 Sample : P3949-01
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BREAKER-2546-BASE

Quant Time: Sep 12 17:30:36 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF090424.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 04 16:11:43 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.875	152	46106	20.000	ng	-0.02
21) Naphthalene-d8	8.157	136	176633	20.000	ng	-0.02
39) Acenaphthene-d10	9.916	164	95989	20.000	ng	-0.01
64) Phenanthrene-d10	11.398	188	181154	20.000	ng	-0.02
76) Chrysene-d12	14.033	240	95174	20.000	ng	#-0.02
86) Perylene-d12	15.504	264	85325	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.498	112	162750	59.051	ng	0.00
7) Phenol-d6	6.498	99	274630	74.641	ng	-0.02
23) Nitrobenzene-d5	7.440	82	225365	66.376	ng	-0.01
42) 2,4,6-Tribromophenol	10.698	330	42719	51.538	ng	-0.02
45) 2-Fluorobiphenyl	9.234	172	444363	71.460	ng	-0.02
79) Terphenyl-d14	12.986	244	450477	76.421	ng	-0.01

Target Compounds	Qvalue

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

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