

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112224\
 Data File : BF140567.D
 Acq On : 22 Nov 2024 10:44
 Operator : RC/JU
 Sample : P4910-08
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MH-759

Quant Time: Nov 22 11:19:17 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF112124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 21 15:23:48 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.869	152	95916	20.000	ng	0.00
21) Naphthalene-d8	8.151	136	362240	20.000	ng	0.00
39) Acenaphthene-d10	9.904	164	199732	20.000	ng	-0.01
64) Phenanthrene-d10	11.398	188	383423	20.000	ng	0.00
76) Chrysene-d12	14.045	240	235981	20.000	ng	0.00
86) Perylene-d12	15.527	264	192364	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.510	112	749572	133.338	ng	0.01
7) Phenol-d6	6.510	99	888052	119.493	ng	0.00
23) Nitrobenzene-d5	7.433	82	703247	99.302	ng	0.00
42) 2,4,6-Tribromophenol	10.704	330	324941	152.115	ng	0.00
45) 2-Fluorobiphenyl	9.227	172	1269026	94.666	ng	0.00
79) Terphenyl-d14	12.986	244	1517055	100.104	ng	0.00
Target Compounds						
77) Benzidine	12.739	184	32788	4.779	ng	Qvalue 98

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

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