

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF090524\
 Data File : BF139252.D
 Acq On : 05 Sep 2024 10:05
 Operator : RC/JU
 Sample : PB163110BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB163110BL

Quant Time: Sep 05 10:48:24 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	6.887	152	163073	20.000	ng	0.00
21) Naphthalene-d8	8.169	136	612665	20.000	ng	0.00
39) Acenaphthene-d10	9.928	164	333239	20.000	ng	0.00
64) Phenanthrene-d10	11.410	188	624390	20.000	ng	0.00
76) Chrysene-d12	14.051	240	304010	20.000	ng	0.00
86) Perylene-d12	15.527	264	268948	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.516	112	1155636	118.550	ng	0.01
7) Phenol-d6	6.516	99	1527169	117.352	ng	0.00
23) Nitrobenzene-d5	7.451	82	970599	82.416	ng	0.00
42) 2,4,6-Tribromophenol	10.716	330	329308	114.440	ng	0.00
45) 2-Fluorobiphenyl	9.245	172	1709667	79.196	ng	0.00
79) Terphenyl-d14	13.004	244	1647739	87.510	ng	0.00

Target Compounds	Qvalue

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

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