

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF090624\
 Data File : BF139269.D
 Acq On : 06 Sep 2024 10:01
 Operator : RC/JU
 Sample : PB163139BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB163139BL

Quant Time: Sep 06 10:43:41 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	6.886	152	163587	20.000	ng	0.00
21) Naphthalene-d8	8.169	136	609350	20.000	ng	0.00
39) Acenaphthene-d10	9.922	164	331699	20.000	ng	0.00
64) Phenanthrene-d10	11.410	188	608322	20.000	ng	0.00
76) Chrysene-d12	14.051	240	296509	20.000	ng	0.00
86) Perylene-d12	15.527	264	264558	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.516	112	1137784	116.352	ng	0.01
7) Phenol-d6	6.516	99	1489553	114.102	ng	0.00
23) Nitrobenzene-d5	7.451	82	949420	81.056	ng	0.00
42) 2,4,6-Tribromophenol	10.716	330	324649	113.344	ng	0.00
45) 2-Fluorobiphenyl	9.245	172	1685709	78.449	ng	0.00
79) Terphenyl-d14	12.998	244	1615248	87.955	ng	0.00

Target Compounds	Qvalue

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

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