

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101024\
 Data File : BF139738.D
 Acq On : 10 Oct 2024 11:04
 Operator : RC/JU
 Sample : PB164011BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB164011BL

Quant Time: Oct 10 11:53:04 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF100124.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 02 04:58:47 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.869	152	113971	20.000	ng	0.00
21) Naphthalene-d8	8.145	136	437153	20.000	ng	0.00
39) Acenaphthene-d10	9.904	164	256984	20.000	ng	0.00
64) Phenanthrene-d10	11.392	188	521320	20.000	ng	0.00
76) Chrysene-d12	14.033	240	375941	20.000	ng	0.00
86) Perylene-d12	15.510	264	293210	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.504	112	840554	127.830	ng	0.02
7) Phenol-d6	6.504	99	1092425	123.539	ng	0.00
23) Nitrobenzene-d5	7.434	82	719378	83.317	ng	0.00
42) 2,4,6-Tribromophenol	10.698	330	325125	131.058	ng	0.00
45) 2-Fluorobiphenyl	9.228	172	1366299	81.621	ng	0.00
79) Terphenyl-d14	12.980	244	1770482	77.274	ng	0.00

Target Compounds	Qvalue

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

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