

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF100724\
 Data File : BF139687.D
 Acq On : 07 Oct 2024 15:33
 Operator : RC/JU
 Sample : PB163902BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB163902BL

Quant Time: Oct 07 15:55:32 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	89954	20.000	ng	0.00
21) Naphthalene-d8	8.145	136	344839	20.000	ng	0.00
39) Acenaphthene-d10	9.904	164	195984	20.000	ng	0.00
64) Phenanthrene-d10	11.386	188	402197	20.000	ng	0.00
76) Chrysene-d12	14.027	240	291108	20.000	ng	0.00
86) Perylene-d12	15.498	264	222477	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.499	112	536394	103.353	ng	0.01
7) Phenol-d6	6.498	99	720169	103.186	ng	0.00
23) Nitrobenzene-d5	7.428	82	664322	97.538	ng	0.00
42) 2,4,6-Tribromophenol	10.692	330	184092	97.305	ng	0.00
45) 2-Fluorobiphenyl	9.222	172	1208928	94.698	ng	0.00
79) Terphenyl-d14	12.980	244	1573522	88.692	ng	0.00

Target Compounds	Qvalue

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

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