

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102324\
 Data File : BF139974.D
 Acq On : 23 Oct 2024 19:54
 Operator : RC/JU
 Sample : P4468-04
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ETGI-329

Quant Time: Oct 24 01:11:51 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF101824.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 18 15:07:50 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.892	152	141549	20.000	ng	0.00
21) Naphthalene-d8	8.169	136	537624	20.000	ng	0.00
39) Acenaphthene-d10	9.922	164	260732	20.000	ng	0.00
64) Phenanthrene-d10	11.410	188	391613	20.000	ng	0.00
76) Chrysene-d12	14.051	240	328805	20.000	ng	0.00
86) Perylene-d12	15.527	264	301905	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.516	112	1138779	125.945	ng	0.02
7) Phenol-d6	6.510	99	1321009	112.804	ng	0.00
23) Nitrobenzene-d5	7.451	82	973311	100.339	ng	0.00
42) 2,4,6-Tribromophenol	10.710	330	326192	133.751	ng	0.00
45) 2-Fluorobiphenyl	9.245	172	1608344	101.948	ng	0.00
79) Terphenyl-d14	12.998	244	1478682	73.280	ng	0.00

Target Compounds	Qvalue

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

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