

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102324\
Data File : BF139973.D
Acq On : 23 Oct 2024 19:25
Operator : RC/JU
Sample : P4468-06
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
ETGI-345

Quant Time: Oct 24 01:11:38 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	147918	20.000	ng	0.00
21) Naphthalene-d8	8.169	136	541314	20.000	ng	0.00
39) Acenaphthene-d10	9.922	164	268595	20.000	ng	0.00
64) Phenanthrene-d10	11.410	188	403583	20.000	ng	0.00
76) Chrysene-d12	14.051	240	340738	20.000	ng	0.00
86) Perylene-d12	15.527	264	309184	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.516	112	1098393	116.248	ng	0.02
7) Phenol-d6	6.510	99	1252340	102.336	ng	0.00
23) Nitrobenzene-d5	7.451	82	962374	98.535	ng	0.00
42) 2,4,6-Tribromophenol	10.710	330	327504	130.357	ng	0.00
45) 2-Fluorobiphenyl	9.245	172	1575128	96.920	ng	0.00
79) Terphenyl-d14	12.998	244	1450770	69.379	ng	0.00
Target Compounds						Qvalue
77) Benzidine	12.745	184	117123	20.904	ng	99

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

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