

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101824\
Data File : BF139864.D
Acq On : 18 Oct 2024 20:05
Operator : RC/JU
Sample : P4426-07
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PAD-4

Quant Time: Oct 18 23:08:54 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	128568	20.000	ng	0.00
21) Naphthalene-d8	8.169	136	429920	20.000	ng	0.00
39) Acenaphthene-d10	9.922	164	184268	20.000	ng	0.00
64) Phenanthrene-d10	11.410	188	333846	20.000	ng	0.00
76) Chrysene-d12	14.051	240	286044	20.000	ng	0.00
86) Perylene-d12	15.527	264	194650	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.498	112	319297	38.879	ng	0.00
7) Phenol-d6	6.498	99	525323	49.388	ng	-0.01
23) Nitrobenzene-d5	7.445	82	404245	52.114	ng	0.00
42) 2,4,6-Tribromophenol	10.710	330	51333	29.783	ng	0.00
45) 2-Fluorobiphenyl	9.245	172	715511	64.174	ng	0.00
79) Terphenyl-d14	12.998	244	910676	51.878	ng	0.00

Target Compounds	Qvalue

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

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