

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101024\
Data File : BF139760.D
Acq On : 10 Oct 2024 21:47
Operator : RC/JU
Sample : SP6638
Misc : 8270 SURROGATES
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SP6638

Quant Time: Oct 10 23:21:04 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	6.863	152	80871	20.000	ng	0.00
21) Naphthalene-d8	8.145	136	315403	20.000	ng	0.00
39) Acenaphthene-d10	9.898	164	176404	20.000	ng	-0.01
64) Phenanthrene-d10	11.386	188	324448	20.000	ng	0.00
76) Chrysene-d12	14.027	240	174947	20.000	ng	0.00
86) Perylene-d12	15.504	264	176647	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.498	112	704334	150.954	ng	0.01
7) Phenol-d6	6.498	99	914065	145.677	ng	0.00
23) Nitrobenzene-d5	7.428	82	593609	95.289	ng	0.00
42) 2,4,6-Tribromophenol	10.692	330	226938	133.266	ng	0.00
45) 2-Fluorobiphenyl	9.222	172	1103928	96.071	ng	0.00
79) Terphenyl-d14	12.974	244	1049620	98.444	ng	0.00

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

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