

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101124\
 Data File : BF139784.D
 Acq On : 11 Oct 2024 19:37
 Operator : RC/JU
 Sample : P4364-08 5X
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-03-2.0-2.5

Quant Time: Oct 12 02:02:04 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	6.869	152	58355	20.000	ng	0.00
21) Naphthalene-d8	8.151	136	188402	20.000	ng	0.00
39) Acenaphthene-d10	9.904	164	96219	20.000	ng	0.00
64) Phenanthrene-d10	11.392	188	202295	20.000	ng	0.00
76) Chrysene-d12	14.033	240	124604	20.000	ng	0.00
86) Perylene-d12	15.510	264	93962	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.487	112	70786	21.025	ng	0.00
7) Phenol-d6	6.504	99	78207	17.273	ng	0.00
23) Nitrobenzene-d5	7.428	82	44789	12.036	ng	0.00
42) 2,4,6-Tribromophenol	10.692	330	17648	19.000	ng	0.00
45) 2-Fluorobiphenyl	9.222	172	86852	13.857	ng	0.00
79) Terphenyl-d14	12.974	244	117776	15.509	ng	0.00
Target Compounds						
						Qvalue
71) Phenanthrene	11.416	178	20496	2.149	ng	97
75) Fluoranthene	12.604	202	31501	3.021	ng	97
78) Pyrene	12.833	202	34489	3.095	ng	99

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

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