

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF092024\
 Data File : BF139512.D
 Acq On : 20 Sep 2024 20:17
 Operator : RC/JU
 Sample : P4103-17
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-23

Quant Time: Sep 20 23:22:57 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091324.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 14 02:58:23 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.881	152	78495	20.000	ng	0.00
21) Naphthalene-d8	8.157	136	271926	20.000	ng	# 0.00
39) Acenaphthene-d10	9.910	164	125895	20.000	ng	0.00
64) Phenanthrene-d10	11.398	188	219540	20.000	ng	0.00
76) Chrysene-d12	14.033	240	129376	20.000	ng	# 0.00
86) Perylene-d12	15.510	264	96172	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.510	112	621394	129.143	ng	0.02
7) Phenol-d6	6.510	99	782507	123.900	ng	0.00
23) Nitrobenzene-d5	7.439	82	523366	90.481	ng	0.00
42) 2,4,6-Tribromophenol	10.704	330	159310	118.861	ng	0.00
45) 2-Fluorobiphenyl	9.233	172	825028	91.911	ng	0.00
79) Terphenyl-d14	12.980	244	846386	107.381	ng	0.00
Target Compounds						
38) 1-Methylnaphthalene	8.969	142	22757	2.572	ng	98
70) Pentachlorophenol	11.204	266	3278	2.140	ng	99

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

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