

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF121824\
 Data File : BF140903.D
 Acq On : 18 Dec 2024 14:09
 Operator : RC/JU
 Sample : P5270-07DL2 20X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW12-20241212DL2

Quant Time: Dec 18 14:38:34 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF121624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 16 16:59:50 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.845	152	61168	20.000	ng	0.00
21) Naphthalene-d8	8.128	136	250458	20.000	ng	0.00
39) Acenaphthene-d10	9.880	164	151708	20.000	ng	0.00
64) Phenanthrene-d10	11.374	188	301660	20.000	ng	0.00
76) Chrysene-d12	14.033	240	164695	20.000	ng	0.00
86) Perylene-d12	15.533	264	181077	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.510	112	10519	2.831	ng	0.02
7) Phenol-d6	6.528	99	9418	1.914	ng	0.03
23) Nitrobenzene-d5	7.416	82	22844	4.563	ng	0.00
42) 2,4,6-Tribromophenol	10.686	330	10257	5.717	ng	0.00
45) 2-Fluorobiphenyl	9.198	172	57863	5.244	ng	0.00
79) Terphenyl-d14	12.957	244	58661	5.616	ng	0.00
Target Compounds						
						Qvalue
31) Naphthalene	8.151	128	701165	50.964	ng	100
37) 2-Methylnaphthalene	8.839	142	82011	9.011	ng	99
38) 1-Methylnaphthalene	8.939	142	301927	33.545	ng	100
46) 1,1'-Biphenyl	9.304	154	38702	3.208	ng	98
52) Acenaphthene	9.916	154	48726	5.624	ng	99
58) Fluorene	10.433	166	27581m	2.490	ng	
71) Phenanthrene	11.398	178	50501	3.255	ng	99

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

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