

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF121724\
 Data File : BF140885.D
 Acq On : 17 Dec 2024 15:12
 Operator : RC/JU
 Sample : P5270-07DL 10X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW12-20241212DL

Quant Time: Dec 17 15:37:42 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	6.851	152	51871	20.000	ng	0.00
21) Naphthalene-d8	8.128	136	191262	20.000	ng	0.00
39) Acenaphthene-d10	9.881	164	103044	20.000	ng	0.00
64) Phenanthrene-d10	11.369	188	183841	20.000	ng	0.00
76) Chrysene-d12	14.022	240	133693	20.000	ng	-0.01
86) Perylene-d12	15.510	264	142721	20.000	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.510	112	18952	6.015	ng	0.02
7) Phenol-d6	6.522	99	13698	3.282	ng	0.02
23) Nitrobenzene-d5	7.416	82	35254	9.222	ng	0.00
42) 2,4,6-Tribromophenol	10.681	330	13526	11.099	ng	0.00
45) 2-Fluorobiphenyl	9.198	172	77247	10.307	ng	0.00
79) Terphenyl-d14	12.951	244	71093	8.385	ng	0.00
Target Compounds						
						Qvalue
31) Naphthalene	8.151	128	1006135	95.766	ng	100
37) 2-Methylnaphthalene	8.839	142	109006	15.685	ng	98
38) 1-Methylnaphthalene	8.939	142	404096	58.792	ng	99
46) 1,1'-Biphenyl	9.304	154	51877	6.331	ng	97
52) Acenaphthene	9.916	154	64096	10.893	ng	98
58) Fluorene	10.434	166	35159	4.674	ng	95
71) Phenanthrene	11.398	178	56926	6.021	ng	99

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

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