

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF070324\
 Data File : BF138456.D
 Acq On : 03 Jul 2024 20:07
 Operator : CG\JU
 Sample : P3073-02DL2 10X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW12-20240625DL2

Quant Time: Jul 04 01:00:31 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF062824.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jun 29 02:44:30 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.875	152	107443	20.000	ng	0.00
21) Naphthalene-d8	8.157	136	372851	20.000	ng	0.00
39) Acenaphthene-d10	9.904	164	158995	20.000	ng	-0.01
64) Phenanthrene-d10	11.392	188	264250	20.000	ng	0.00
76) Chrysene-d12	14.027	240	195711	20.000	ng	0.00
86) Perylene-d12	15.498	264	140163	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.492	112	52613	8.119	ng	0.00
7) Phenol-d6	6.504	99	46863	5.368	ng	-0.01
23) Nitrobenzene-d5	7.428	82	67686	9.430	ng	-0.02
42) 2,4,6-Tribromophenol	10.692	330	19014	14.637	ng	-0.01
45) 2-Fluorobiphenyl	9.222	172	116629	11.674	ng	-0.01
79) Terphenyl-d14	12.974	244	136126	9.532	ng	0.00
Target Compounds						
						Qvalue
31) Naphthalene	8.180	128	1312109	69.612	ng	99
37) 2-Methylnaphthalene	8.863	142	113925	9.352	ng	97
38) 1-Methylnaphthalene	8.963	142	408470	34.575	ng	98
46) 1,1'-Biphenyl	9.327	154	60246	5.075	ng	97
52) Acenaphthene	9.939	154	61596	7.067	ng	98
58) Fluorene	10.451	166	24954	2.615	ng	97
71) Phenanthrene	11.416	178	61574	4.772	ng	99

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

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