

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF070524\
 Data File : BF138467.D
 Acq On : 05 Jul 2024 14:27
 Operator : CG\JU
 Sample : P3073-02DL2 10X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW12-20240625DL2

Quant Time: Jul 05 14:50:02 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	6.869	152	83792	20.000	ng	0.00
21) Naphthalene-d8	8.151	136	329952	20.000	ng	0.00
39) Acenaphthene-d10	9.904	164	181442	20.000	ng	-0.01
64) Phenanthrene-d10	11.392	188	291188	20.000	ng	0.00
76) Chrysene-d12	14.027	240	173018	20.000	ng	0.00
86) Perylene-d12	15.498	264	172152	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.493	112	42084	8.327	ng	0.00
7) Phenol-d6	6.504	99	37384	5.491	ng	-0.01
23) Nitrobenzene-d5	7.428	82	62989	9.917	ng	-0.02
42) 2,4,6-Tribromophenol	10.692	330	24957	16.835	ng	-0.01
45) 2-Fluorobiphenyl	9.222	172	120993	10.612	ng	-0.01
79) Terphenyl-d14	12.974	244	115492	9.148	ng	0.00
Target Compounds						
						Qvalue
31) Naphthalene	8.181	128	1317854	79.007	ng	99
37) 2-Methylnaphthalene	8.863	142	124439	11.543	ng	99
38) 1-Methylnaphthalene	8.963	142	445806	42.641	ng	99
46) 1,1'-Biphenyl	9.328	154	71463	5.275	ng	97
52) Acenaphthene	9.939	154	80116	8.054	ng	98
58) Fluorene	10.451	166	36212	3.325	ng	93
71) Phenanthrene	11.416	178	78364	5.512	ng	98

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

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