

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF010824\
 Data File : BF137006.D
 Acq On : 08 Jan 2024 21:06
 Operator : CG\JU
 Sample : P1038-02DL 2X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-2DL

Quant Time: Jan 09 00:44:06 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122023.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 21 01:54:25 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.781	152	51141	20.000	ng	-0.01
21) Naphthalene-d8	8.063	136	189767	20.000	ng	0.00
39) Acenaphthene-d10	9.822	164	85286	20.000	ng	0.00
64) Phenanthrene-d10	11.316	188	135408	20.000	ng	0.00
76) Chrysene-d12	13.969	240	107902	20.000	ng	# 0.00
86) Perylene-d12	15.451	264	132771	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.428	112	138442	42.238	ng	0.00
7) Phenol-d6	6.434	99	119788	28.205	ng	0.00
23) Nitrobenzene-d5	7.345	82	198145	53.430	ng	-0.01
42) 2,4,6-Tribromophenol	10.616	330	64777	64.491	ng	0.00
45) 2-Fluorobiphenyl	9.139	172	385481	60.791	ng	-0.01
79) Terphenyl-d14	12.904	244	306222	39.660	ng	-0.01
Target Compounds						
						Qvalue
31) Naphthalene	8.081	128	85003	8.154	ng	# 93
37) 2-Methylnaphthalene	8.775	142	312783	46.619	ng	99
38) 1-Methylnaphthalene	8.875	142	194968	29.619	ng	98
46) 1,1'-Biphenyl	9.239	154	50850	7.105	ng	98
52) Acenaphthene	9.851	154	20597	3.996	ng	# 82
58) Fluorene	10.369	166	32863	5.300	ng	# 93
71) Phenanthrene	11.339	178	91528	13.174	ng	96

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

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