

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF060122\
 Data File : BF128633.D
 Acq On : 01 Jun 2022 20:23
 Operator : CG\JU
 Sample : N3087-08DL2 25X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-14DL2

Quant Time: Jun 02 09:37:41 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF052622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 27 01:51:59 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.834	152	100710	20.000	ng	0.00
21) Naphthalene-d8	8.116	136	340662	20.000	ng	-0.01
39) Acenaphthene-d10	9.869	164	155570	20.000	ng	-0.01
64) Phenanthrene-d10	11.351	188	261065	20.000	ng	-0.02
76) Chrysene-d12	13.998	240	251449	20.000	ng	#-0.01
86) Perylene-d12	15.462	264	236534	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.510	112	11574	1.904	ng	0.06
7) Phenol-d6	6.534	99	8354	1.078	ng	0.06
23) Nitrobenzene-d5	7.392	82	22759	3.854	ng	-0.02
42) 2,4,6-Tribromophenol	10.657	330	6251	4.221	ng	-0.02
45) 2-Fluorobiphenyl	9.186	172	42084	4.395	ng	-0.02
79) Terphenyl-d14	12.939	244	35698	2.743	ng	-0.02
Target Compounds						
						Qvalue
31) Naphthalene	8.139	128	1348607	80.885	ng	99
37) 2-Methylnaphthalene	8.828	142	154942	14.722	ng	100
38) 1-Methylnaphthalene	8.928	142	177890	17.379	ng	99
52) Acenaphthene	9.898	154	33019	4.062	ng	92
71) Phenanthrene	11.380	178	25162	2.035	ng	# 94

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

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