

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091022\
 Data File : BN021677.D
 Acq On : 10 Sep 2022 01:10
 Operator : CG/JU
 Sample : N4537-04
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 MW1

Quant Time: Sep 10 02:12:26 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN082922.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 09 23:19:14 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.062	152	5014	0.400	ng	0.00
7) Naphthalene-d8	10.888	136	17111	0.400	ng	#-0.01
13) Acenaphthene-d10	14.718	164	11346	0.400	ng	# 0.00
19) Phenanthrene-d10	17.460	188	21229	0.400	ng	#-0.01
29) Chrysene-d12	21.655	240	15841	0.400	ng	# 0.00
35) Perylene-d12	24.165	264	10078	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.592	112	1904	0.194	ng	0.00
5) Phenol-d6	7.217	99	1443	0.115	ng	0.00
8) Nitrobenzene-d5	9.233	82	4274	0.370	ng	0.00
11) 2-Methylnaphthalene-d10	12.475	152	9317	0.242	ng	0.00
14) 2,4,6-Tribromophenol	16.208	330	2116	0.568	ng	0.00
15) 2-Fluorobiphenyl	13.339	172	12276	0.247	ng	0.00
27) Fluoranthene-d10	19.493	212	20569	0.377	ng	0.00
31) Terphenyl-d14	20.079	244	13464	0.378	ng	0.00
Target Compounds						
9) Naphthalene	10.941	128	4498	0.089	ng	# 69
18) Fluorene	15.733	166	4550	0.098	ng	# 8
34) Bis(2-ethylhexyl)phtha...	21.548	149	6449	0.290	ng	# 97

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

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