

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100622\
 Data File : BN022019.D
 Acq On : 06 Oct 2022 14:36
 Operator : CG/JU
 Sample : N4785-08
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PT-ACIDS-WP

Quant Time: Oct 07 00:40:16 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN100322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 04 01:32:21 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.978	152	4496	0.400	ng	0.00
7) Naphthalene-d8	10.805	136	14306	0.400	ng	-0.01
13) Acenaphthene-d10	14.653	164	7007	0.400	ng	0.00
19) Phenanthrene-d10	17.400	188	14020	0.400	ng	#-0.01
29) Chrysene-d12	21.609	240	9833	0.400	ng	0.00
35) Perylene-d12	24.090	264	5282	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.508	112	2016422	193.607	ng	0.00
5) Phenol-d6	7.148	99	2722592	201.465	ng	0.00
8) Nitrobenzene-d5	9.161	82	1369039	119.026	ng	0.00
11) 2-Methylnaphthalene-d10	12.417	152	2034	0.077	ng	0.00
14) 2,4,6-Tribromophenol	16.146	330	674061	252.228	ng	0.00
15) 2-Fluorobiphenyl	13.274	172	2628726	85.864	ng	-0.01
27) Fluoranthene-d10	19.445	212	19	0.001	ng	0.00
31) Terphenyl-d14	20.041	244	1944780	98.377	ng	0.00
Target Compounds						
12) 2-Methylnaphthalene	12.107	142	1595605	249.202	ng	# 80
20) 4,6-Dinitro-2-methylph...	15.774	198	310421	112.948	ng	# 52

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100622\
Data File : BN022019.D
Acq On : 06 Oct 2022 14:36
Operator : CG/JU
Sample : N4785-08
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
PT-ACIDS-WP

Quant Time: Oct 07 00:40:16 2022
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN100322.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Oct 04 01:32:21 2022
Response via : Initial Calibration

