

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100622\
 Data File : BN022020.D
 Acq On : 06 Oct 2022 15:14
 Operator : CG/JU
 Sample : N4785-08DL 25X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PT-ACIDS-WPDL

Quant Time: Oct 07 00:40:27 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	7.986	152	4034	0.400	ng	0.00
7) Naphthalene-d8	10.816	136	12660	0.400	ng	# 0.00
13) Acenaphthene-d10	14.653	164	6332	0.400	ng	0.00
19) Phenanthrene-d10	17.400	188	12862	0.400	ng	#-0.01
29) Chrysene-d12	21.609	240	9216	0.400	ng	0.00
35) Perylene-d12	24.090	264	6505	0.400	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.508	112	69946	7.485	ng	0.00
5) Phenol-d6	7.133	99	94790	7.818	ng	-0.01
8) Nitrobenzene-d5	9.151	82	43619	4.285	ng	-0.01
11) 2-Methylnaphthalene-d10	12.380	152	39	0.002	ng	-0.03
14) 2,4,6-Tribromophenol	16.146	330	20126	8.334	ng	0.00
15) 2-Fluorobiphenyl	13.274	172	120546	4.357	ng	-0.01
27) Fluoranthene-d10	19.445	212	4	0.000	ng	0.00
31) Terphenyl-d14	20.033	244	103984	5.612	ng	0.00
Target Compounds						
12) 2-Methylnaphthalene	12.104	142	56610	9.991	ng	# 80
20) 4,6-Dinitro-2-methylph...	15.774	198	4568	2.001	ng	# 55

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

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