

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN033022\
 Data File : BN019242.D
 Acq On : 30 Mar 2022 15:16
 Operator : CG/JU
 Sample : N1870-08
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 PT-ACIDS-WP

Quant Time: Mar 30 15:49:08 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN033022.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 30 12:57:13 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.833	152	3484	0.400	ng	0.00
7) Naphthalene-d8	10.637	136	9281	0.400	ng	# 0.00
13) Acenaphthene-d10	14.474	164	4946	0.400	ng	0.00
19) Phenanthrene-d10	17.215	188	9646	0.400	ng	0.00
29) Chrysene-d12	21.405	240	6107	0.400	ng	# 0.00
35) Perylene-d12	23.767	264	4740	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.384	112	1462244	173.652	ng	0.00
5) Phenol-d6	6.995	99	1833746	200.520	ng	0.00
8) Nitrobenzene-d5	8.982	82	946624	133.988	ng	-0.01
11) 2-Methylnaphthalene-d10	12.235	152	2	0.000	ng	0.00
14) 2,4,6-Tribromophenol	15.964	330	557821	225.892	ng	0.00
15) 2-Fluorobiphenyl	13.095	172	1967765	93.009	ng	-0.01
27) Fluoranthene-d10	19.265	212	108	0.004	ng	0.02
31) Terphenyl-d14	19.843	244	1552485	113.395	ng	0.00
Target Compounds						
20) 4,6-Dinitro-2-methylph...	15.596	198	175642	91.868	ng	# 73
24) Pentachlorophenol	16.867	266	730395	249.502	ng	99

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

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