Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN033022\
Data File : BN019243.D
Acq On : 30 Mar 2022 16:45
Operator : CG/JU
Sample : N1870-08DL 100X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
PT-ACIDS-WPDL

Quant Time: Mar 30 17:17:06 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	3511	0.400	ng	0.00
7) Naphthalene-d8	10.637	136	8962	0.400	ng	# 0.00
13) Acenaphthene-d10	14.474	164	4786	0.400	ng	0.00
19) Phenanthrene-d10	17.215	188	9441	0.400	ng	0.00
29) Chrysene-d12	21.413	240	6205	0.400	ng	# 0.00
35) Perylene-d12	23.767	264	4666	0.400	ng	# 0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.384	112	12564	1.481	ng	0.00
5) Phenol-d6	6.995	99	12354	1.341	ng	0.00
8) Nitrobenzene-d5	8.993	82	5911	0.866	ng	0.00
11) 2-Methylnaphthalene-d10	12.238	152	59	0.004	ng	0.00
14) 2,4,6-Tribromophenol	15.964	330	3025	1.266	ng	0.00
15) 2-Fluorobiphenyl	13.095	172	20628	1.008	ng	-0.01
27) Fluoranthene-d10	19.261	212	140	0.005	ng	0.01
31) Terphenyl-d14	19.843	244	17647	1.269	ng	0.00
Target Compounds						Qvalue
20) 4,6-Dinitro-2-methylph...	15.618	198	470	0.388	ng	# 63
24) Pentachlorophenol	16.877	266	3513	1.226	ng	99

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

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