Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN040921\
Data File : BN014256.D
Acq On : 09 Apr 2021 18:49
Operator : CG/JU
Sample : M1557-08DL 100X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : PT-ACIDS-WPDL

Quant Time: Apr 09 19:20:36 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN040921.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Apr 09 15:06:44 2021
Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.716	152	5792	0.40	ng	0.00
7) Naphthalene-d8	10.492	136	21429	0.40	ng	0.00
13) Acenaphthene-d10	14.346	164	11710	0.40	ng	0.00
19) Phenanthrene-d10	17.092	188	25733	0.40	ng	0.00
29) Chrysene-d12	21.282	240	23378	0.40	ng	# 0.00
36) Perylene-d12	23.516	264	22807	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.316	112	21509	1.56	ng	0.00
5) Phenol-d6	6.879	99	25806	1.50	ng	0.00
8) Nitrobenzene-d5	8.857	82	14724	1.08	ng	0.00
11) 2-Methylnaphthalene-d10	0.000	152	0d	0.00	ng	
14) 2,4,6-Tribromophenol	15.830	330	6178	1.35	ng	0.00
15) 2-Fluorobiphenyl	12.963	172	51546	1.20	ng	0.00
27) Fluoranthene-d10	0.000	212	0d	0.00	ng	
31) Terphenyl-d14	19.729	244	69004	1.34	ng	0.00
Target Compounds						Qvalue
20) 4,6-Dinitro-2-methylph...	15.462	198	4379	1.44	ng	# 57
24) Pentachlorophenol	16.739	266	6719	1.31	ng	99

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

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