Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN011823\
Data File : BN023707.D
Acq On : 18 Jan 2023 17:31
Operator : CG/JU
Sample : N3980-07
Misc : LOD-MDL-WATER 0.2ng
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
LOD-MDL-WATER-01-QT3-2022

Quant Time: Jan 19 07:53:45 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\8270-SIM-BN011823.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jan 19 07:52:57 2023
Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.992	152	4818	0.400	ng	0.00
7) Naphthalene-d8	10.818	136	12545	0.400	ng	# 0.01
13) Acenaphthene-d10	14.645	164	6300	0.400	ng	0.00
19) Phenanthrene-d10	17.402	188	13683	0.400	ng	# 0.01
29) Chrysene-d12	21.589	240	10712	0.400	ng	0.00
35) Perylene-d12	24.065	264	8246	0.400	ng	0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.529	112	2969	0.293	ng	0.00
5) Phenol-d6	7.154	99	3525	0.293	ng	0.00
8) Nitrobenzene-d5	9.164	82	2534	0.255	ng	0.01
11) 2-Methylnaphthalene-d10	12.405	152	8591	0.495	ng	0.00
14) 2,4,6-Tribromophenol	16.148	330	648	0.242	ng	0.01
15) 2-Fluorobiphenyl	13.276	172	6946	0.259	ng	0.01
27) Fluoranthene-d10	19.431	212	13277	0.414	ng	0.00
31) Terphenyl-d14	20.022	244	4856	0.287	ng	0.00
Target Compounds						Qvalue

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

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