

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061822\
 Data File : VN072930.D
 Acq On : 18 Jun 2022 11:17
 Operator : JC\MD
 Sample : N3334-12
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 WC-14A-1A-8-GR

Quant Time: Jun 20 08:02:51 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N061722W.M
 Quant Title : SW846 8260
 QLast Update : Sat Jun 18 06:22:45 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.016	168	299169	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.898	114	527318	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.680	117	462083	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.610	152	207002	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.369	65	180599	44.651	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	89.300%
35) Dibromofluoromethane	7.951	113	144465	44.324	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	88.640%
50) Toluene-d8	10.375	98	519212	42.811	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	85.620%#
62) 4-Bromofluorobenzene	12.663	95	188344	43.188	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	86.380%
Target Compounds						
					Qvalue	
16) Acetone	4.228	43	117418	52.812	ug/l	92
20) Methylene Chloride	5.016	84	79291	22.861	ug/l	90
25) 2-Butanone	7.269	43	52137	14.778	ug/l #	87
43) Isopropyl Acetate	8.492	43	40978	4.064	ug/l #	88
68) m/p-Xylenes	11.886	106	6396	1.027	ug/l	83
69) o-Xylene	12.216	106	10080	1.619	ug/l	96
84) 1,2,4-Trimethylbenzene	13.304	105	26063	1.989	ug/l	92

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

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