

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN082322\
 Data File : VN074070.D
 Acq On : 23 Aug 2022 16:01
 Operator : JC\MD
 Sample : N4291-01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 Y-7

Quant Time: Aug 24 01:21:34 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\624N072722W.M
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS
 QLast Update : Wed Jul 27 18:07:26 2022
 Response via : Initial Calibration

Manual Integrations
 APPROVED

Reviewed By :Krupa Patel 08/24/2022
 Supervised By :Mahesh Dadoda 08/24/2022

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

Internal Standards						
1) Bromochloromethane	7.593	128	69984	30.000	ug/l	0.00
28) 1,4-Difluorobenzene	8.904	114	361534	30.000	ug/l	0.00
57) Chlorobenzene-d5	11.686	117	319229	30.000	ug/l	0.00
System Monitoring Compounds						
27) 1,2-Dichloroethane-d4	8.369	65	183902	30.392	ug/l	0.00
Spiked Amount 30.000	Range 91 - 110		Recovery = 101.300%			
60) 4-Bromofluorobenzene	12.675	95	148595	26.289	ug/l	0.00
Spiked Amount 30.000	Range 63 - 112		Recovery = 87.633%			
63) Toluene-d8	10.381	98	508868	28.310	ug/l	0.00
Spiked Amount 30.000	Range 91 - 112		Recovery = 94.367%			
Target Compounds						
15) Acetone	4.234	58	2641m	2.764	ug/l	Qvalue
18) Methylene Chloride	5.016	84	1405	0.253	ug/l	92
25) Chloroform	7.745	83	324677	33.529	ug/l	100
37) Trichloroethene	9.151	130	1219	0.227	ug/l	90
41) Bromodichloromethane	9.698	83	13943	1.903	ug/l	91
61) Tetrachloroethene	10.916	164	4495	0.919	ug/l #	77

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

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