

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN100824\
 Data File : VN084349.D
 Acq On : 08 Oct 2024 15:17
 Operator : JC\MD
 Sample : P4334-02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 002-35TH-AVE(SEP)

Quant Time: Oct 09 02:31:09 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\624N100824W.M
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS
 QLast Update : Wed Oct 09 02:21:08 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Bromochloromethane	7.812	128	30781	30.000	ug/l	0.00
28) 1,4-Difluorobenzene	9.100	114	158595	30.000	ug/l	0.00
57) Chlorobenzene-d5	11.865	117	149399	30.000	ug/l	0.00
System Monitoring Compounds						
27) 1,2-Dichloroethane-d4	8.577	65	77489	31.269	ug/l	0.00
Spiked Amount	30.000	Range	91 - 110	Recovery	=	104.233%
60) 4-Bromofluorobenzene	12.847	95	65501	26.565	ug/l	0.00
Spiked Amount	30.000	Range	63 - 112	Recovery	=	88.533%
63) Toluene-d8	10.565	98	201772	29.153	ug/l	0.00
Spiked Amount	30.000	Range	91 - 112	Recovery	=	97.167%
Target Compounds						
					Qvalue	
15) Acetone	4.424	58	280218	909.997	ug/l	95
25) Chloroform	7.965	83	4083	1.052	ug/l	99
30) 2-Butanone	7.488	43	33526	24.187	ug/l #	95
58) 4-Methyl-2-Pentanone	10.441	43	11054	3.972	ug/l #	68
62) Toluene	10.629	91	355507	40.647	ug/l	100
82) p-Isopropyltoluene	13.729	119	7489	1.014	ug/l	97
91) Naphthalene	15.641	128	5304	0.780	ug/l #	94

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

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