

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN100824\
 Data File : VN084351.D
 Acq On : 08 Oct 2024 16:05
 Operator : JC\MD
 Sample : P4337-02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 16 Sample Multiplier: 1

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

Quant Time: Oct 09 02:32:00 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	7.818	128	30720	30.000	ug/l	0.00
28) 1,4-Difluorobenzene	9.100	114	157276	30.000	ug/l	0.00
57) Chlorobenzene-d5	11.865	117	146077	30.000	ug/l	0.00
System Monitoring Compounds						
27) 1,2-Dichloroethane-d4	8.577	65	75843	30.665	ug/l	0.00
Spiked Amount	30.000	Range	91 - 110	Recovery	=	102.233%
60) 4-Bromofluorobenzene	12.847	95	64761	26.862	ug/l	0.00
Spiked Amount	30.000	Range	63 - 112	Recovery	=	89.533%
63) Toluene-d8	10.565	98	197805	29.230	ug/l	0.00
Spiked Amount	30.000	Range	91 - 112	Recovery	=	97.433%
Target Compounds						
					Qvalue	
15) Acetone	4.424	58	295470	961.432	ug/l	93
25) Chloroform	7.965	83	4337	1.119	ug/l	91
30) 2-Butanone	7.483	43	35142	25.565	ug/l	98
58) 4-Methyl-2-Pentanone	10.441	43	11940	4.388	ug/l	94
62) Toluene	10.629	91	419130	49.012	ug/l	99
82) p-Isopropyltoluene	13.729	119	8370	1.159	ug/l	92
91) Naphthalene	15.641	128	6334	0.952	ug/l #	96

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

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