

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN102324\
 Data File : VN084473.D
 Acq On : 23 Oct 2024 15:26
 Operator : JC\MD
 Sample : P4464-01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 001-WILLETS-PT-BLVD(SEP)

Quant Time: Oct 23 15:48:04 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\624N100824W.M
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS
 QLast Update : Wed Oct 09 02:22:28 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Bromochloromethane	7.812	128	30612	30.000	ug/l	0.00
28) 1,4-Difluorobenzene	9.100	114	143930	30.000	ug/l	0.00
57) Chlorobenzene-d5	11.865	117	140820	30.000	ug/l	0.00
System Monitoring Compounds						
27) 1,2-Dichloroethane-d4	8.577	65	74713	30.315	ug/l	0.00
Spiked Amount	30.000	Range	91 - 110	Recovery	=	101.067%
60) 4-Bromofluorobenzene	12.847	95	59657	25.669	ug/l	0.00
Spiked Amount	30.000	Range	63 - 112	Recovery	=	85.567%
63) Toluene-d8	10.565	98	189009	28.973	ug/l	0.00
Spiked Amount	30.000	Range	91 - 112	Recovery	=	96.567%
Target Compounds						
					Qvalue	
15) Acetone	4.424	58	195025	636.832	ug/l	93
25) Chloroform	7.965	83	7599	1.968	ug/l	84
30) 2-Butanone	7.483	43	40993	32.587	ug/l	99
58) 4-Methyl-2-Pentanone	10.441	43	16207	6.178	ug/l	95
62) Toluene	10.629	91	911396	110.554	ug/l	98
67) m/p-Xylenes	12.070	106	1380	0.406	ug/l	98
82) p-Isopropyltoluene	13.729	119	9378	1.348	ug/l	98
91) Naphthalene	15.641	128	4984	0.777	ug/l	96

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

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