

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX080824\
 Data File : VX042945.D
 Acq On : 08 Aug 2024 11:32
 Operator : JC/MD
 Sample : P3505-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 001-WILLETS-PT-BLVD(JUNE)

Quant Time: Aug 08 14:51:45 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\624X080624W.M
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS
 QLast Update : Wed Aug 07 04:53:59 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Bromochloromethane	4.891	128	42291	30.000	ug/l	0.00
28) 1,4-Difluorobenzene	6.757	114	240439	30.000	ug/l	0.00
57) Chlorobenzene-d5	10.055	117	217465	30.000	ug/l	0.00
System Monitoring Compounds						
27) 1,2-Dichloroethane-d4	5.952	65	107533	27.536	ug/l	0.00
Spiked Amount	30.000	Range	91 - 110	Recovery	=	91.800%
60) 4-Bromofluorobenzene	11.079	95	97218	27.402	ug/l	0.00
Spiked Amount	30.000	Range	63 - 112	Recovery	=	91.333%
63) Toluene-d8	8.647	98	293296	29.858	ug/l	0.00
Spiked Amount	30.000	Range	91 - 112	Recovery	=	99.533%
Target Compounds						
					Qvalue	
3) Chloromethane	1.295	50	1378	0.604	ug/l	93
15) Acetone	2.386	58	41228	75.617	ug/l	94
18) Methylene Chloride	2.782	84	1695	0.577	ug/l #	86
25) Chloroform	5.087	83	170886	31.046	ug/l	97
30) 2-Butanone	4.581	43	15141	6.636	ug/l #	77
41) Bromodichloromethane	7.824	83	9635	2.545	ug/l #	96
62) Toluene	8.714	91	898112	84.321	ug/l	100
82) p-Isopropyltoluene	12.012	119	9011	0.869	ug/l	97

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

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