

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX090322\
 Data File : VX031114.D
 Acq On : 04 Sep 2022 04:21
 Operator : JC/MD
 Sample : N4438-03
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 KTS2-WC-S-MH-02-0-20

Quant Time: Sep 06 01:30:41 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X090122W.M
 Quant Title : SW846 8260
 QLast Update : Fri Sep 02 02:49:21 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.562	168	219190	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	357039	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	322085	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	143021	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	135838	48.164	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	96.320%
35) Dibromofluoromethane	5.391	113	131930	50.634	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	101.260%
50) Toluene-d8	8.653	98	466525	48.650	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	97.300%
62) 4-Bromofluorobenzene	11.079	95	163759	48.350	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	96.700%
Target Compounds						
					Qvalue	
16) Acetone	2.392	43	23345	21.442	ug/l	99
18) Methyl Acetate	2.709	43	3636	1.162	ug/l #	77
20) Methylene Chloride	2.794	84	18140	4.869	ug/l #	84
37) Ethyl Acetate	4.721	43	17251	5.440	ug/l #	94
43) Isopropyl Acetate	6.348	43	79301	16.001	ug/l #	89

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

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