

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX090322\
 Data File : VX031112.D
 Acq On : 04 Sep 2022 03:34
 Operator : JC/MD
 Sample : N4438-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 KTS2-WC-S-TR-01-02-03

Quant Time: Sep 06 01:29:57 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	220870	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	347922	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	324934	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	149253	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	137991	48.556	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	97.120%
35) Dibromofluoromethane	5.391	113	130543	51.415	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	102.820%
50) Toluene-d8	8.653	98	467364	50.015	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	100.020%
62) 4-Bromofluorobenzene	11.079	95	166305	50.389	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	100.780%
Target Compounds						
					Qvalue	
16) Acetone	2.386	43	23860	21.748	ug/l	95
20) Methylene Chloride	2.794	84	15329	3.714	ug/l #	83
37) Ethyl Acetate	4.721	43	18700	6.052	ug/l	96
43) Isopropyl Acetate	6.342	43	88909	18.410	ug/l #	90

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

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