

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX071023\
 Data File : VX036493.D
 Acq On : 10 Jul 2023 19:17
 Operator : JC/MD
 Sample : 03555-10
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 ETGI-289

Quant Time: Jul 11 01:33:25 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X062323W.M
 Quant Title : SW846 8260
 QLast Update : Fri Jun 23 15:03:28 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.550	168	191209	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	328969	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	303333	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	134213	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	125996	45.970	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	91.940%
35) Dibromofluoromethane	5.385	113	103173	47.564	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	95.120%
50) Toluene-d8	8.647	98	405573	50.060	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	100.120%
62) 4-Bromofluorobenzene	11.079	95	140888	49.370	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	98.740%
Target Compounds						
					Qvalue	
16) Acetone	2.380	43	69262	59.122	ug/l	93
20) Methylene Chloride	2.788	84	9122	3.616	ug/l #	82
25) 2-Butanone	4.562	43	50388	29.935	ug/l	99
37) Ethyl Acetate	4.721	43	13413	4.070	ug/l #	91
43) Isopropyl Acetate	6.342	43	121945	23.196	ug/l	96
51) 4-Methyl-2-Pentanone	8.574	43	209793	61.940	ug/l #	83

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

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