

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX071023\
 Data File : VX036491.D
 Acq On : 10 Jul 2023 18:30
 Operator : JC/MD
 Sample : 03555-06
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 ETGI-369

Quant Time: Jul 11 01:32:42 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	177304	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	302788	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	286272	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	129345	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	117280	46.145	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	92.300%
35) Dibromofluoromethane	5.385	113	95582	47.875	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	95.740%
50) Toluene-d8	8.647	98	370826	49.729	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	99.460%
62) 4-Bromofluorobenzene	11.079	95	133135	50.688	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	101.380%
Target Compounds						
					Qvalue	
16) Acetone	2.380	43	66303	61.034	ug/l	90
20) Methylene Chloride	2.788	84	7217	3.086	ug/l	97
25) 2-Butanone	4.562	43	48951	31.362	ug/l	93
37) Ethyl Acetate	4.708	43	11642	3.838	ug/l	98
43) Isopropyl Acetate	6.336	43	122341	25.283	ug/l	98
51) 4-Methyl-2-Pentanone	8.574	43	219401	70.378	ug/l #	82

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

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