

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX090623\
 Data File : VX037559.D
 Acq On : 06 Sep 2023 17:03
 Operator : JC/MD
 Sample : 04287-05
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 EB01-090123

Quant Time: Sep 06 22:47:28 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X081723W.M
 Quant Title : SW846 8260
 QLast Update : Fri Aug 18 09:30:38 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.550	168	129404	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	215258	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	191244	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	78221	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	74394	39.877	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	79.760%
35) Dibromofluoromethane	5.385	113	65366	45.851	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	91.700%
50) Toluene-d8	8.647	98	250063	48.357	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	96.720%
62) 4-Bromofluorobenzene	11.079	95	85803	42.222	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	84.440%
Target Compounds						
					Qvalue	
16) Acetone	2.386	43	4853	6.649	ug/l	90
18) Methyl Acetate	2.709	43	929	0.338	ug/l	99
20) Methylene Chloride	2.788	84	3477	2.130	ug/l #	77
25) 2-Butanone	4.574	43	2238	2.037	ug/l	95
27) cis-1,2-Dichloroethene	4.507	96	551	0.309	ug/l	68
44) Trichloroethene	7.129	130	14261	8.658	ug/l	99
51) 4-Methyl-2-Pentanone	8.580	43	6627	3.323	ug/l	92

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

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