

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070523\
 Data File : VX036419.D
 Acq On : 05 Jul 2023 21:15
 Operator : JC/MD
 Sample : 03507-08
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 CP-3

Quant Time: Jul 06 04:58:52 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	5.556	168	175528	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	298753	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	270664	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	123294	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	113886	45.263	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery =	90.520%		
35) Dibromofluoromethane	5.391	113	93581	47.506	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	95.020%		
50) Toluene-d8	8.647	98	363402	49.392	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery =	98.780%		
62) 4-Bromofluorobenzene	11.079	95	128441	49.561	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	99.120%		
Target Compounds						
					Qvalue	
16) Acetone	2.404	43	241457	224.519	ug/l	95
20) Methylene Chloride	2.788	84	12083	5.218	ug/l	98
25) 2-Butanone	4.574	43	120726	78.130	ug/l	96
37) Ethyl Acetate	4.727	43	50971	17.031	ug/l	97
43) Isopropyl Acetate	6.342	43	183938	38.527	ug/l	98
51) 4-Methyl-2-Pentanone	8.574	43	490567	159.486	ug/l	87

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

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