

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070523\
 Data File : VX036418.D
 Acq On : 05 Jul 2023 20:51
 Operator : JC/MD
 Sample : 03507-05
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 CP-2

Quant Time: Jul 06 04:58:24 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	180633	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	302109	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	269093	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	126716	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	113630	43.885	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery =	87.780%		
35) Dibromofluoromethane	5.391	113	93005	46.689	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	93.380%		
50) Toluene-d8	8.653	98	361331	48.565	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery =	97.120%		
62) 4-Bromofluorobenzene	11.079	95	128979	49.216	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	98.440%		
Target Compounds						
					Qvalue	
16) Acetone	2.404	43	232409	209.998	ug/l	96
20) Methylene Chloride	2.788	84	9152	3.841	ug/l	93
25) 2-Butanone	4.580	43	121827	76.614	ug/l	96
37) Ethyl Acetate	4.720	43	54413	17.979	ug/l	99
43) Isopropyl Acetate	6.342	43	192781	39.930	ug/l	97
51) 4-Methyl-2-Pentanone	8.573	43	521979	167.813	ug/l #	87

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

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