

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070523\
 Data File : VX036416.D
 Acq On : 05 Jul 2023 20:04
 Operator : JC/MD
 Sample : 03492-05
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 TB-TP2

Quant Time: Jul 06 04:57:35 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	5.549	168	183966	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	311663	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	282989	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	126722	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	119370	45.267	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery =	90.540%		
35) Dibromofluoromethane	5.385	113	96535	46.975	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	93.960%		
50) Toluene-d8	8.653	98	371830	48.444	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery =	96.880%		
62) 4-Bromofluorobenzene	11.079	95	133086	49.226	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	98.460%		
Target Compounds						
					Qvalue	
16) Acetone	2.404	43	76685	68.035	ug/l	99
20) Methylene Chloride	2.794	84	6683	2.754	ug/l	86
25) 2-Butanone	4.580	43	39469	24.372	ug/l	98
37) Ethyl Acetate	4.720	43	23020	7.373	ug/l	97
43) Isopropyl Acetate	6.342	43	156304	31.382	ug/l	98
51) 4-Methyl-2-Pentanone	8.573	43	234444	73.062	ug/l #	76

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

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