

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX071023\
 Data File : VX036492.D
 Acq On : 10 Jul 2023 18:53
 Operator : JC/MD
 Sample : 03555-08
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 VNJ-203

Quant Time: Jul 11 01:33:03 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	170455	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	6.763	114	294579	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	274729	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	125464	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	113282	46.363	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery =	92.720%		
35) Dibromofluoromethane	5.385	113	92896	47.826	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	95.660%		
50) Toluene-d8	8.647	98	357966	49.342	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery =	98.680%		
62) 4-Bromofluorobenzene	11.079	95	129018	50.489	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	100.980%		
Target Compounds						
						Qvalue
16) Acetone	2.380	43	62595	59.936	ug/l	98
20) Methylene Chloride	2.788	84	7593	3.377	ug/l	92
25) 2-Butanone	4.556	43	45376	30.240	ug/l	97
37) Ethyl Acetate	4.721	43	10371	3.514	ug/l	97
43) Isopropyl Acetate	6.336	43	119608	25.407	ug/l	97
51) 4-Methyl-2-Pentanone	8.574	43	211009	69.572	ug/l #	81

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

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