

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070523\
 Data File : VX036409.D
 Acq On : 05 Jul 2023 17:18
 Operator : JC/MD
 Sample : PB153983ZHE#10
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PB153983ZHE#10

Quant Time: Jul 06 04:54:31 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	176849	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	6.763	114	297049	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	262297	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	121916	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	111906	44.144	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery =	88.280%		
35) Dibromofluoromethane	5.385	113	92047	46.995	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	93.980%		
50) Toluene-d8	8.653	98	352717	48.215	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery =	96.420%		
62) 4-Bromofluorobenzene	11.079	95	125474	48.694	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	97.380%		
Target Compounds						
						Qvalue
16) Acetone	2.404	43	219344	202.434	ug/l	94
18) Methyl Acetate	2.709	43	4993	1.121	ug/l	96
20) Methylene Chloride	2.794	84	7834	3.358	ug/l	91
25) 2-Butanone	4.574	43	117647	75.569	ug/l	93
37) Ethyl Acetate	4.721	43	53945	18.128	ug/l	98
43) Isopropyl Acetate	6.342	43	200301	42.195	ug/l	97
51) 4-Methyl-2-Pentanone	8.574	43	527045	172.328	ug/l #	87

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

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