

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070523\
 Data File : VX036406.D
 Acq On : 05 Jul 2023 16:07
 Operator : JC/MD
 Sample : PB153983ZHE#07
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PB153983ZHE#07

Quant Time: Jul 06 04:53:12 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	163227	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	273878	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	246188	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	115204	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	104404	44.622	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery =	89.240%		
35) Dibromofluoromethane	5.391	113	85713	47.463	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	94.920%		
50) Toluene-d8	8.652	98	329569	48.862	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery =	97.720%		
62) 4-Bromofluorobenzene	11.079	95	117659	49.524	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	99.040%		
Target Compounds						
					Qvalue	
16) Acetone	2.398	43	219924	219.908	ug/l	93
18) Methyl Acetate	2.708	43	4467	1.086	ug/l	94
20) Methylene Chloride	2.788	84	7117	3.305	ug/l	95
25) 2-Butanone	4.568	43	120567	83.907	ug/l	96
37) Ethyl Acetate	4.720	43	53614	19.541	ug/l	98
43) Isopropyl Acetate	6.342	43	201829	46.113	ug/l	99
51) 4-Methyl-2-Pentanone	8.573	43	526305	186.645	ug/l #	87

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

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