

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070523\
 Data File : VX036403.D
 Acq On : 05 Jul 2023 14:56
 Operator : JC/MD
 Sample : PB153983ZHE#04
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PB153983ZHE#04

Quant Time: Jul 06 04:51:53 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	5.550	168	166443	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	283310	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	253524	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	113746	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	108296	45.391	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	90.780%
35) Dibromofluoromethane	5.385	113	90233	48.303	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	96.600%
50) Toluene-d8	8.653	98	342132	49.036	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	98.080%
62) 4-Bromofluorobenzene	11.079	95	119047	48.440	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	96.880%
Target Compounds						
					Qvalue	
16) Acetone	2.385	43	103560	101.552	ug/l	93
18) Methyl Acetate	2.709	43	4424	1.055	ug/l	97
20) Methylene Chloride	2.788	84	7228	3.292	ug/l #	85
25) 2-Butanone	4.556	43	46692	31.867	ug/l	100
37) Ethyl Acetate	4.720	43	31137	10.971	ug/l	96
43) Isopropyl Acetate	6.342	43	178107	39.339	ug/l	98
51) 4-Methyl-2-Pentanone	8.573	43	309947	106.258	ug/l #	79

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

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