

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070523\
 Data File : VX036410.D
 Acq On : 05 Jul 2023 17:42
 Operator : JC/MD
 Sample : PB153983ZHE#11
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PB153983ZHE#11

Quant Time: Jul 06 04:55:00 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	5.555	168	172493	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	6.763	114	288034	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	254964	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	112598	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	109238	44.180	ug/l	0.00
Spiked Amount	50.000	Range 78 - 117	Recovery	=	88.360%	
35) Dibromofluoromethane	5.385	113	90723	47.769	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	=	95.540%	
50) Toluene-d8	8.652	98	349592	49.283	ug/l	0.00
Spiked Amount	50.000	Range 92 - 112	Recovery	=	98.560%	
62) 4-Bromofluorobenzene	11.079	95	119807	47.950	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	=	95.900%	
Target Compounds						
						Qvalue
16) Acetone	2.398	43	218449	206.699	ug/l	97
18) Methyl Acetate	2.715	43	4676	1.076	ug/l	96
20) Methylene Chloride	2.788	84	7024	3.087	ug/l	89
25) 2-Butanone	4.568	43	116976	77.035	ug/l	94
37) Ethyl Acetate	4.720	43	52842	18.313	ug/l	97
43) Isopropyl Acetate	6.342	43	198731	43.174	ug/l	98
51) 4-Methyl-2-Pentanone	8.573	43	526829	177.648	ug/l #	87

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

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