

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070523\
 Data File : VX036400.D
 Acq On : 05 Jul 2023 13:45
 Operator : JC/MD
 Sample : PB153983ZHE#01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PB153983ZHE#01

Quant Time: Jul 06 04:50:26 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	164246	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	276209	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	254006	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	116244	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	102763	43.648	ug/l	0.00
Spiked Amount	50.000	Range 78 - 117	Recovery	=	87.300%	
35) Dibromofluoromethane	5.391	113	87730	48.170	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	=	96.340%	
50) Toluene-d8	8.647	98	329223	48.399	ug/l	0.00
Spiked Amount	50.000	Range 92 - 112	Recovery	=	96.800%	
62) 4-Bromofluorobenzene	11.079	95	121823	50.844	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	=	101.680%	
Target Compounds						
					Qvalue	
16) Acetone	2.386	43	100429	99.799	ug/l	92
20) Methylene Chloride	2.788	84	7231	3.337	ug/l	90
25) 2-Butanone	4.562	43	43596	30.152	ug/l	95
37) Ethyl Acetate	4.721	43	31791	11.490	ug/l	95
43) Isopropyl Acetate	6.342	43	173051	39.205	ug/l	97
51) 4-Methyl-2-Pentanone	8.574	43	299785	105.416	ug/l #	80

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

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