

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX071023\
 Data File : VX036481.D
 Acq On : 10 Jul 2023 14:33
 Operator : JC/MD
 Sample : PB154051ZHE#03
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PB154051ZHE#03

Quant Time: Jul 11 01:29:05 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	179504	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	309559	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	282943	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	129345	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	120612	46.875	ug/l	0.00
Spiked Amount	50.000	Range 78 - 117	Recovery	=	93.740%	
35) Dibromofluoromethane	5.385	113	98119	48.071	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	=	96.140%	
50) Toluene-d8	8.647	98	379611	49.794	ug/l	0.00
Spiked Amount	50.000	Range 92 - 112	Recovery	=	99.580%	
62) 4-Bromofluorobenzene	11.079	95	133393	49.675	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	=	99.360%	
Target Compounds						
					Qvalue	
16) Acetone	2.380	43	169826	154.415	ug/l	90
20) Methylene Chloride	2.788	84	11006	4.648	ug/l	98
25) 2-Butanone	4.556	43	99280	62.828	ug/l	98
37) Ethyl Acetate	4.715	43	40764	13.145	ug/l	98
43) Isopropyl Acetate	6.336	43	176695	35.718	ug/l	97
51) 4-Methyl-2-Pentanone	8.574	43	439542	137.909	ug/l #	86

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

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