

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX071023\
 Data File : VX036483.D
 Acq On : 10 Jul 2023 15:21
 Operator : JC/MD
 Sample : PB154051ZHE#05
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PB154051ZHE#05

Quant Time: Jul 11 01:29:49 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	176537	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	303922	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	276749	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	125498	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	117691	46.508	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery =	93.020%		
35) Dibromofluoromethane	5.385	113	94203	47.008	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	94.020%		
50) Toluene-d8	8.647	98	367963	49.161	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery =	98.320%		
62) 4-Bromofluorobenzene	11.079	95	130801	49.613	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	99.220%		
Target Compounds						
					Qvalue	
16) Acetone	2.380	43	173874	160.753	ug/l	93
20) Methylene Chloride	2.788	84	10157	4.361	ug/l	90
25) 2-Butanone	4.556	43	98760	63.549	ug/l	99
37) Ethyl Acetate	4.715	43	41088	13.495	ug/l	99
43) Isopropyl Acetate	6.342	43	178515	36.755	ug/l	98
51) 4-Methyl-2-Pentanone	8.574	43	437754	139.895	ug/l #	87

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

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