

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX071023\
 Data File : VX036484.D
 Acq On : 10 Jul 2023 15:44
 Operator : JC/MD
 Sample : PB154051ZHE#06
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PB154051ZHE#06

Quant Time: Jul 11 01:30:10 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	5.549	168	171001	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	299583	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	274569	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	125112	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	115113	46.962	ug/l	0.00
Spiked Amount	50.000	Range 78 - 117	Recovery	=	93.920%	
35) Dibromofluoromethane	5.385	113	93124	47.143	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	=	94.280%	
50) Toluene-d8	8.652	98	361692	49.023	ug/l	0.00
Spiked Amount	50.000	Range 92 - 112	Recovery	=	98.040%	
62) 4-Bromofluorobenzene	11.079	95	131563	50.625	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	=	101.240%	
Target Compounds						
					Qvalue	
16) Acetone	2.379	43	172006	164.174	ug/l	95
20) Methylene Chloride	2.788	84	10965	4.861	ug/l	93
25) 2-Butanone	4.562	43	97875	65.018	ug/l	97
37) Ethyl Acetate	4.714	43	40260	13.415	ug/l	99
43) Isopropyl Acetate	6.336	43	168902	35.279	ug/l	97
51) 4-Methyl-2-Pentanone	8.573	43	424608	137.660	ug/l #	87

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

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