

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX071023\
 Data File : VX036487.D
 Acq On : 10 Jul 2023 16:55
 Operator : JC/MD
 Sample : PB154051ZHE#01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PB154051ZHE#01

Quant Time: Jul 11 01:31:16 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	174386	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	302317	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	278453	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	130453	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	116188	46.481	ug/l	0.00
Spiked Amount	50.000	Range 78 - 117	Recovery	=	92.960%	
35) Dibromofluoromethane	5.385	113	94327	47.320	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	=	94.640%	
50) Toluene-d8	8.647	98	373019	50.101	ug/l	0.00
Spiked Amount	50.000	Range 92 - 112	Recovery	=	100.200%	
62) 4-Bromofluorobenzene	11.079	95	132552	50.544	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	=	101.080%	
Target Compounds						
					Qvalue	
16) Acetone	2.379	43	175039	163.826	ug/l	94
20) Methylene Chloride	2.788	84	8986	3.906	ug/l	86
25) 2-Butanone	4.556	43	99098	64.553	ug/l	93
37) Ethyl Acetate	4.714	43	39407	13.012	ug/l	97
43) Isopropyl Acetate	6.342	43	175829	36.394	ug/l	98
51) 4-Methyl-2-Pentanone	8.573	43	435912	140.046	ug/l #	86

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

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