

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX071023\
 Data File : VX036478.D
 Acq On : 10 Jul 2023 13:23
 Operator : JC/MD
 Sample : PB154051TB
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PB154051TB

Quant Time: Jul 11 01:27:19 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	178427	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	6.763	114	309249	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	284244	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	129358	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	117677	46.010	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	92.020%
35) Dibromofluoromethane	5.385	113	96351	47.252	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	94.500%
50) Toluene-d8	8.647	98	378479	49.695	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	99.400%
62) 4-Bromofluorobenzene	11.079	95	134718	50.219	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	100.440%
Target Compounds						
						Qvalue
16) Acetone	2.380	43	80825	73.934	ug/l	# 89
20) Methylene Chloride	2.794	84	7135	3.031	ug/l	# 89
25) 2-Butanone	4.556	43	56877	36.211	ug/l	97
37) Ethyl Acetate	4.721	43	15406	4.973	ug/l	96
43) Isopropyl Acetate	6.336	43	146410	29.625	ug/l	98
51) 4-Methyl-2-Pentanone	8.574	43	276884	86.961	ug/l	# 81

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

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