

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070523\
 Data File : VX036399.D
 Acq On : 05 Jul 2023 13:21
 Operator : JC/MD
 Sample : PB153983TB
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PB153983TB

Quant Time: Jul 06 04:49:53 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	175404	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	288777	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	266330	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	120247	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	110409	43.913	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	87.820%
35) Dibromofluoromethane	5.385	113	92640	48.652	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	97.300%
50) Toluene-d8	8.646	98	352186	49.521	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	99.040%
62) 4-Bromofluorobenzene	11.079	95	124884	49.853	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	99.700%
Target Compounds						
					Qvalue	
16) Acetone	2.385	43	96256	89.567	ug/l	92
20) Methylene Chloride	2.794	84	7493	3.238	ug/l #	84
25) 2-Butanone	4.568	43	43667	28.280	ug/l	93
37) Ethyl Acetate	4.720	43	32607	11.272	ug/l	98
43) Isopropyl Acetate	6.342	43	174144	37.735	ug/l	97
51) 4-Methyl-2-Pentanone	8.573	43	302407	101.710	ug/l #	80

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

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