

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070523\
 Data File : VX036404.D
 Acq On : 05 Jul 2023 15:19
 Operator : JC/MD
 Sample : PB153983ZHE#05
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PB153983ZHE#05

Quant Time: Jul 06 04:52:20 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	169874	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	287783	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	256014	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	114988	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	110308	45.301	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	90.600%
35) Dibromofluoromethane	5.385	113	90686	47.791	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	95.580%
50) Toluene-d8	8.653	98	345696	48.776	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	97.560%
62) 4-Bromofluorobenzene	11.079	95	120364	48.215	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	96.420%
Target Compounds						
					Qvalue	
16) Acetone	2.385	43	102922	98.887	ug/l	92
18) Methyl Acetate	2.709	43	4638	1.084	ug/l	96
20) Methylene Chloride	2.788	84	7133	3.183	ug/l	87
25) 2-Butanone	4.562	43	45557	30.464	ug/l	97
37) Ethyl Acetate	4.714	43	32225	11.178	ug/l	96
43) Isopropyl Acetate	6.342	43	178507	38.814	ug/l	97
51) 4-Methyl-2-Pentanone	8.573	43	303249	102.346	ug/l #	80

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

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