

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX032322\
 Data File : VX027671.D
 Acq On : 23 Mar 2022 16:37
 Operator : JC/MD
 Sample : N1908-10
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 WC-14A-1B

Quant Time: Mar 23 17:29:06 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X031722W.M
 Quant Title : SW846 8260
 QLast Update : Thu Mar 17 15:02:59 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Pentafluorobenzene	5.556	168	367959	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	598062	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	520643	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	249746	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	212911	46.970	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	93.940%
35) Dibromofluoromethane	5.391	113	193877	50.386	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	100.780%
50) Toluene-d8	8.653	98	711586	48.871	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	97.740%
62) 4-Bromofluorobenzene	11.079	95	264360	49.396	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	98.800%
Target Compounds						
					Qvalue	
16) Acetone	2.380	43	46117	24.960	ug/l	92
18) Methyl Acetate	2.703	43	23770	5.368	ug/l	99
20) Methylene Chloride	2.788	84	8368	2.106	ug/l	92
25) 2-Butanone	4.568	43	9896	3.560	ug/l	100
43) Isopropyl Acetate	6.348	43	25159	2.984	ug/l	98
69) o-Xylene	10.646	106	36217	5.168	ug/l	98
73) Isopropylbenzene	10.963	105	20969	1.153	ug/l	99
84) 1,2,4-Trimethylbenzene	11.756	105	76444	5.007	ug/l	99

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

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