

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX032322\
 Data File : VX027670.D
 Acq On : 23 Mar 2022 16:14
 Operator : JC/MD
 Sample : N1908-08
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 WC-14A-2

Quant Time: Mar 23 17:28:38 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	359325	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	582104	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	549812	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	255572	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	195898	44.255	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	88.520%
35) Dibromofluoromethane	5.391	113	186776	49.871	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	99.740%
50) Toluene-d8	8.653	98	683759	48.247	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	96.500%
62) 4-Bromofluorobenzene	11.079	95	282251	54.185	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	108.380%
Target Compounds						
					Qvalue	
16) Acetone	2.380	43	28030	15.535	ug/l	97
18) Methyl Acetate	2.709	43	17046	3.942	ug/l	98
20) Methylene Chloride	2.788	84	8847	2.281	ug/l	98
25) 2-Butanone	4.568	43	9522	3.508	ug/l	95
43) Isopropyl Acetate	6.342	43	25911	3.157	ug/l	98
69) o-Xylene	10.646	106	11246	1.520	ug/l	100
84) 1,2,4-Trimethylbenzene	11.756	105	29968	1.918	ug/l	98

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

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