

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX082923\
 Data File : VX037340.D
 Acq On : 29 Aug 2023 14:27
 Operator : JC/MD
 Sample : 04207-03
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 OW-06B-79-082823-H

Quant Time: Aug 30 02:06:47 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X081723W.M
 Quant Title : SW846 8260
 QLast Update : Fri Aug 18 09:30:38 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.550	168	140066	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	229358	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	199458	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	86727	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	84570	41.881	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	83.760%
35) Dibromofluoromethane	5.385	113	70293	46.276	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	92.560%
50) Toluene-d8	8.647	98	266083	48.292	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	96.580%
62) 4-Bromofluorobenzene	11.079	95	95040	43.892	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	87.780%
Target Compounds						
					Qvalue	
12) 1,1-Dichloroethene	2.319	96	15249	10.576	ug/l	83
16) Acetone	2.386	43	2408	3.048	ug/l	92
17) Carbon Disulfide	2.508	76	7877	2.148	ug/l	98
18) Methyl Acetate	2.709	43	1961	0.659	ug/l #	83
20) Methylene Chloride	2.788	84	6250	3.538	ug/l #	82
21) trans-1,2-Dichloroethene	3.093	96	3155	1.925	ug/l	86
24) 1,1-Dichloroethane	3.611	63	3744	1.304	ug/l	95
27) cis-1,2-Dichloroethene	4.489	96	32392	16.764	ug/l	81
30) Chloroform	5.099	83	1083	0.336	ug/l #	62
40) Benzene	6.037	78	1774	0.287	ug/l	91
42) 1,2-Dichloroethane	6.092	62	2295	0.946	ug/l	96
44) Trichloroethene	7.129	130	12099	6.894	ug/l	99

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

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