

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072122\
 Data File : VX030151.D
 Acq On : 20 Jul 2022 23:04
 Operator : JC/MD
 Sample : N3781-02 50X
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 ORA-0019

Quant Time: Jul 21 03:34:58 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X071222W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jul 13 09:05:23 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.562	168	181672	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	328638	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	289867	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	126220	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	117481	53.128	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 106.260%	
35) Dibromofluoromethane	5.391	113	101354	48.276	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 96.560%	
50) Toluene-d8	8.653	98	387454	49.070	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 98.140%	
62) 4-Bromofluorobenzene	11.086	95	130583	47.656	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 95.320%	
Target Compounds						
					Qvalue	
13) Acrolein	2.252	56	2120	18.651	ug/l #	79
16) Acetone	2.386	43	615468	590.679	ug/l	96
25) 2-Butanone	4.574	43	10607	5.976	ug/l #	89
51) 4-Methyl-2-Pentanone	8.580	43	13607	3.924	ug/l	97
52) Toluene	8.720	92	127299	23.487	ug/l	99
67) Ethyl Benzene	10.202	91	9336	0.954	ug/l	92
68) m/p-Xylenes	10.305	106	15213	3.965	ug/l	89
69) o-Xylene	10.647	106	6245	1.626	ug/l	99

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

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