

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX123121\
 Data File : VX026180.D
 Acq On : 30 Dec 2021 19:53
 Operator : JC/MD
 Sample : M5257-02
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 D3-12292021

Quant Time: Dec 31 07:11:30 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X123021W.M
 Quant Title : SW846 8260
 QLast Update : Thu Dec 30 08:02:21 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.562	168	135663	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	233799	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	211042	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	91952	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	102277	52.112	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 104.220%	
35) Dibromofluoromethane	5.391	113	79587	52.360	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 104.720%	
50) Toluene-d8	8.653	98	283360	50.118	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 100.240%	
62) 4-Bromofluorobenzene	11.085	95	97286	48.336	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 96.680%	
Target Compounds						
					Qvalue	
6) Chloroethane	1.690	64	36576	29.800	ug/l	98
16) Acetone	2.385	43	20996	19.744	ug/l	98
20) Methylene Chloride	2.794	84	5842	2.539	ug/l #	89
25) 2-Butanone	4.580	43	8342	5.632	ug/l	91
43) Isopropyl Acetate	6.348	43	14847	3.205	ug/l	95
67) Ethyl Benzene	10.195	91	15780	1.904	ug/l	99
68) m/p-Xylenes	10.305	106	23190	7.492	ug/l	90
69) o-Xylene	10.646	106	3455	1.141	ug/l	95
84) 1,2,4-Trimethylbenzene	11.756	105	11827	1.996	ug/l	94

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

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