

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX123121\
 Data File : VX026169.D
 Acq On : 30 Dec 2021 15:37
 Operator : JC/MD
 Sample : M5207-02
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 SUMP-122021

Quant Time: Dec 31 07:09:10 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	134112	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	230317	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	204134	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	88165	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	99507	51.287	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	102.580%
35) Dibromofluoromethane	5.391	113	77791	51.952	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	103.900%
50) Toluene-d8	8.653	98	271884	48.815	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	97.640%
62) 4-Bromofluorobenzene	11.085	95	93556	47.186	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	94.380%
Target Compounds						
					Qvalue	
16) Acetone	2.386	43	3577	3.403	ug/l	100
17) Carbon Disulfide	2.514	76	1206	0.283	ug/l #	76
27) cis-1,2-Dichloroethene	4.507	96	1444	0.783	ug/l	92
44) Trichloroethene	7.129	130	949	0.512	ug/l	87
64) Tetrachloroethene	9.275	164	61827	35.891	ug/l	96

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

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