

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX032921\
 Data File : VX021569.D
 Acq On : 29 Mar 2021 12:09
 Operator : JC/MD
 Sample : M1787-12
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 RE-122D1-20210319

Manual Integrations
 APPROVED

MMDadoda
 3/31/2021 9:41:44 AM

Quant Time: Mar 30 03:25:38 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X032421W.M
 Quant Title : SW846 8260
 QLast Update : Wed Mar 24 12:25:26 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.623	168	63470	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.829	114	114756	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.094	117	120005	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.059	152	53849	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.029	65	53601	50.76	ug/l	0.00
Spiked Amount	50.000	Range 78 - 117	Recovery	=	101.52%	
35) Dibromofluoromethane	5.464	113	40662	49.88	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	=	99.76%	
50) Toluene-d8	8.694	98	148823	51.23	ug/l	0.00
Spiked Amount	50.000	Range 92 - 112	Recovery	=	102.46%	
62) 4-Bromofluorobenzene	11.118	95	60223	53.33	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	=	106.66%	
Target Compounds						Qvalue
9) 1,1,2-Trichlorotrifluo...	2.364	101	2346	3.59	ug/l	96
12) 1,1-Dichloroethene	2.358	96	690	1.08	ug/l #	90
16) Acetone	2.423	43	1691	4.23	ug/l	99
21) trans-1,2-Dichloroethene	3.146	96	368	0.52	ug/l	97
27) cis-1,2-Dichloroethene	4.558	96	2110	2.59	ug/l	99
30) Chloroform	5.176	83	794	0.55	ug/l	99
38) Carbon Tetrachloride	5.758	117	697m	0.58	ug/l	
44) Trichloroethene	7.188	130	410479	487.16	ug/l	97
49) 1,4-Dioxane	7.723	88	309	15.33	ug/l #	60
55) 1,1,2-Trichloroethane	9.194	97	263	0.32	ug/l #	86
64) Tetrachloroethene	9.318	164	1466	1.82	ug/l	97

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

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