

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX091721\
 Data File : VX024316.D
 Acq On : 17 Sep 2021 17:55
 Operator : JC/MD
 Sample : M3782-04
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 RE132D4-20210914

Quant Time: Sep 20 01:59:27 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091421W.M
 Quant Title : SW846 8260
 QLast Update : Tue Sep 14 12:18:04 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.562	168	144816	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	246849	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	236826	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	102608	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.970	65	106963	51.797	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	103.600%
35) Dibromofluoromethane	5.397	113	84061	50.778	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	101.560%
50) Toluene-d8	8.653	98	304282	50.454	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	100.900%
62) 4-Bromofluorobenzene	11.085	95	112232	47.820	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	95.640%
Target Compounds						
						Qvalue
2) Dichlorodifluoromethane	1.172	85	915	0.657	ug/l	92
3) Chloromethane	1.300	50	550	0.378	ug/l	96
7) Trichlorofluoromethane	1.892	101	733	0.302	ug/l #	79
9) 1,1,2-Trichlorotrifluo...	2.331	101	10499	7.364	ug/l	94
12) 1,1-Dichloroethene	2.319	96	1977	1.481	ug/l	96
16) Acetone	2.392	43	2620	2.976	ug/l	91
27) cis-1,2-Dichloroethene	4.507	96	4362	2.584	ug/l	96
30) Chloroform	5.111	83	1210	0.407	ug/l #	79
38) Carbon Tetrachloride	5.696	117	1116	0.437	ug/l	98
44) Trichloroethene	7.135	130	467745	255.751	ug/l	96
49) 1,4-Dioxane	7.671	88	1149	26.126	ug/l	89
64) Tetrachloroethene	9.281	164	10094	5.502	ug/l	90

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

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