

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120221\
 Data File : VX025422.D
 Acq On : 02 Dec 2021 12:46
 Operator : JC/MD
 Sample : M4853-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 GWBR3D-BA-472-493-112921

Quant Time: Dec 03 03:05:43 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X111721W.M
 Quant Title : SW846 8260
 QLast Update : Wed Nov 17 15:36:16 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	110149	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	183529	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	165619	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	74610	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	63253	54.479	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 108.960%	
35) Dibromofluoromethane	5.391	113	56756	50.402	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 100.800%	
50) Toluene-d8	8.653	98	211426	49.937	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 99.880%	
62) 4-Bromofluorobenzene	11.085	95	74365	46.428	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 92.860%	
Target Compounds						
					Qvalue	
12) 1,1-Dichloroethene	2.319	96	10421	10.208	ug/l	89
16) Acetone	2.380	43	18795	29.133	ug/l	97
21) trans-1,2-Dichloroethene	3.099	96	1934	1.729	ug/l #	78
24) 1,1-Dichloroethane	3.611	63	6497	3.379	ug/l	98
25) 2-Butanone	4.568	43	6984	7.488	ug/l	96
27) cis-1,2-Dichloroethene	4.489	96	654033	500.770	ug/l	81
42) 1,2-Dichloroethane	6.098	62	2862	1.813	ug/l	97
44) Trichloroethene	7.129	130	125481	89.051	ug/l	94
52) Toluene	8.720	92	41799	13.199	ug/l	96

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

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