

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX113021\
 Data File : VX025392.D
 Acq On : 30 Nov 2021 13:03
 Operator : JC/MD
 Sample : M4826-03
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 GWBR3D-B-425-446-112321

Quant Time: Dec 01 02:20:53 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	5.549	168	128437	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.757	114	209427	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	187404	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	86084	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	71663	52.934	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 105.860%	
35) Dibromofluoromethane	5.379	113	64624	50.292	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 100.580%	
50) Toluene-d8	8.646	98	241072	49.898	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 99.800%	
62) 4-Bromofluorobenzene	11.079	95	84534	46.250	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 92.500%	
Target Compounds						
12) 1,1-Dichloroethene	2.318	96	3869	3.250	ug/l	96
16) Acetone	2.373	43	63762	84.760	ug/l	97
21) trans-1,2-Dichloroethene	3.093	96	1157	0.887	ug/l #	82
24) 1,1-Dichloroethane	3.611	63	2285	1.019	ug/l #	95
25) 2-Butanone	4.562	43	994	0.914	ug/l	97
27) cis-1,2-Dichloroethene	4.483	96	104981	68.935	ug/l	80
42) 1,2-Dichloroethane	6.092	62	1462	0.812	ug/l	97
44) Trichloroethene	7.122	130	63369	39.410	ug/l	92
52) Toluene	8.720	92	27679	7.660	ug/l	98
64) Tetrachloroethene	9.274	164	497	0.322	ug/l #	83

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

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