

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU111121\
 Data File : VU045745.D
 Acq On : 11 Nov 2021 17:46
 Operator : SY/MD
 Sample : M4606-08
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 MH-121.5T-20211109

Quant Time: Nov 12 06:11:27 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U110521W.M
 Quant Title : SW846 8260
 QLast Update : Fri Nov 05 13:35:16 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.375	168	116756	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.253	114	215855	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.420	117	226821	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.815	152	113326	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.703	65	97263	55.625	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 111.240%	
35) Dibromofluoromethane	5.292	113	76863	53.335	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 106.660%	
50) Toluene-d8	7.899	98	288098	53.761	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 107.520%	
62) 4-Bromofluorobenzene	10.635	95	108346	50.419	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 100.840%	
Target Compounds						
						Qvalue
16) Acetone	2.626	43	1201	1.349	ug/l	96
27) cis-1,2-Dichloroethene	4.671	96	38505	24.532	ug/l	88
44) Trichloroethene	6.546	130	26849	18.000	ug/l	96

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

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