

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092721\
 Data File : VU044912.D
 Acq On : 28 Sep 2021 01:31
 Operator : SY/MD
 Sample : M3859-03DL 5X
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 1029-MW-03ADL

Quant Time: Sep 28 06:30:55 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U092421W.M
 Quant Title : SW846 8260
 QLast Update : Sat Sep 25 09:30:46 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.375	168	199198	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.253	114	347297	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.420	117	319540	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.816	152	152312	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.707	65	156139	44.912	ug/l	0.00
Spiked Amount 50.000			Recovery	=	89.820%	
35) Dibromofluoromethane	5.292	113	107536	45.135	ug/l	0.00
Spiked Amount 50.000			Recovery	=	90.280%	
50) Toluene-d8	7.903	98	434422	48.279	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.560%	
62) 4-Bromofluorobenzene	10.636	95	160878	48.136	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.280%	
Target Compounds					Qvalue	
16) Acetone	2.633	43	3589	1.830	ug/l	95
21) trans-1,2-Dichloroethene	3.353	96	2646	1.030	ug/l	88
27) cis-1,2-Dichloroethene	4.668	96	42597	13.639	ug/l	88
44) Trichloroethene	6.546	130	39910	14.781	ug/l	97
64) Tetrachloroethene	8.555	164	113516	52.410	ug/l	96

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

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