

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092721\
 Data File : VU044921.D
 Acq On : 28 Sep 2021 05:01
 Operator : SY/MD
 Sample : M3915-18
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 RW4-20210922

Quant Time: Sep 28 06:32:19 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	193516	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.253	114	346145	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.420	117	313990	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.815	152	147451	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.703	65	155462	46.030	ug/l	0.00
Spiked Amount	50.000		Recovery	=	92.060%	
35) Dibromofluoromethane	5.292	113	105904	44.598	ug/l	0.00
Spiked Amount	50.000		Recovery	=	89.200%	
50) Toluene-d8	7.902	98	428888	47.822	ug/l	0.00
Spiked Amount	50.000		Recovery	=	95.640%	
62) 4-Bromofluorobenzene	10.635	95	155837	46.783	ug/l	0.00
Spiked Amount	50.000		Recovery	=	93.560%	
Target Compounds						
						Qvalue
9) 1,1,2-Trichlorotrifluo...	2.581	101	11235	4.772	ug/l	97
12) 1,1-Dichloroethene	2.578	96	3899	1.673	ug/l	88
16) Acetone	2.629	43	2559	1.343	ug/l	98
27) cis-1,2-Dichloroethene	4.671	96	3596	1.185	ug/l	92
30) Chloroform	5.089	83	1585	0.304	ug/l	94
44) Trichloroethene	6.546	130	1607276	597.248	ug/l	96
49) 1,4-Dioxane	6.983	88	1154	12.689	ug/l #	81
64) Tetrachloroethene	8.555	164	9002	4.230	ug/l	97

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

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