

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092721\
 Data File : VU044920.D
 Acq On : 28 Sep 2021 04:38
 Operator : SY/MD
 Sample : M3915-17
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 RE109D3-20210922

Quant Time: Sep 28 06:32:10 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	190810	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.253	114	335034	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.420	117	309901	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.815	152	145084	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.703	65	152031	45.653	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.300%	
35) Dibromofluoromethane	5.292	113	101690	44.244	ug/l	0.00
Spiked Amount 50.000			Recovery	=	88.480%	
50) Toluene-d8	7.899	98	418236	48.181	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.360%	
62) 4-Bromofluorobenzene	10.635	95	160905	49.907	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.820%	
Target Compounds					Qvalue	
9) 1,1,2-Trichlorotrifluo...	2.578	101	2802	1.207	ug/l	98
16) Acetone	2.636	43	2032	1.082	ug/l	100
27) cis-1,2-Dichloroethene	4.665	96	1287	0.430	ug/l	85
44) Trichloroethene	6.546	130	108571	41.682	ug/l	99
64) Tetrachloroethene	8.555	164	1720	0.819	ug/l #	82

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

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