

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092521\
 Data File : VU044866.D
 Acq On : 26 Sep 2021 03:30
 Operator : SY/MD
 Sample : M3859-04
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 1030-MW-11

Quant Time: Sep 27 03:15:08 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	5.379	168	182863	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.253	114	321388	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.420	117	301230	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.816	152	136939	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.707	65	153631	48.138	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.280%	
35) Dibromofluoromethane	5.295	113	103627	47.001	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.000%	
50) Toluene-d8	7.903	98	412462	49.533	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.060%	
62) 4-Bromofluorobenzene	10.636	95	143042	46.250	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.500%	
Target Compounds					Qvalue	
16) Acetone	2.639	43	16480	9.155	ug/l	97
25) 2-Butanone	4.723	43	6568	2.405	ug/l	98
27) cis-1,2-Dichloroethene	4.674	96	48706	16.988	ug/l	88
29) Tetrahydrofuran	5.070	42	1839	1.070	ug/l	87
44) Trichloroethene	6.546	130	57178	22.883	ug/l	98
64) Tetrachloroethene	8.559	164	785818	384.865	ug/l	98

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

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