

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092521\
 Data File : VU044849.D
 Acq On : 25 Sep 2021 20:05
 Operator : SY/MD
 Sample : M3859-03
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 1029-MW-03A

Quant Time: Sep 27 01:35:29 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	175911	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.253	114	305422	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.420	117	289265	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.816	152	129639	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.706	65	147815	48.146	ug/l	0.00
Spiked Amount	50.000		Recovery	=	96.300%	
35) Dibromofluoromethane	5.295	113	102371	48.859	ug/l	0.00
Spiked Amount	50.000		Recovery	=	97.720%	
50) Toluene-d8	7.903	98	397732	50.261	ug/l	0.00
Spiked Amount	50.000		Recovery	=	100.520%	
62) 4-Bromofluorobenzene	10.636	95	135674	46.161	ug/l	0.00
Spiked Amount	50.000		Recovery	=	92.320%	
Target Compounds						
					Qvalue	
16) Acetone	2.629	43	14287	8.250	ug/l	100
21) trans-1,2-Dichloroethene	3.359	96	11994	5.289	ug/l	97
27) cis-1,2-Dichloroethene	4.671	96	208073	75.441	ug/l	88
44) Trichloroethene	6.546	130	197148	83.026	ug/l	97
64) Tetrachloroethene	8.558	164	602815	307.449	ug/l	98

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

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