

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092821\
 Data File : VU044974.D
 Acq On : 29 Sep 2021 04:53
 Operator : SY/MD
 Sample : M3952-02
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 MW-1

Quant Time: Sep 29 05:51:44 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U092821W.M
 Quant Title : SW846 8260
 QLast Update : Tue Sep 28 12:32:13 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.379	168	189923	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.253	114	341056	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.420	117	333765	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.816	152	164119	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.707	65	158787	50.225	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery = 100.460%			
35) Dibromofluoromethane	5.295	113	108573	50.464	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 100.920%			
50) Toluene-d8	7.903	98	440686	51.840	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery = 103.680%			
62) 4-Bromofluorobenzene	10.636	95	165846	47.254	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 94.500%			
Target Compounds					Qvalue	
16) Acetone	2.633	43	2777	1.480	ug/l	98
27) cis-1,2-Dichloroethene	4.671	96	3855	1.325	ug/l	98
40) Benzene	5.780	78	6593	0.621	ug/l	95
52) Toluene	7.973	92	12496	1.880	ug/l	100
64) Tetrachloroethene	8.559	164	150486	69.081	ug/l	98
68) m/p-Xylenes	9.694	106	2258	0.451	ug/l	89

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

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