

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092821\
 Data File : VU044978.D
 Acq On : 29 Sep 2021 06:27
 Operator : SY/MD
 Sample : M3952-06
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 53 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 MW-4

Quant Time: Sep 29 07:59:16 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	192448	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.253	114	342491	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.420	117	332149	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.816	152	166410	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.707	65	160029	49.954	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	99.900%
35) Dibromofluoromethane	5.295	113	111361	51.543	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	103.080%
50) Toluene-d8	7.903	98	443677	51.973	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	103.940%
62) 4-Bromofluorobenzene	10.636	95	166668	47.290	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	94.580%
Target Compounds						
					Qvalue	
16) Acetone	2.633	43	2047	1.076	ug/l	91
27) cis-1,2-Dichloroethene	4.681	96	2791	0.947	ug/l	86
30) Chloroform	5.099	83	2701	0.541	ug/l	96
39) Methylcyclohexane	6.768	83	5789	1.317	ug/l	98
40) Benzene	5.781	78	11433	1.072	ug/l	100
44) Trichloroethene	6.549	130	678	0.270	ug/l	96
64) Tetrachloroethene	8.559	164	15057	6.946	ug/l	97
67) Ethyl Benzene	9.575	91	3619	0.282	ug/l	99
68) m/p-Xylenes	9.700	106	5733	1.152	ug/l	95
78) n-propylbenzene	10.909	91	3086	0.204	ug/l	90
80) 1,3,5-Trimethylbenzene	11.092	105	3540	0.326	ug/l	92
84) 1,2,4-Trimethylbenzene	11.472	105	9949	0.926	ug/l	97
95) Naphthalene	14.092	128	17433	1.420	ug/l	98

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

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