

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092721\
 Data File : VU044897.D
 Acq On : 27 Sep 2021 18:29
 Operator : SY/MD
 Sample : M3888-13
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 20898

Quant Time: Sep 27 18:53:10 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	201255	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.250	114	359761	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.423	117	349226	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.815	152	179821	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.703	65	161174	45.887	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.780%	
35) Dibromofluoromethane	5.292	113	111012	44.980	ug/l	0.00
Spiked Amount 50.000			Recovery	=	89.960%	
50) Toluene-d8	7.899	98	454547	48.765	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.540%	
62) 4-Bromofluorobenzene	10.639	95	192979	55.741	ug/l	0.00
Spiked Amount 50.000			Recovery	=	111.480%	
Target Compounds					Qvalue	
11) Tert butyl alcohol	3.253	59	91930	89.671	ug/l	99
16) Acetone	2.633	43	335363	169.267	ug/l	99
18) Methyl Acetate	2.951	43	121145	18.302	ug/l	98
25) 2-Butanone	4.713	43	105425	35.069	ug/l	100
40) Benzene	5.774	78	90869	7.808	ug/l	99
52) Toluene	7.973	92	60158	8.646	ug/l	100
67) Ethyl Benzene	9.574	91	13859	1.071	ug/l	97
68) m/p-Xylenes	9.697	106	29242	5.908	ug/l	97
69) o-Xylene	10.105	106	15356	3.209	ug/l	99
70) Styrene	10.121	104	31926	4.508	ug/l	100
84) 1,2,4-Trimethylbenzene	11.471	105	28120	2.638	ug/l	90
95) Naphthalene	14.089	128	504317	36.818	ug/l	100

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

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