

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092521\
 Data File : VU044877.D
 Acq On : 26 Sep 2021 07:48
 Operator : SY/MD
 Sample : M3888-13
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 52 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 20898

Quant Time: Sep 27 03:19:29 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	5.378	168	217142	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.253	114	380391	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.420	117	351757	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.815	152	170490	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.706	65	169253	44.661	ug/l	0.00
Spiked Amount 50.000			Recovery	=	89.320%	
35) Dibromofluoromethane	5.295	113	117666	45.091	ug/l	0.00
Spiked Amount 50.000			Recovery	=	90.180%	
50) Toluene-d8	7.902	98	482727	48.980	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.960%	
62) 4-Bromofluorobenzene	10.635	95	177911	48.601	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.200%	
Target Compounds					Qvalue	
11) Tert butyl alcohol	3.182	59	84381	76.286	ug/l	99
16) Acetone	2.626	43	396216	185.350	ug/l	99
18) Methyl Acetate	2.951	43	24994	3.500	ug/l	97
25) 2-Butanone	4.703	43	118927	36.666	ug/l	100
40) Benzene	5.777	78	96593	7.850	ug/l	98
52) Toluene	7.973	92	94719	12.875	ug/l	100
59) 2-Hexanone	8.687	43	5962	1.258	ug/l	100
67) Ethyl Benzene	9.574	91	23470	1.801	ug/l	97
68) m/p-Xylenes	9.693	106	50782	10.187	ug/l	95
69) o-Xylene	10.102	106	22669	4.703	ug/l	97
70) Styrene	10.118	104	31367	4.422	ug/l	97
78) n-propylbenzene	10.915	91	11793	0.804	ug/l #	68
80) 1,3,5-Trimethylbenzene	11.092	105	9461	0.923	ug/l	95
84) 1,2,4-Trimethylbenzene	11.471	105	47266	4.677	ug/l	90
95) Naphthalene	14.089	128	538477	41.345	ug/l	100

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

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