

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092521\
 Data File : VU044876.D
 Acq On : 26 Sep 2021 07:25
 Operator : SY/MD
 Sample : M3888-12
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 51 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 40459

Quant Time: Sep 27 03:19:05 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	217208	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.253	114	383068	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.420	117	355377	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.816	152	165920	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.707	65	174257	45.968	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.940%	
35) Dibromofluoromethane	5.295	113	119821	45.596	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.200%	
50) Toluene-d8	7.903	98	486480	49.016	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.040%	
62) 4-Bromofluorobenzene	10.636	95	175185	47.522	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.040%	
Target Compounds						
					Qvalue	
11) Tert butyl alcohol	3.199	59	11038	9.976	ug/l	96
16) Acetone	2.636	43	37641	17.603	ug/l	100
25) 2-Butanone	4.719	43	4845	1.493	ug/l #	85
29) Tetrahydrofuran	5.063	42	5014	2.456	ug/l	98
40) Benzene	5.780	78	7578	0.612	ug/l	99
52) Toluene	7.973	92	92873	12.536	ug/l	99
67) Ethyl Benzene	9.575	91	33742	2.562	ug/l	100
68) m/p-Xylenes	9.697	106	57759	11.468	ug/l	98
69) o-Xylene	10.105	106	23092	4.742	ug/l	98
78) n-propylbenzene	10.912	91	8256	0.579	ug/l	99
80) 1,3,5-Trimethylbenzene	11.092	105	12111	1.214	ug/l	97
84) 1,2,4-Trimethylbenzene	11.472	105	56226	5.717	ug/l	97
95) Naphthalene	14.089	128	77936	6.952	ug/l	99

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

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