

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU061521\
 Data File : VU044225.D
 Acq On : 15 Jun 2021 14:04
 Operator : SY/MD
 Sample : M2674-12 5X
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 TMR-DS-06-060921

Quant Time: Jun 15 14:23:05 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U060921W.M
 Quant Title : SW846 8260
 QLast Update : Thu Jun 10 05:11:37 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.385	168	192469	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.260	114	326611	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.423	117	322170	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.819	152	176343	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.713	65	136705	44.664	ug/l	0.00
Spiked Amount	50.000		Recovery	=	89.320%	
35) Dibromofluoromethane	5.298	113	107630	47.372	ug/l	0.00
Spiked Amount	50.000		Recovery	=	94.740%	
50) Toluene-d8	7.906	98	423663	48.554	ug/l	0.00
Spiked Amount	50.000		Recovery	=	97.100%	
62) 4-Bromofluorobenzene	10.639	95	169126	48.366	ug/l	0.00
Spiked Amount	50.000		Recovery	=	96.740%	
Target Compounds						
					Qvalue	
16) Acetone	2.649	43	23916	13.104	ug/l	93
18) Methyl Acetate	2.964	43	25072	6.368	ug/l	98
25) 2-Butanone	4.723	43	12959	5.073	ug/l	95
29) Tetrahydrofuran	5.076	42	2532	1.823	ug/l	99
51) 4-Methyl-2-Pentanone	7.803	43	4664	1.044	ug/l	93
52) Toluene	7.977	92	50699	7.943	ug/l	99
59) 2-Hexanone	8.694	43	14793	4.148	ug/l	89
67) Ethyl Benzene	9.578	91	41526	3.356	ug/l	99
68) m/p-Xylenes	9.700	106	40025	8.660	ug/l	99
69) o-Xylene	10.108	106	27702	6.321	ug/l	98
84) 1,2,4-Trimethylbenzene	11.475	105	36645	3.472	ug/l	97
95) Naphthalene	14.092	128	132414	12.913	ug/l	98

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

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