

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU061521\
 Data File : VU044222.D
 Acq On : 15 Jun 2021 12:55
 Operator : SY/MD
 Sample : M2674-09 5X
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 TMR-DS-03-060921

Quant Time: Jun 15 13:24:20 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	5.382	168	192528	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.256	114	325977	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.427	117	335361	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.819	152	177049	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.710	65	136174	44.477	ug/l	0.00
Spiked Amount 50.000			Recovery	=	88.960%	
35) Dibromofluoromethane	5.298	113	107098	47.230	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.460%	
50) Toluene-d8	7.906	98	421144	48.360	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.720%	
62) 4-Bromofluorobenzene	10.642	95	173479	49.708	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.420%	
Target Compounds					Qvalue	
16) Acetone	2.658	43	25889	14.358	ug/l	93
18) Methyl Acetate	2.964	43	11807	2.998	ug/l	99
20) Methylene Chloride	3.054	84	6892	1.331	ug/l	97
25) 2-Butanone	4.736	43	14729	5.764	ug/l	97
29) Tetrahydrofuran	5.083	42	3101	2.232	ug/l	95
40) Benzene	5.781	78	13782	1.282	ug/l	99
52) Toluene	7.977	92	88378	13.874	ug/l	100
59) 2-Hexanone	8.710	43	11376	3.196	ug/l	80
67) Ethyl Benzene	9.578	91	62077	4.819	ug/l	97
68) m/p-Xylenes	9.700	106	58078	12.072	ug/l	100
69) o-Xylene	10.108	106	36096	7.913	ug/l	98
80) 1,3,5-Trimethylbenzene	11.095	105	12920	1.225	ug/l	100
84) 1,2,4-Trimethylbenzene	11.475	105	42398	4.001	ug/l	99
95) Naphthalene	14.092	128	122386	12.112	ug/l	98

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

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