

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU041321\
 Data File : VU043053.D
 Acq On : 13 Apr 2021 15:50
 Operator : SY/MD
 Sample : M1983-06
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 31019

Quant Time: Apr 14 03:51:06 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U041221W.M
 Quant Title : SW846 8260
 QLast Update : Mon Apr 12 13:17:29 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.382	168	134454	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.256	114	224257	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.423	117	215189	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.819	152	112002	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.710	65	110277	48.79	ug/l	0.00
Spiked Amount	50.000		Recovery	=	97.58%	
35) Dibromofluoromethane	5.298	113	77193	45.81	ug/l	0.00
Spiked Amount	50.000		Recovery	=	91.62%	
50) Toluene-d8	7.906	98	281160	47.15	ug/l	0.00
Spiked Amount	50.000		Recovery	=	94.30%	
62) 4-Bromofluorobenzene	10.639	95	109649	45.92	ug/l	0.00
Spiked Amount	50.000		Recovery	=	91.84%	
Target Compounds						
						Qvalue
16) Acetone	2.639	43	319527	240.20	ug/l	100
18) Methyl Acetate	2.957	43	3377	1.20	ug/l #	88
25) 2-Butanone	4.716	43	98592	48.02	ug/l	100
29) Tetrahydrofuran	5.067	42	6324	4.89	ug/l #	80
51) 4-Methyl-2-Pentanone	7.800	43	10423	2.43	ug/l	93
52) Toluene	7.976	92	3929	0.93	ug/l	93
59) 2-Hexanone	8.693	43	16493	4.67	ug/l	95
69) o-Xylene	10.105	106	1126	0.37	ug/l	87
84) 1,2,4-Trimethylbenzene	11.475	105	7055	1.03	ug/l	77
95) Naphthalene	14.092	128	91932	13.47	ug/l #	92

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

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