

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU041421\
 Data File : VU043089.D
 Acq On : 14 Apr 2021 13:13
 Operator : SY/MD
 Sample : M1983-03ME
 Misc : 2.23g/5mL/100uL/5.00mL/MSVOA_U/MEOH
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 41019ME

Manual Integrations
 APPROVED

MMDadoda
 4/15/2021 5:36:51 PM

Quant Time: Apr 15 02:03:46 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	5.378	168	201848	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.256	114	324484	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.436	117	302654	50.00	ug/l	0.01
72) 1,4-Dichlorobenzene-d4	11.828	152	154243	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.713	65	157689	46.47	ug/l	0.00
Spiked Amount 50.000			Recovery	=	92.94%	
35) Dibromofluoromethane	5.314	113	104518	42.87	ug/l	0.02
Spiked Amount 50.000			Recovery	=	85.74%	
50) Toluene-d8	7.909	98	409925	47.51	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.02%	
62) 4-Bromofluorobenzene	10.655	95	152270	44.07	ug/l	0.02
Spiked Amount 50.000			Recovery	=	88.14%	
Target Compounds						Qvalue
16) Acetone	2.671	43	1130874	566.68	ug/l	98
25) 2-Butanone	4.758	43	49008m	15.90	ug/l	
39) Methylcyclohexane	6.761	83	25094	6.63	ug/l	97
51) 4-Methyl-2-Pentanone	7.877	43	119601m	19.26	ug/l	
52) Toluene	7.980	92	3583241	587.62	ug/l	100
59) 2-Hexanone	8.754	43	7552	1.48	ug/l	99
68) m/p-Xylenes	9.709	106	5658	1.31	ug/l	92
84) 1,2,4-Trimethylbenzene	11.481	105	18165	1.93	ug/l	87
86) p-Isopropyltoluene	11.806	119	106568	10.99	ug/l	98
95) Naphthalene	14.095	128	9455	3.54	ug/l #	83

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

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41019ME

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