

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU041421\
 Data File : VU043093.D
 Acq On : 14 Apr 2021 14:45
 Operator : SY/MD
 Sample : M1983-03MEDL 10X
 Misc : 2.23g/5mL/100uL/5.00mL/MSVOA_U/MEOH
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 41019MEDL

Manual Integrations
 APPROVED

MMDadoda
 4/15/2021 5:36:54 PM

Quant Time: Apr 15 02:05:12 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	185027	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.256	114	306862	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.423	117	288449	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.819	152	147476	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.710	65	147736	47.50	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.00%	
35) Dibromofluoromethane	5.298	113	103913	45.07	ug/l	0.00
Spiked Amount 50.000			Recovery	=	90.14%	
50) Toluene-d8	7.906	98	383231	46.97	ug/l	0.00
Spiked Amount 50.000			Recovery	=	93.94%	
62) 4-Bromofluorobenzene	10.639	95	146719	44.91	ug/l	0.00
Spiked Amount 50.000			Recovery	=	89.82%	
Target Compounds						Qvalue
16) Acetone	2.642	43	95408	51.88	ug/l	99
25) 2-Butanone	4.729	43	3316m	1.17	ug/l	
52) Toluene	7.976	92	332884	57.73	ug/l	99
86) p-Isopropyltoluene	11.799	119	10799	1.17	ug/l	99

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

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