

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU041421\
 Data File : VU043088.D
 Acq On : 14 Apr 2021 12:50
 Operator : SY/MD
 Sample : M1983-02ME
 Misc : 2.79g/5mL/100uL/5.00mL/MSVOA_U/MEOH
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 16873ME

Quant Time: Apr 15 02:03:22 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	175867	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.256	114	291084	50.00	ug/l	0.00
63) Chlorobenzene-d5	9.436	117	270459	50.00	ug/l	0.01
72) 1,4-Dichlorobenzene-d4	11.828	152	142612	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.716	65	142497	48.20	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.40%	
35) Dibromofluoromethane	5.311	113	88314	40.38	ug/l	0.01
Spiked Amount 50.000			Recovery	=	80.76%	
50) Toluene-d8	7.909	98	365131	47.18	ug/l	0.00
Spiked Amount 50.000			Recovery	=	94.36%	
62) 4-Bromofluorobenzene	10.655	95	133808	43.17	ug/l	0.02
Spiked Amount 50.000			Recovery	=	86.34%	
Target Compounds						
					Qvalue	
16) Acetone	2.671	43	332540	191.05	ug/l	99
39) Methylcyclohexane	6.761	83	103183	30.39	ug/l	96
52) Toluene	7.980	92	1538130	281.18	ug/l	100
67) Ethyl Benzene	9.587	91	10735	1.03	ug/l	99
68) m/p-Xylenes	9.713	106	23927	6.18	ug/l	98
69) o-Xylene	10.118	106	8167	2.16	ug/l	89
78) n-propylbenzene	10.922	91	13020	1.14	ug/l	98
80) 1,3,5-Trimethylbenzene	11.102	105	17837	2.11	ug/l	99
84) 1,2,4-Trimethylbenzene	11.481	105	51846	5.97	ug/l	100
86) p-Isopropyltoluene	11.806	119	84579	9.44	ug/l	99
95) Naphthalene	14.095	128	19953	4.57	ug/l #	96

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

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