

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN070822\
 Data File : VN073397.D
 Acq On : 08 Jul 2022 16:55
 Operator : JC\MD
 Sample : N3593-03ME
 Misc : 5.21g/5mL/100uL/5.00mL/MSVOA_N/MEOH
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 20220483C-1CME

Quant Time: Jul 09 05:28:03 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N062822W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jun 29 01:54:03 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.010	168	270178	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.898	114	446132	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.686	117	455279	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.610	152	245446	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.363	65	164821	42.258	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	84.520%
35) Dibromofluoromethane	7.945	113	138263	46.582	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	93.160%
50) Toluene-d8	10.374	98	481486	45.673	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	91.340%
62) 4-Bromofluorobenzene	12.674	95	248076	61.235	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	122.460%
Target Compounds						
					Qvalue	
16) Acetone	4.245	43	8800	4.020	ug/l #	84
17) Carbon Disulfide	4.398	76	25463	3.088	ug/l #	95
18) Methyl Acetate	4.769	43	39947	6.511	ug/l #	86

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

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