

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN071322\
 Data File : VN073450.D
 Acq On : 13 Jul 2022 15:25
 Operator : JC\MD
 Sample : N3619-01
 Misc : 1.22g/10mL/100uL/5.00mL/MSVOA_N/MEOH
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 31653

Quant Time: Jul 14 07:03:31 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	268337	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.898	114	444696	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.686	117	439525	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.615	152	238524	50.000	ug/l	0.00

System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.363	65	168839	43.586	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	87.180%
35) Dibromofluoromethane	7.939	113	137838	46.588	ug/l	-0.01
Spiked Amount	50.000	Range	75 - 124	Recovery	=	93.180%
50) Toluene-d8	10.374	98	499494	47.534	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	95.060%
62) 4-Bromofluorobenzene	12.668	95	231612	57.355	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	114.720%

Target Compounds					Qvalue	
18) Methyl Acetate	4.781	43	13539	2.222	ug/l	# 86
31) Cyclohexane	7.980	56	53181	8.383	ug/l	# 47
39) Methylcyclohexane	9.386	83	390115	68.831	ug/l	95
52) Toluene	10.439	92	54557	6.283	ug/l	99
69) o-Xylene	12.221	106	29400	4.228	ug/l	85
80) 1,3,5-Trimethylbenzene	12.998	105	185306	11.038	ug/l	96
84) 1,2,4-Trimethylbenzene	13.304	105	148982	8.888	ug/l	72
85) sec-Butylbenzene	13.439	105	42006	2.078	ug/l	78
86) p-Isopropyltoluene	13.557	119	56512	3.461	ug/l	96
89) n-Butylbenzene	13.880	91	62777	4.614	ug/l	# 72

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

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