

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN051920\
 Data File : VN061485.D
 Acq On : 19 May 2020 15:56
 Operator : JC/MD
 Sample : L2663-09
 Misc : 1.00µ/10mL/100uL/5.00mL/MSVOA_N/MEOH
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 1364

Quant Time: May 20 05:25:07 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N051820W.M
 Quant Title : SW846 8260
 QLast Update : Mon May 18 13:09:55 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.62	168	313719	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	555604	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	528878	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	289731	50.00	ug/l	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
33) 1,2-Dichloroethane-d4	7.99	65	196588	41.99	ug/l	0.00
Spiked Amount			Recovery	=	83.98%	
35) Dibromofluoromethane	7.54	113	143199	49.20	ug/l	0.00
Spiked Amount			Recovery	=	98.40%	
50) Toluene-d8	10.06	98	672041	49.78	ug/l	0.00
Spiked Amount			Recovery	=	99.56%	
62) 4-Bromofluorobenzene	12.38	95	271292	53.79	ug/l	0.00
Spiked Amount			Recovery	=	107.58%	

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
10) Methyl Iodide	3.89	142	2859	0.766	ug/l #	88
16) Acetone	3.78	43	45018	27.395	ug/l	98
29) Tetrahydrofuran	7.18	42	1738	1.231	ug/l #	74
51) 4-Methyl-2-Pentanone	9.95	43	15735	3.515	ug/l	96
52) Toluene	10.12	92	6514	0.648	ug/l	89
67) Ethyl Benzene	11.48	91	18610	0.945	ug/l #	86
68) m/p-Xylenes	11.59	106	37209	4.979	ug/l	100
69) o-Xylene	11.92	106	37409	5.173	ug/l	96
84) 1,2,4-Trimethylbenzene	13.01	105	29820	1.565	ug/l	96
85) sec-Butylbenzene	13.15	105	7762	0.352	ug/l	93

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

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