

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN111621\
 Data File : VN069511.D
 Acq On : 16 Nov 2021 16:24
 Operator : JC/MD
 Sample : M4619-01
 Misc : 5.04g/10mL/100uL/5.00mL/MSVOA_N/MEOH
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 HR-01-111021

Manual Integrations
 APPROVED

Reviewed By : John Carlone 11/17/2021
 Supervised By : Mahesh Dadoda 11/18/2021

Quant Time: Nov 17 01:49:49 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N110921W.M
 Quant Title : SW846 8260
 QLast Update : Tue Nov 09 18:18:22 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.078	168	1341367	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.963	114	2340106	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.746	117	2221353	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.675	152	1035320	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.432	65	984432	50.616	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery = 101.240%			
35) Dibromofluoromethane	8.016	113	650629	46.358	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery = 92.720%			
50) Toluene-d8	10.440	98	2929439	53.070	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery = 106.140%			
62) 4-Bromofluorobenzene	12.733	95	1060310	52.288	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery = 104.580%			
Target Compounds						
						Qvalue
18) Methyl Acetate	4.857	43	25361	0.743	ug/l	98
20) Methylene Chloride	5.068	84	10083m	0.678	ug/l	
59) 2-Hexanone	11.092	43	27014	9.064	ug/l	91
69) o-Xylene	12.286	106	8422	0.301	ug/l	98
80) 1,3,5-Trimethylbenzene	13.058	105	47119	0.707	ug/l	98
84) 1,2,4-Trimethylbenzene	13.372	105	19597	0.305	ug/l	91
85) sec-Butylbenzene	13.501	105	23668	0.290	ug/l	100
86) p-Isopropyltoluene	13.616	119	31560	0.487	ug/l	97

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

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