

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU053122\
 Data File : VU048769.D
 Acq On : 01 Jun 2022 04:10
 Operator : SY/MD
 Sample : N3099-01
 Misc : 5.24g/10.0mL/100uL/5.0mL/MSVOA_U/MEOH
 ALS Vial : 45 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 20220417-COMPOSITE-F

Manual Integrations
 APPROVED

Reviewed By : John Carlone 06/01/2022
 Supervised By : Mahesh Dadoda 06/02/2022

Quant Time: Jun 01 05:47:29 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U052622W.M
 Quant Title : SW846 8260
 QLast Update : Fri May 27 06:42:31 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.363	168	289666	50.000	ug/l	-0.01
34) 1,4-Difluorobenzene	6.244	114	489918	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.420	117	469326	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.819	152	231735	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.700	65	227103	52.205	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery = 104.400%			
35) Dibromofluoromethane	5.292	113	170132	48.676	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 97.360%			
50) Toluene-d8	7.896	98	635785	50.745	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery = 101.500%			
62) 4-Bromofluorobenzene	10.639	95	238812	50.400	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 100.800%			
Target Compounds						
					Qvalue	
12) 1,1-Dichloroethene	2.507	96	3030	1.028	ug/l #	78
16) Acetone	2.655	43	7848m	3.540	ug/l	
18) Methyl Acetate	2.964	43	28296	5.083	ug/l	95
24) 1,1-Dichloroethane	3.848	63	22630	3.421	ug/l	99
25) 2-Butanone	4.739	43	59744	18.758	ug/l	98
27) cis-1,2-Dichloroethene	4.652	96	21346	5.468	ug/l	87
29) Tetrahydrofuran	5.070	42	4894m	2.439	ug/l	
65) Chlorobenzene	9.452	112	10124	1.020	ug/l	96

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

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