

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU053122\
 Data File : VU048765.D
 Acq On : 01 Jun 2022 02:36
 Operator : SY/MD
 Sample : N3099-03 20X
 Misc : 3.17g/10.0mL/100uL/5.0mL/MSVOA_U/MEOH
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 20220426-AMENDED-COMPOSITE-F

Manual Integrations
 APPROVED

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

Quant Time: Jun 01 05:45:58 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	5.362	168	269323	50.000	ug/l	-0.01
34) 1,4-Difluorobenzene	6.243	114	468944	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.420	117	450673	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.816	152	219189	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.700	65	216065	53.419	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery = 106.840%			
35) Dibromofluoromethane	5.289	113	159058	47.543	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 95.080%			
50) Toluene-d8	7.896	98	581795	48.513	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery = 97.020%			
62) 4-Bromofluorobenzene	10.642	95	218355	48.144	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 96.280%			
Target Compounds						
16) Acetone	2.671	43	8680m	4.211	ug/l	Qvalue
18) Methyl Acetate	2.957	43	188104	36.346	ug/l	98
24) 1,1-Dichloroethane	3.848	63	7885	1.282	ug/l	98
25) 2-Butanone	4.735	43	71267	24.067	ug/l	98
27) cis-1,2-Dichloroethene	4.655	96	10723	2.955	ug/l	85
29) Tetrahydrofuran	5.083	42	5303m	2.843	ug/l	

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

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