

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU053122\
 Data File : VU048767.D
 Acq On : 01 Jun 2022 03:23
 Operator : SY/MD
 Sample : N3099-04 20X
 Misc : 3.73g/10.0mL/100uL/5.0mL/MSVOA_U/MEOH
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 20220429-AMENDED-COMPOSITE-I

Quant Time: Jun 01 05:46:43 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

Manual Integrations
 APPROVED

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

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

Internal Standards						
1) Pentafluorobenzene	5.362	168	301854	50.000	ug/l	-0.01
34) 1,4-Difluorobenzene	6.243	114	516261	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.420	117	496857	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.816	152	244565	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.700	65	227231	50.125	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery = 100.260%			
35) Dibromofluoromethane	5.292	113	168576	45.770	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 91.540%			
50) Toluene-d8	7.896	98	616632	46.705	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery = 93.420%			
62) 4-Bromofluorobenzene	10.639	95	234333	46.931	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 93.860%			
Target Compounds						
16) Acetone	2.658	43	9467	4.098	ug/l #	82
18) Methyl Acetate	2.954	43	200087	34.495	ug/l	98
25) 2-Butanone	4.735	43	69122	20.827	ug/l	98
29) Tetrahydrofuran	5.092	42	4909m	2.348	ug/l	

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

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