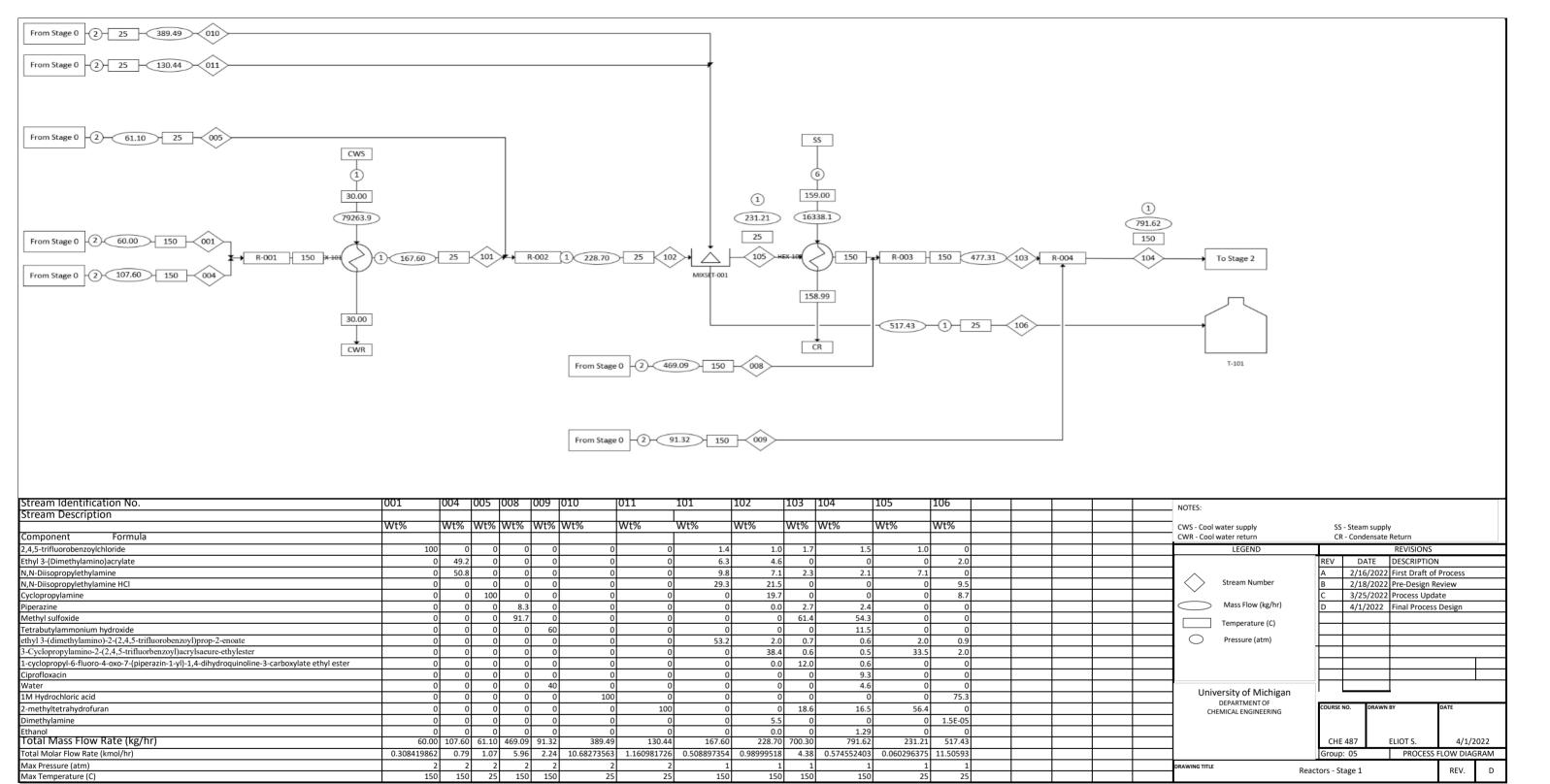
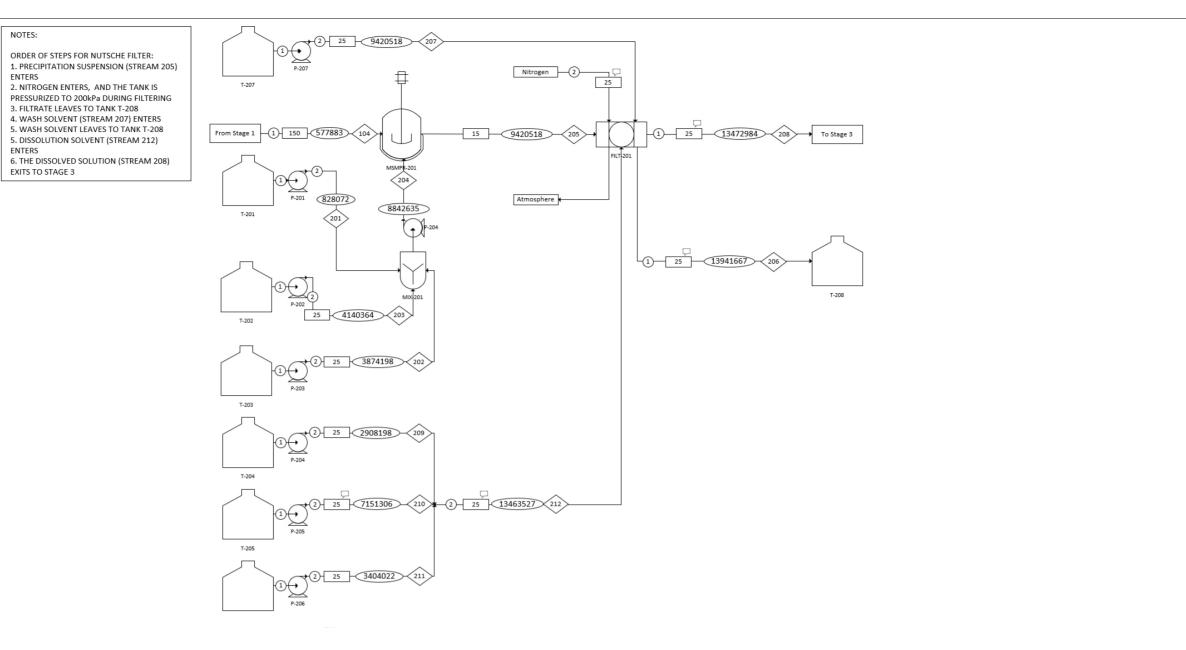


Stream Identification No.	001	002	003	004	005	006	007	800	009	010	011					NOTES:				
Stream Description																NOTES.				
	Wt%	Wt%	Wt%	Wt%	Wt%	Wt%	Wt%	Wt%	Wt%	Wt%	Wt%					CWS - Cool water supply	SS - Stea	m supply		
Component Formula																CWR - Cool water return	CR - Con	densate Ret	urn	
2,4,5-trifluorobenzoylchloride C7H2F3Cl	100) (0	0	0	0	0	0	())				LEGEND			REVISIONS	
Ethyl 3-(Dimethylamino)acrylate C7H13O2N	C	100	(49.2	. 0	0	0	0	0	()					REV	DATE	DESCRIPTIO	N
N,N-Diisopropylethylamine	C		100	50.8	0	0	0	0	0	()				Stream Number	Α	2/16/2022	First Draft of	f Process
Cyclopropylamine	C) ((0	100	0	0	0	0	(_	В	2/18/2022	Pre-Design F	Review
Piperazine	C) ((0	0	100	0	8.3	0	()				Mass Flow (kg/hr)	С	3/25/2022	Process Upd	late
Methyl sulfoxide	C) ((0	0	0	100	91.7	0	()					D	4/1/2022	Final Proces	s Design
Tetrabutylammonium hydroxide	C) ((0	0	0	0	0	60	()				Temperature (C)				
Water	C) ((0	0	0	0	0	40	()				Pressure (atm)				
1M HCl	C	0	(0	0	0	0	0	0	100))								
2-methyltetrahydrofuran	C	0	(0	0	0	0	0	0	(10)				University of Michigan	COURSE NO	DRAWN	I BY	DATE
Total Mass Flow Rate (kg/hr)	60.00	52.99	54.61	107.60	61.10	39.09	430.00	469.09	91.32	389.49	130.4	1				DEPARTMENT OF	CHE 4	187	ELIOT S.	3/25/2022
Total Molar Flow Rate (kmol/hr)	0.31	0.37	0.42	0.79	1.07	0.45	5.50	5.96	2.24	10.68274	1.16098	2				CHEMICAL ENGINEERING	Group:	05	PROCESS FL	OW DIAGRAM
Max Pressure (atm)	2	2 2	. 2	2	2	1	2	2	2	2		2				DRAWING TITLE Feed Storage and Preparation - Stage 0			REV. C	
Max Temperature (C)	150) 25	25	150	25	25	25	150	150	25	2	5				Teed Storage and Preparation - Stage 0				ILV.



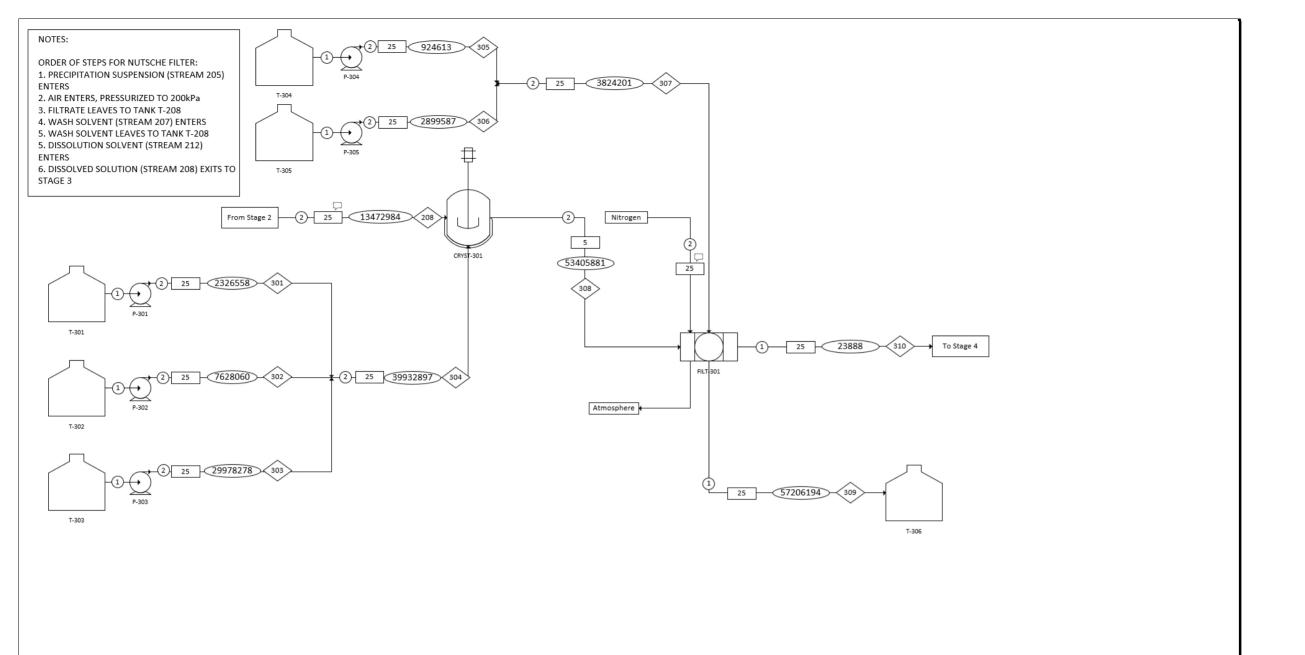
377722.6



Stream Identification No.	206	207	208	209	210	211	212		NOTES:					
Stream Description									1.0.125					
	Wt%		CWS - Cool water supply	SS - Stear	REVISIONS / DATE DESCRIPTION 2/16/2022 First Draft of Process 2/18/2022 Pre-Design Review 3/25/2022 Process Update 4/1/2022 Final Process Design									
Component Formula									CWR - Cool water return	REV DATE DESCRIPTION A 2/16/2022 First Draft of Process B 2/18/2022 Pre-Design Review C 3/25/2022 Process Update D 4/1/2022 Final Process Design COURSE NO. DRAWN BY DATE CHE 487 ELIOT S. 4/1/2022 Group: 05 PROCESS FLOW DIAGRAM				
Acetic Acid	6.0	0	0	0	0	0	0		LEGEND		F	REVISIONS		
Water	29.7	0	53.1	0	100	0	53.1		_	REV [DATE [DESCRIPTION		
Acetonitrile	27.8	0	0	0	0	0	0		Stream Number	A 2/1	L6/2022 F	irst Draft of F	Process	
Acetone	32.5	100	0	0	0	0	0			B 2/1	L8/2022 F	Pre-Design Re	view	
Formic Acid	0.0	0	21.6	100	0	0	21.6		Mass Flow (kg/hr)	C 3/2	25/2022 P	Process Updat	te	
Isopropanol	0.0	0	0	0	0	0	0			D 4/1	/2022 F	inal Process I	Design	
ciprofloxacin	0.008	0	0.2	0	0	0	0		Temperature (C)					
Tetrabutylammonium hydroxide	0.5	0	0	0	0	0	0		O Brossuro (atm)					
0.35 M Hydrochloric acid	0	0	25.3	0	0	100	25.3		Pressure (atm)					
2,4,5-trifluorobenzoylchloride	0.06	0	0.0	0	0	0	0.0							
N,N-Diisopropylethylamine	0.09	0	0.0	0	0	0	0.0							
Piperazine	0.10	0	0.0	0	0	0	0.0							
Methyl sulfoxide	2.3	0	0.0	0	0	0	0.0							
trifluorobenzoyl)prop-2-enoate	0	0	0.025	0	0	0	0.0							
trifluorbenzoyl)acrylsaeure-ethylester	0	0	0.02	0	0	0	0		University of Michigan		DRAWN BY	' ['	DATE	
1-cyclopropyl-6-fluoro-4-oxo-7-(piperazin-1-	0	0	0.02	0	0	0	0		DEPARTMENT OF					
2-methyltetrahydrofuran	0.7	0	0	0	0	0	0		CHEMICAL ENGINEERING					
Ethanol	0.05		0	0	0	0	0							
Total Mass Flow Rate (kg/batch)			13472984.38			3404022.07	13463527.05			CHE 487	El			
Total Molar Flow Rate (kmol/hr)	314.8163939	106.8579409	50.17983412	86.54862147	543.7561484	127.8947868	758.1995567			Group: 05		PROCESS FL	OW DIAGRAM	
Max Pressure (atm)									DRAWING TITLE Primary 6	ecovery - St	200 2		REV C	
Max Temperature ©	25	25	25	25	25	25	25		DEPARTMENT OF CHEMICAL ENGINEERING CHE 487 ELIOT S. 4/1/2022 Group: 05 PROCESS FLOW DIAGRAM DRAWING TITLE					

NOTES:

ENTERS



Stream Identification No.	309	310								NOTES:						
Stream Description										1						
•	Wt%	Wt%								CWS - Cool water supply SS - Steam supply						
Component Formula										CWR - Cool water return CR - Condensate Return						
Acetic Acid	0	0								LEGEND			REVISIONS			
Water	27.4	29.9									REV	DATE	DESCRIPTION	N		
Acetonitrile	0	0								Stream Number	Α	2/16/202	2 First Draft of	Process		
Acetone	5.1	0									В	2/18/202	2 Pre-Design R	eview		
Formic Acid	9.2	0								Mass Flow (kg/hr)						
Isopropanol	52.4	0									D	4/1/2022	Final Process	Design		
Ciprofloxacin HCl	0.0173	70.1								Temperature (C)						
Tetrabutylammonium hydroxide	0	0														
Hydrochloric Acid	5.950	0								Pressure (atm)						
trifluorobenzoyl)prop-2-enoate	0.006	0								University of Michigan	COURSE N	DRAWI	N BY	DATE		
trifluorbenzoyl)acrylsaeure-ethylester	0.005	0								DEPARTMENT OF						
1-cyclopropyl-6-fluoro-4-oxo-7-(piperazin-1-										CHEMICAL ENGINEERING						
Total Mass Flow Rate (kg/batch)	57206194.3									CHE 487 ELIOT				3/25/202		
Total Molar Flow Rate	1520.809476	0.605472									Group:	05	PROCESS FL	OW DIAGRA		
Max Pressure (atm) DRAWING TITLE							DRAWING TITLE Crystalli	Crystallization - Stage 3								
Max Temperature							l	I		Crystallization - Stage 5						

