



# Full wwPDB X-ray Structure Validation Report ⓘ

May 16, 2020 – 02:32 pm BST

PDB ID : 2FXR  
Title : human beta tryptase II complexed with activated ketone inhibitor CRA-29382  
Authors : Katz, B.A.  
Deposited on : 2006-02-06  
Resolution : 2.50 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.11

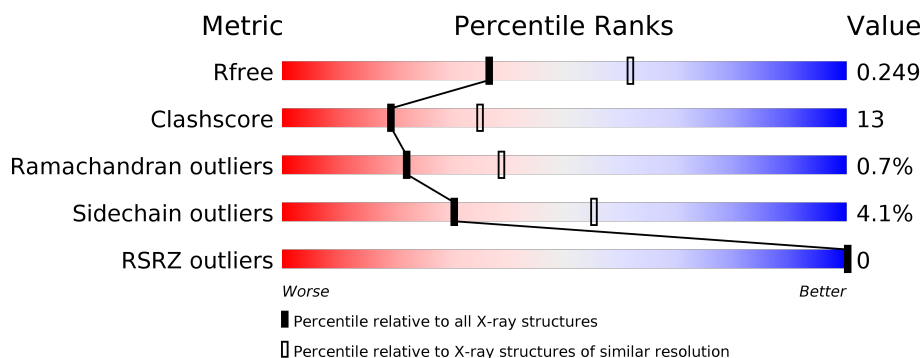
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*





The reported resolution of this entry is 2.50 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	245	
1	B	245	
1	C	245	
1	D	245	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	C3A	A	999	X	-	-	-
2	C3A	B	998	X	-	-	-
2	C3A	C	997	X	-	-	-
2	C3A	D	996	X	-	-	-

## 2 Entry composition [i](#)

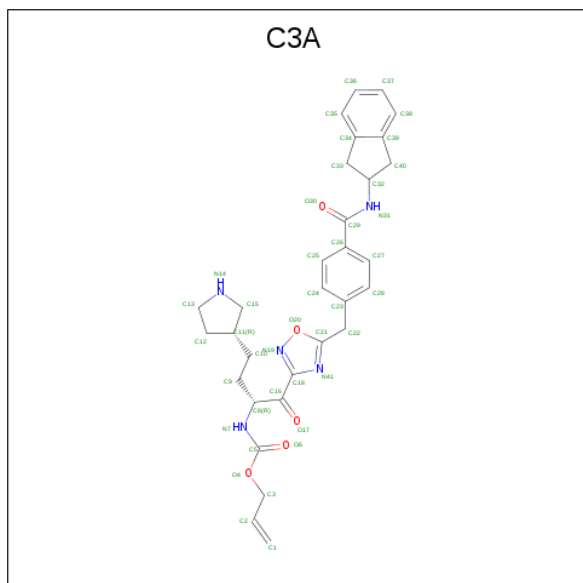
There are 3 unique types of molecules in this entry. The entry contains 7923 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Tryptase beta-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	243	Total	C	N	O	S	0	0	0
			1917	1229	338	338	12			
1	B	243	Total	C	N	O	S	0	0	0
			1920	1230	338	340	12			
1	C	243	Total	C	N	O	S	0	0	0
			1920	1230	338	340	12			
1	D	243	Total	C	N	O	S	0	0	0
			1920	1230	338	340	12			

- Molecule 2 is ALLYL {(1S)-1-[(5-{4-[(2,3-DIHYDRO-1H-INDEN-2-YLAMINO)CARBONYL]BENZYL}-1,2,4-OXADIAZOL-3-YL)CARBONYL]-3-PYRROLIDIN-3-YLPROPYL}CARBAMATE (three-letter code: C3A) (formula: C<sub>31</sub>H<sub>35</sub>N<sub>5</sub>O<sub>5</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			41	31	5	5		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	B	1	Total 41	C 31	N 5	O 5	0	0
2	C	1	Total 41	C 31	N 5	O 5	0	0
2	D	1	Total 41	C 31	N 5	O 5	0	0

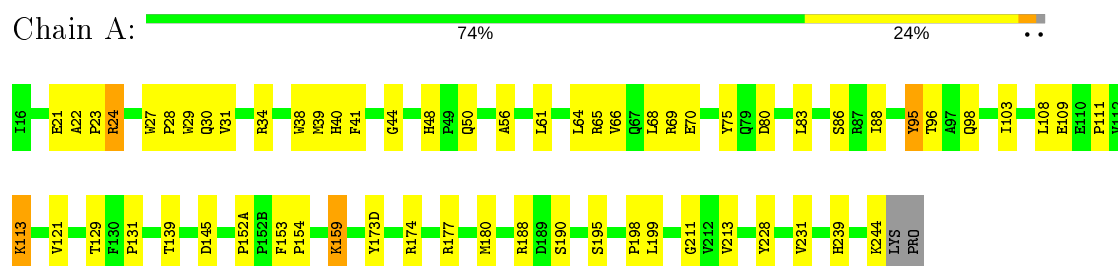
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	25	Total 25	O 25	0	0
3	B	16	Total 16	O 16	0	0
3	C	29	Total 29	O 29	0	0
3	D	12	Total 12	O 12	0	0

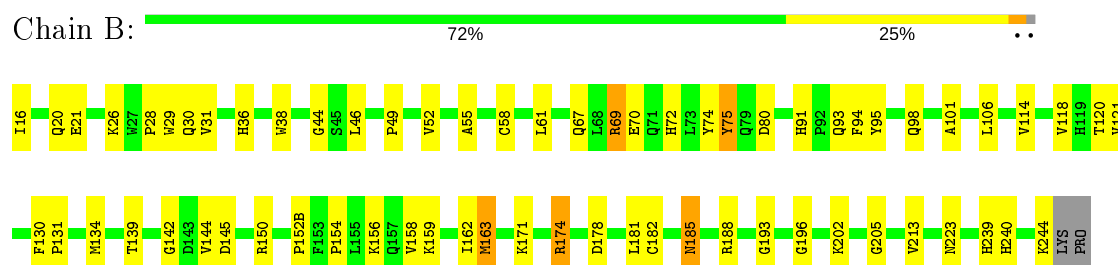
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

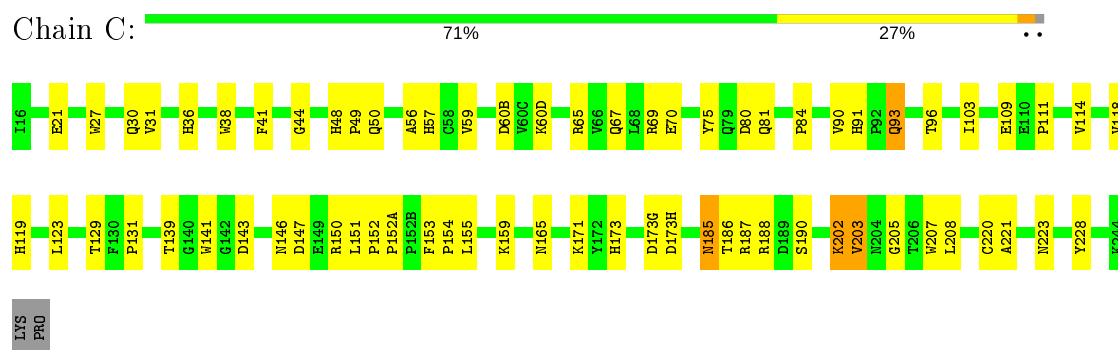
#### • Molecule 1: Trypsin beta-2



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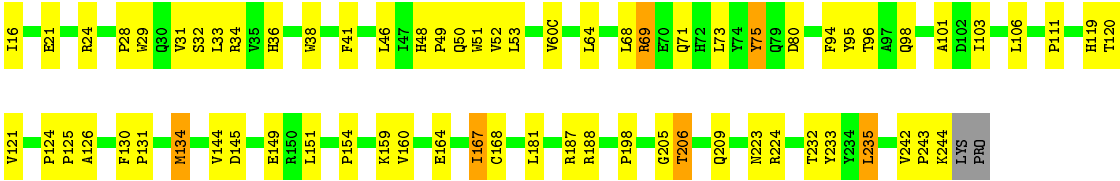


#### • Molecule 1: Trypsin beta-2



#### • Molecule 1: Trypsin beta-2





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 31	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	78.20 Å 78.20 Å 165.52 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	42.78 – 2.50 42.77 – 2.50	Depositor EDS
% Data completeness (in resolution range)	92.7 (42.78-2.50) 92.7 (42.77-2.50)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.11 (at 2.51 Å)	Xtriage
Refinement program	CNS	Depositor
R, $R_{free}$	0.220 , 0.259 0.211 , 0.249	Depositor DCC
$R_{free}$ test set	3931 reflections (10.06%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	27.6	Xtriage
Anisotropy	0.357	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 6.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.016 for -h,-k,l 0.480 for h,-h-k,-l 0.019 for -k,-h,-l	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	7923	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.11% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup> Intensities estimated from amplitudes.

<sup>2</sup> Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.



## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: C3A

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z  > 5$	RMSZ	$\# Z  > 5$
1	A	0.37	0/1982	0.66	0/2716
1	B	0.35	0/1985	0.66	0/2720
1	C	0.37	0/1985	0.65	0/2720
1	D	0.37	0/1985	0.64	0/2720
All	All	0.36	0/7937	0.65	0/10876

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	2
1	D	0	1
All	All	0	4

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	95	TYR	Sidechain
1	B	75	TYR	Sidechain
1	B	95	TYR	Sidechain
1	D	95	TYR	Sidechain

## 5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1917	0	1852	55	0
1	B	1920	0	1854	52	0
1	C	1920	0	1854	48	0
1	D	1920	0	1854	51	0
2	A	41	0	35	4	0
2	B	41	0	35	1	0
2	C	41	0	35	3	0
2	D	41	0	35	0	0
3	A	25	0	0	0	0
3	B	16	0	0	0	0
3	C	29	0	0	0	0
3	D	12	0	0	0	0
All	All	7923	0	7554	198	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 13.

All (198) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:24:ARG:HH11	1:A:24:ARG:HB3	1.26	0.98
1:B:185:ASN:ND2	1:B:188:ARG:H	1.68	0.92
1:D:232:THR:HA	1:D:235:LEU:HD22	1.52	0.90
1:B:239:HIS:HB3	1:B:244:LYS:HE2	1.53	0.88
1:B:185:ASN:HD21	1:B:188:ARG:H	1.24	0.84
1:B:91:HIS:HD2	1:B:93:GLN:H	1.26	0.83
1:C:203:VAL:HG21	1:C:208:LEU:HD12	1.63	0.81
1:C:202:LYS:HE3	1:C:205:GLY:HA2	1.62	0.80
1:A:24:ARG:NH1	1:A:24:ARG:HB3	1.97	0.79
1:C:129:THR:HG23	1:C:131:PRO:HD3	1.64	0.79
1:A:48:HIS:HD2	1:A:50:GLN:H	1.30	0.78
1:A:129:THR:HG23	1:A:131:PRO:HD3	1.66	0.77
1:C:146:ASN:HD21	1:C:220:CYS:H	1.34	0.74
1:A:75:TYR:HB2	1:B:75:TYR:CE1	2.24	0.73
1:B:174:ARG:HG2	1:B:174:ARG:HH11	1.54	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:205:GLY:O	1:D:206:THR:HG23	1.92	0.69
1:C:21:GLU:HG3	1:C:154:PRO:HB2	1.74	0.69
1:A:21:GLU:HG3	1:A:154:PRO:HB2	1.72	0.69
1:D:164:GLU:O	1:D:167:ILE:HG22	1.93	0.69
1:C:150:ARG:HH11	1:C:150:ARG:HG2	1.58	0.68
1:B:174:ARG:HH22	1:B:178:ASP:CG	1.97	0.67
1:C:202:LYS:HE3	1:C:205:GLY:CA	2.23	0.67
1:C:114:VAL:HG13	1:C:119:HIS:HB3	1.76	0.67
1:A:61:LEU:CD2	1:A:88:ILE:HG12	2.26	0.66
1:B:67:GLN:NE2	1:B:70:GLU:HG3	2.12	0.64
1:C:185:ASN:ND2	1:C:188:ARG:H	1.96	0.64
1:A:48:HIS:CD2	1:A:50:GLN:H	2.15	0.62
1:D:53:LEU:HD11	1:D:103:ILE:HD11	1.81	0.62
1:D:159:LYS:O	1:D:188:ARG:HD3	1.99	0.61
1:B:131:PRO:O	1:B:134:MET:HB2	2.01	0.61
1:C:150:ARG:NH1	1:C:150:ARG:HG2	2.17	0.60
1:B:181:LEU:HD23	1:B:182:CYS:N	2.16	0.60
1:A:98:GLN:HE22	2:A:999:C3A:C1	2.15	0.59
1:A:211:GLY:HA2	1:A:231:VAL:HG23	1.82	0.59
1:B:188:ARG:HH11	1:B:188:ARG:HG2	1.67	0.59
1:A:69:ARG:HG2	1:A:69:ARG:HH11	1.69	0.58
1:C:84:PRO:HB2	1:C:109:GLU:OE2	2.03	0.58
1:B:21:GLU:HG3	1:B:154:PRO:HB2	1.86	0.57
1:B:174:ARG:HG2	1:B:174:ARG:NH1	2.18	0.57
1:C:185:ASN:C	1:C:185:ASN:HD22	2.08	0.57
1:B:29:TRP:CG	1:B:121:VAL:HB	2.40	0.56
1:C:143:ASP:OD1	1:C:150:ARG:HD3	2.05	0.56
1:C:59:VAL:O	1:C:60(D):LYS:HE2	2.04	0.56
1:C:67:GLN:NE2	1:C:70:GLU:HG3	2.21	0.56
1:D:48:HIS:CD2	1:D:50:GLN:H	2.24	0.56
1:C:152(A):PRO:HG3	1:C:153:PHE:CZ	2.41	0.55
1:D:167:ILE:O	1:D:167:ILE:HD13	2.06	0.55
1:D:52:VAL:HB	1:D:106:LEU:HB2	1.88	0.55
1:B:185:ASN:HD22	1:B:185:ASN:H	1.53	0.55
1:D:188:ARG:HG2	1:D:188:ARG:HH11	1.70	0.55
1:A:159:LYS:O	1:A:188:ARG:HD3	2.07	0.55
1:B:185:ASN:HD21	1:B:188:ARG:N	1.99	0.54
1:C:114:VAL:HG13	1:C:119:HIS:CB	2.37	0.54
1:C:41:PHE:CZ	2:C:997:C3A:H332	2.42	0.54
1:B:36:HIS:HB2	1:B:38:TRP:CH2	2.42	0.54
1:B:46:LEU:O	1:B:120:THR:HA	2.07	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:75:TYR:HB2	1:B:75:TYR:HE1	1.73	0.53
1:B:185:ASN:N	1:B:185:ASN:HD22	2.06	0.53
1:A:61:LEU:HD21	1:A:88:ILE:HG12	1.91	0.53
1:C:69:ARG:HH11	1:C:69:ARG:HG2	1.74	0.53
1:B:213:VAL:HG13	2:B:998:C3A:H121	1.91	0.53
1:C:31:VAL:HG22	1:C:44:GLY:C	2.29	0.53
1:A:41:PHE:CZ	2:A:999:C3A:H332	2.44	0.53
1:D:34:ARG:HH21	1:D:73:LEU:HD22	1.74	0.53
1:A:177:ARG:HB2	1:A:180:MET:HE3	1.91	0.52
1:C:203:VAL:CG2	1:C:208:LEU:HD12	2.36	0.52
1:B:98:GLN:HG2	1:C:96:THR:HG21	1.91	0.52
1:D:21:GLU:HG3	1:D:154:PRO:HB2	1.93	0.52
1:C:190:SER:OG	2:C:997:C3A:H131	2.10	0.51
1:C:70:GLU:OE1	1:C:80:ASP:OD1	2.28	0.51
1:D:28:PRO:HG3	1:D:119:HIS:CE1	2.44	0.51
1:A:56:ALA:HB2	1:A:103:ILE:O	2.10	0.51
1:C:48:HIS:HD2	1:C:50:GLN:H	1.58	0.51
1:A:98:GLN:HE22	2:A:999:C3A:H12	1.76	0.50
1:D:94:PHE:HB2	1:D:101:ALA:O	2.11	0.50
1:D:126:ALA:HA	1:D:235:LEU:HD23	1.92	0.50
1:C:91:HIS:CE1	1:C:93:GLN:HB2	2.47	0.50
1:D:188:ARG:HG2	1:D:188:ARG:NH1	2.26	0.50
1:B:188:ARG:NH1	1:B:188:ARG:HG2	2.26	0.50
1:B:30:GLN:NE2	1:B:139:THR:OG1	2.44	0.50
1:C:56:ALA:HB2	1:C:103:ILE:O	2.11	0.50
1:A:66:VAL:HG11	1:A:108:LEU:CD2	2.41	0.50
1:C:141:TRP:CZ2	1:C:155:LEU:HD13	2.47	0.50
1:D:223:ASN:O	1:D:224:ARG:HD3	2.12	0.49
1:D:151:LEU:HD21	1:D:154:PRO:O	2.12	0.49
1:D:187:ARG:HH21	1:D:187:ARG:HG3	1.78	0.49
1:B:29:TRP:CD2	1:B:121:VAL:HB	2.47	0.49
1:A:239:HIS:HB3	1:A:244:LYS:HE2	1.94	0.49
1:C:185:ASN:ND2	1:C:187:ARG:H	2.10	0.49
1:B:74:TYR:HA	1:B:80:ASP:OD1	2.12	0.49
1:B:144:VAL:O	1:B:145:ASP:HB2	2.11	0.49
1:D:131:PRO:HG2	1:D:134:MET:HG2	1.94	0.49
1:A:27:TRP:N	1:A:28:PRO:HD3	2.28	0.49
1:C:173:HIS:CD2	1:C:173(G):ASP:HA	2.48	0.49
1:B:52:VAL:HB	1:B:106:LEU:HB2	1.93	0.48
1:C:185:ASN:HD22	1:C:188:ARG:H	1.59	0.48
1:C:75:TYR:CE2	1:D:75:TYR:HB2	2.49	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:27:TRP:CD1	1:C:139:THR:HG21	2.49	0.48
1:D:36:HIS:HB2	1:D:38:TRP:CH2	2.49	0.47
1:A:31:VAL:HG22	1:A:44:GLY:C	2.35	0.47
1:B:31:VAL:HG22	1:B:44:GLY:C	2.35	0.47
1:A:64:LEU:CD2	1:A:88:ILE:HD11	2.44	0.47
1:D:29:TRP:CD2	1:D:121:VAL:HB	2.49	0.47
1:B:171:LYS:HD3	1:B:223:ASN:HA	1.97	0.47
1:A:50:GLN:HG3	1:A:111:PRO:HB3	1.96	0.47
1:A:239:HIS:CG	1:A:244:LYS:HE2	2.50	0.47
1:A:30:GLN:HE22	1:A:198:PRO:HD2	1.80	0.46
1:B:118:VAL:O	1:B:118:VAL:HG23	2.15	0.46
1:D:68:LEU:O	1:D:69:ARG:C	2.53	0.46
1:A:86:SER:HB3	1:A:109:GLU:HG2	1.97	0.46
1:B:144:VAL:HG12	1:B:156:LYS:HE2	1.96	0.46
1:B:28:PRO:HA	1:B:69:ARG:HG3	1.96	0.46
1:C:49:PRO:O	1:C:111:PRO:HA	2.16	0.46
1:D:24:ARG:HA	1:D:71:GLN:NE2	2.31	0.46
1:A:113:LYS:N	1:A:113:LYS:HD2	2.30	0.46
1:A:95:TYR:CD2	1:A:96:THR:HG23	2.51	0.46
1:D:69:ARG:HH22	1:D:71:GLN:HE21	1.64	0.46
1:A:152(A):PRO:HA	1:A:153:PHE:N	2.31	0.46
1:C:36:HIS:HD2	1:C:38:TRP:CZ2	2.34	0.46
1:C:171:LYS:HD2	1:C:223:ASN:HD22	1.81	0.46
1:A:27:TRP:CD1	1:A:139:THR:HG21	2.51	0.46
1:B:16:ILE:O	1:B:144:VAL:HA	2.16	0.46
1:D:149:GLU:O	1:D:149:GLU:HG3	2.15	0.45
1:D:48:HIS:HD2	1:D:50:GLN:H	1.60	0.45
1:B:91:HIS:CD2	1:B:93:GLN:H	2.18	0.45
1:B:240:HIS:O	1:B:240:HIS:ND1	2.49	0.45
1:B:61:LEU:HD12	1:B:61:LEU:N	2.32	0.45
1:D:46:LEU:O	1:D:120:THR:HA	2.17	0.45
1:A:177:ARG:HG3	1:A:180:MET:HE1	1.99	0.44
1:A:195:SER:HA	1:A:213:VAL:HG12	1.99	0.44
1:D:34:ARG:HD2	1:D:38:TRP:O	2.17	0.44
1:C:30:GLN:NE2	1:C:139:THR:OG1	2.51	0.44
1:A:173(D):TYR:CZ	1:D:60(C):VAL:HA	2.52	0.44
1:B:185:ASN:ND2	1:B:185:ASN:N	2.64	0.44
1:A:66:VAL:HG11	1:A:108:LEU:HD21	1.99	0.44
1:A:64:LEU:HD23	1:A:88:ILE:HD11	1.98	0.44
1:A:228:TYR:CD1	1:A:228:TYR:N	2.85	0.44
1:B:158:VAL:HG21	1:B:188:ARG:HB3	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:50:GLN:OE1	1:D:111:PRO:HG3	2.17	0.44
1:A:29:TRP:CG	1:A:121:VAL:HB	2.53	0.44
1:A:177:ARG:O	1:A:180:MET:HB2	2.19	0.43
1:A:34:ARG:HA	1:A:39:MET:O	2.18	0.43
1:D:232:THR:CA	1:D:235:LEU:HD22	2.37	0.43
1:B:94:PHE:HB2	1:B:101:ALA:O	2.17	0.43
1:B:162:ILE:HG22	1:B:163:MET:N	2.34	0.43
1:A:38:TRP:CZ3	1:A:65:ARG:HG3	2.53	0.43
1:A:70:GLU:OE1	1:A:80:ASP:OD1	2.36	0.43
1:C:146:ASN:O	1:C:147:ASP:HB2	2.19	0.43
1:B:159:LYS:O	1:B:188:ARG:HD3	2.18	0.43
1:B:67:GLN:HE21	1:B:70:GLU:HG3	1.80	0.43
1:C:146:ASN:HD21	1:C:220:CYS:N	2.10	0.43
1:B:44:GLY:HA2	1:B:196:GLY:O	2.19	0.43
1:D:16:ILE:O	1:D:144:VAL:HA	2.18	0.43
1:A:34:ARG:HD3	1:A:40:HIS:HA	2.01	0.43
1:A:69:ARG:HG2	1:A:69:ARG:NH1	2.33	0.43
1:A:173(D):TYR:CE2	1:D:60(C):VAL:HA	2.54	0.42
1:A:64:LEU:O	1:A:65:ARG:HG2	2.19	0.42
1:A:61:LEU:HD23	1:A:88:ILE:HG12	1.99	0.42
1:D:130:PHE:HA	1:D:131:PRO:HD2	1.91	0.42
1:A:98:GLN:HG2	1:D:96:THR:HG21	2.01	0.42
1:B:158:VAL:CG2	1:B:188:ARG:HB3	2.50	0.42
1:B:185:ASN:C	1:B:185:ASN:HD22	2.22	0.42
1:D:187:ARG:NH2	1:D:187:ARG:HG3	2.34	0.42
1:A:48:HIS:HD2	1:A:50:GLN:N	2.07	0.42
1:A:66:VAL:HG12	1:A:83:LEU:O	2.20	0.42
1:C:186:THR:HA	1:C:221:ALA:O	2.19	0.42
1:C:67:GLN:HE21	1:C:70:GLU:HG3	1.84	0.42
1:C:202:LYS:HG2	1:C:207:TRP:CE2	2.54	0.42
1:A:152(A):PRO:HB3	1:A:153:PHE:CZ	2.54	0.42
1:D:31:VAL:CG1	1:D:32:SER:N	2.83	0.42
1:D:34:ARG:O	1:D:64:LEU:HD12	2.20	0.42
1:B:202:LYS:HE2	1:B:205:GLY:HA2	2.01	0.41
1:C:223:ASN:HD22	1:C:223:ASN:HA	1.67	0.41
1:C:57:HIS:NE2	2:C:997:C3A:N41	2.57	0.41
1:A:96:THR:HG21	1:D:98:GLN:HG2	2.02	0.41
1:D:235:LEU:HD12	1:D:235:LEU:HA	1.95	0.41
1:D:48:HIS:O	1:D:50:GLN:N	2.53	0.41
1:B:55:ALA:O	1:B:58:CYS:HB2	2.20	0.41
1:C:56:ALA:HB1	1:C:90:VAL:HG13	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:233:TYR:O	1:D:233:TYR:CD1	2.74	0.41
1:A:21:GLU:CG	1:A:154:PRO:HB2	2.46	0.41
1:C:228:TYR:N	1:C:228:TYR:CD1	2.89	0.41
1:A:22:ALA:HA	1:A:23:PRO:HD3	1.89	0.41
1:C:151:LEU:HA	1:C:152:PRO:HD3	1.88	0.41
1:B:142:GLY:HA2	1:B:193:GLY:HA3	2.02	0.40
1:D:33:LEU:HB3	1:D:41:PHE:CZ	2.56	0.40
1:D:145:ASP:HB3	1:D:149:GLU:HG2	2.03	0.40
1:D:51:TRP:CG	1:D:242:VAL:HG22	2.56	0.40
1:D:69:ARG:HH22	1:D:71:GLN:NE2	2.19	0.40
1:B:162:ILE:HD12	1:B:181:LEU:HD21	2.03	0.40
1:B:72:HIS:HB3	1:B:152(B):PRO:O	2.22	0.40
1:C:159:LYS:O	1:C:188:ARG:HD3	2.21	0.40
1:D:124:PRO:HD3	1:D:209:GLN:O	2.21	0.40
1:D:160:VAL:O	1:D:160:VAL:HG23	2.21	0.40
1:D:29:TRP:CG	1:D:121:VAL:HB	2.56	0.40
1:A:190:SER:OG	2:A:999:C3A:H131	2.22	0.40
1:B:163:MET:HB2	1:B:182:CYS:SG	2.62	0.40
1:D:124:PRO:HA	1:D:125:PRO:HD3	1.92	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles ⓘ

### 5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	241/245 (98%)	222 (92%)	19 (8%)	0	100	100
1	B	241/245 (98%)	216 (90%)	22 (9%)	3 (1%)	13	24
1	C	241/245 (98%)	223 (92%)	17 (7%)	1 (0%)	34	54
1	D	241/245 (98%)	219 (91%)	19 (8%)	3 (1%)	13	24
All	All	964/980 (98%)	880 (91%)	77 (8%)	7 (1%)	22	39

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	243	PRO
1	B	114	VAL
1	D	69	ARG
1	B	130	PHE
1	C	203	VAL
1	B	49	PRO
1	D	49	PRO

### 5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	207/210 (99%)	200 (97%)	7 (3%)	37	63
1	B	208/210 (99%)	201 (97%)	7 (3%)	37	63
1	C	208/210 (99%)	198 (95%)	10 (5%)	25	48
1	D	208/210 (99%)	198 (95%)	10 (5%)	25	48
All	All	831/840 (99%)	797 (96%)	34 (4%)	30	55

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	24	ARG
1	A	68	LEU
1	A	113	LYS
1	A	145	ASP
1	A	159	LYS
1	A	174	ARG
1	A	199	LEU
1	B	20	GLN
1	B	26	LYS
1	B	69	ARG
1	B	150	ARG
1	B	163	MET
1	B	174	ARG

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Mol	Chain	Res	Type
1	B	185	ASN
1	C	60(B)	ASP
1	C	65	ARG
1	C	81	GLN
1	C	93	GLN
1	C	118	VAL
1	C	123	LEU
1	C	165	ASN
1	C	173(H)	ASP
1	C	185	ASN
1	C	202	LYS
1	D	75	TYR
1	D	80	ASP
1	D	134	MET
1	D	167	ILE
1	D	168	CYS
1	D	181	LEU
1	D	198	PRO
1	D	206	THR
1	D	235	LEU
1	D	244	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	30	GLN
1	A	48	HIS
1	A	93	GLN
1	A	119	HIS
1	A	165	ASN
1	A	223	ASN
1	B	20	GLN
1	B	30	GLN
1	B	50	GLN
1	B	91	HIS
1	B	185	ASN
1	B	204	ASN
1	B	221(A)	GLN
1	C	30	GLN
1	C	36	HIS
1	C	48	HIS
1	C	79	GLN

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Mol	Chain	Res	Type
1	C	119	HIS
1	C	146	ASN
1	C	165	ASN
1	C	185	ASN
1	C	223	ASN
1	D	30	GLN
1	D	48	HIS
1	D	71	GLN
1	D	79	GLN
1	D	93	GLN
1	D	223	ASN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

4 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
2	C3A	D	996	1	39,45,45	2.08	8 (20%)	42,61,61	0.93	3 (7%)
2	C3A	C	997	1	39,45,45	2.16	9 (23%)	42,61,61	1.03	3 (7%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	C3A	B	998	1	39,45,45	2.20	11 (28%)	42,61,61	0.86	2 (4%)
2	C3A	A	999	1	39,45,45	2.09	10 (25%)	42,61,61	0.93	2 (4%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	C3A	D	996	1	1/1/6/10	3/28/48/48	0/5/5/5
2	C3A	C	997	1	1/1/6/10	4/28/48/48	0/5/5/5
2	C3A	B	998	1	1/1/6/10	3/28/48/48	0/5/5/5
2	C3A	A	999	1	1/1/6/10	4/28/48/48	0/5/5/5

All (38) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	998	C3A	O17-C16	9.30	1.35	1.22
2	C	997	C3A	O17-C16	9.28	1.35	1.22
2	A	999	C3A	O17-C16	9.09	1.35	1.22
2	D	996	C3A	O17-C16	8.85	1.34	1.22
2	C	997	C3A	C8-C16	4.97	1.61	1.53
2	B	998	C3A	C8-C16	4.45	1.61	1.53
2	A	999	C3A	C8-C16	4.13	1.60	1.53
2	D	996	C3A	C8-C16	4.01	1.60	1.53
2	B	998	C3A	C29-N31	2.96	1.40	1.34
2	B	998	C3A	C25-C26	2.96	1.44	1.39
2	D	996	C3A	C25-C26	2.83	1.44	1.39
2	C	997	C3A	C25-C26	2.51	1.43	1.39
2	A	999	C3A	C25-C26	2.49	1.43	1.39
2	C	997	C3A	C29-N31	2.43	1.39	1.34
2	C	997	C3A	C37-C36	2.39	1.44	1.38
2	A	999	C3A	C28-C23	2.36	1.43	1.38
2	A	999	C3A	C27-C26	2.36	1.43	1.39
2	C	997	C3A	C27-C26	2.33	1.43	1.39
2	D	996	C3A	C37-C36	2.31	1.44	1.38
2	B	998	C3A	C37-C36	2.29	1.44	1.38
2	B	998	C3A	C36-C35	2.23	1.43	1.38
2	B	998	C3A	C28-C23	2.22	1.43	1.38
2	D	996	C3A	C29-N31	2.18	1.38	1.34
2	C	997	C3A	C24-C23	2.17	1.43	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	999	C3A	C37-C36	2.15	1.43	1.38
2	A	999	C3A	C24-C23	2.15	1.43	1.38
2	D	996	C3A	C28-C23	2.14	1.43	1.38
2	A	999	C3A	C29-N31	2.12	1.38	1.34
2	C	997	C3A	C25-C24	2.12	1.42	1.38
2	B	998	C3A	C25-C24	2.11	1.42	1.38
2	B	998	C3A	C27-C26	2.07	1.42	1.39
2	A	999	C3A	C28-C27	2.05	1.42	1.38
2	B	998	C3A	C24-C23	2.05	1.43	1.38
2	D	996	C3A	C24-C23	2.05	1.43	1.38
2	A	999	C3A	C37-C38	2.04	1.43	1.38
2	C	997	C3A	C39-C34	2.02	1.43	1.39
2	B	998	C3A	O4-C5	2.02	1.39	1.35
2	D	996	C3A	C27-C26	2.00	1.42	1.39

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	996	C3A	C34-C33-C32	-2.91	99.95	102.76
2	C	997	C3A	C32-N31-C29	2.84	126.47	122.55
2	C	997	C3A	C39-C40-C32	-2.61	100.24	102.76
2	A	999	C3A	C39-C40-C32	-2.45	100.40	102.76
2	D	996	C3A	C39-C40-C32	-2.32	100.53	102.76
2	B	998	C3A	C34-C33-C32	-2.25	100.59	102.76
2	A	999	C3A	C32-N31-C29	2.11	125.47	122.55
2	B	998	C3A	C9-C10-C11	-2.09	111.72	115.17
2	C	997	C3A	C36-C35-C34	-2.09	117.78	120.89
2	D	996	C3A	C16-C8-N7	-2.03	106.94	109.85

All (4) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	D	996	C3A	C8
2	C	997	C3A	C8
2	B	998	C3A	C8
2	A	999	C3A	C8

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	996	C3A	O17-C16-C8-N7
2	D	996	C3A	N41-C21-C22-C23

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Mol	Chain	Res	Type	Atoms
2	C	997	C3A	O17-C16-C8-N7
2	C	997	C3A	N41-C21-C22-C23
2	B	998	C3A	O17-C16-C8-N7
2	B	998	C3A	N41-C21-C22-C23
2	A	999	C3A	O17-C16-C8-N7
2	A	999	C3A	N41-C21-C22-C23
2	C	997	C3A	C1-C2-C3-O4
2	A	999	C3A	C1-C2-C3-O4
2	D	996	C3A	C18-C16-C8-N7
2	C	997	C3A	C18-C16-C8-N7
2	B	998	C3A	C18-C16-C8-N7
2	A	999	C3A	C18-C16-C8-N7

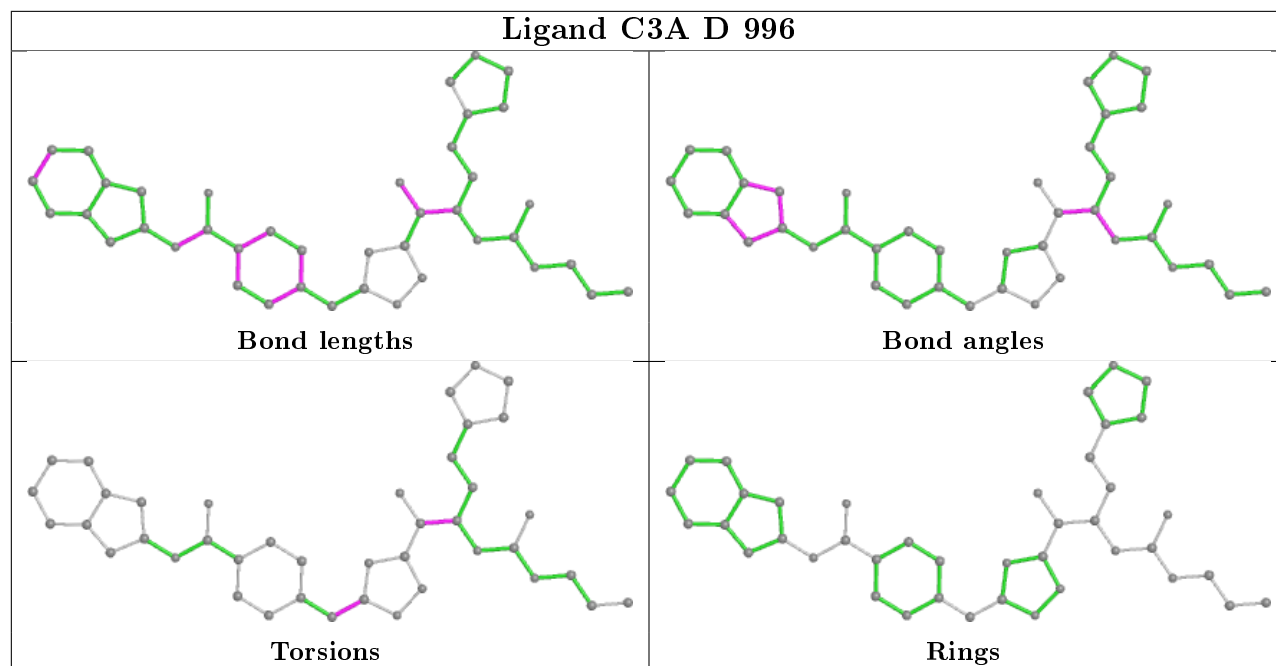
There are no ring outliers.

3 monomers are involved in 8 short contacts:

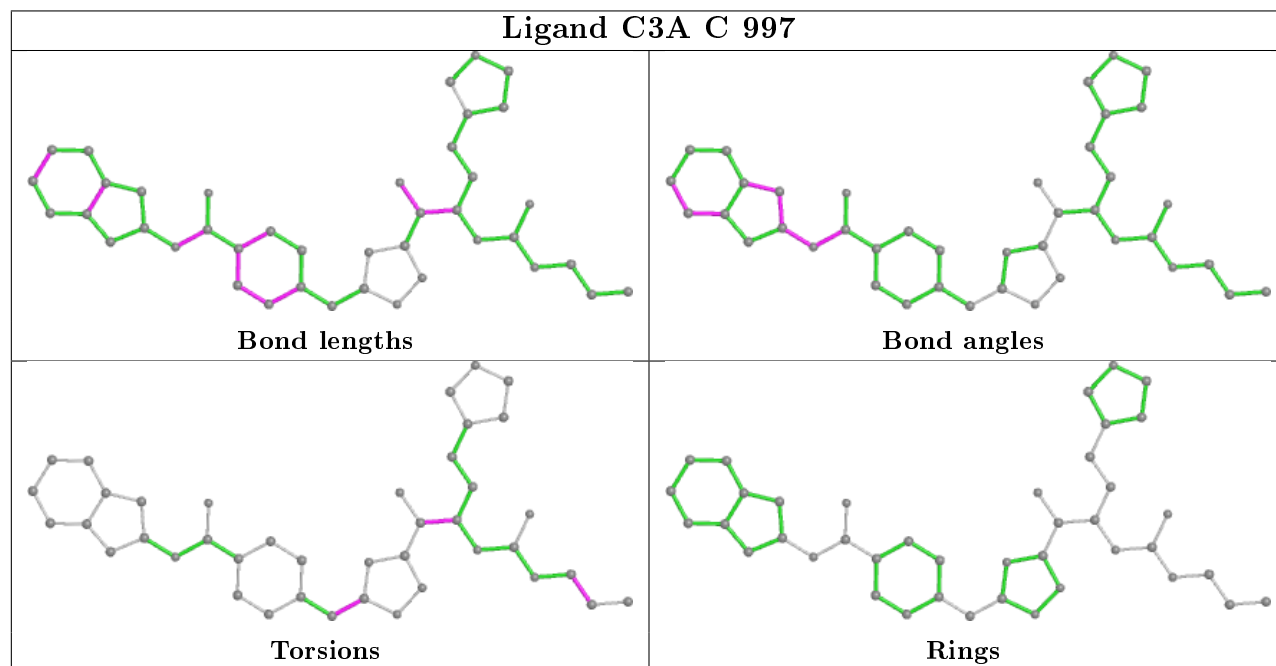
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	997	C3A	3	0
2	B	998	C3A	1	0
2	A	999	C3A	4	0

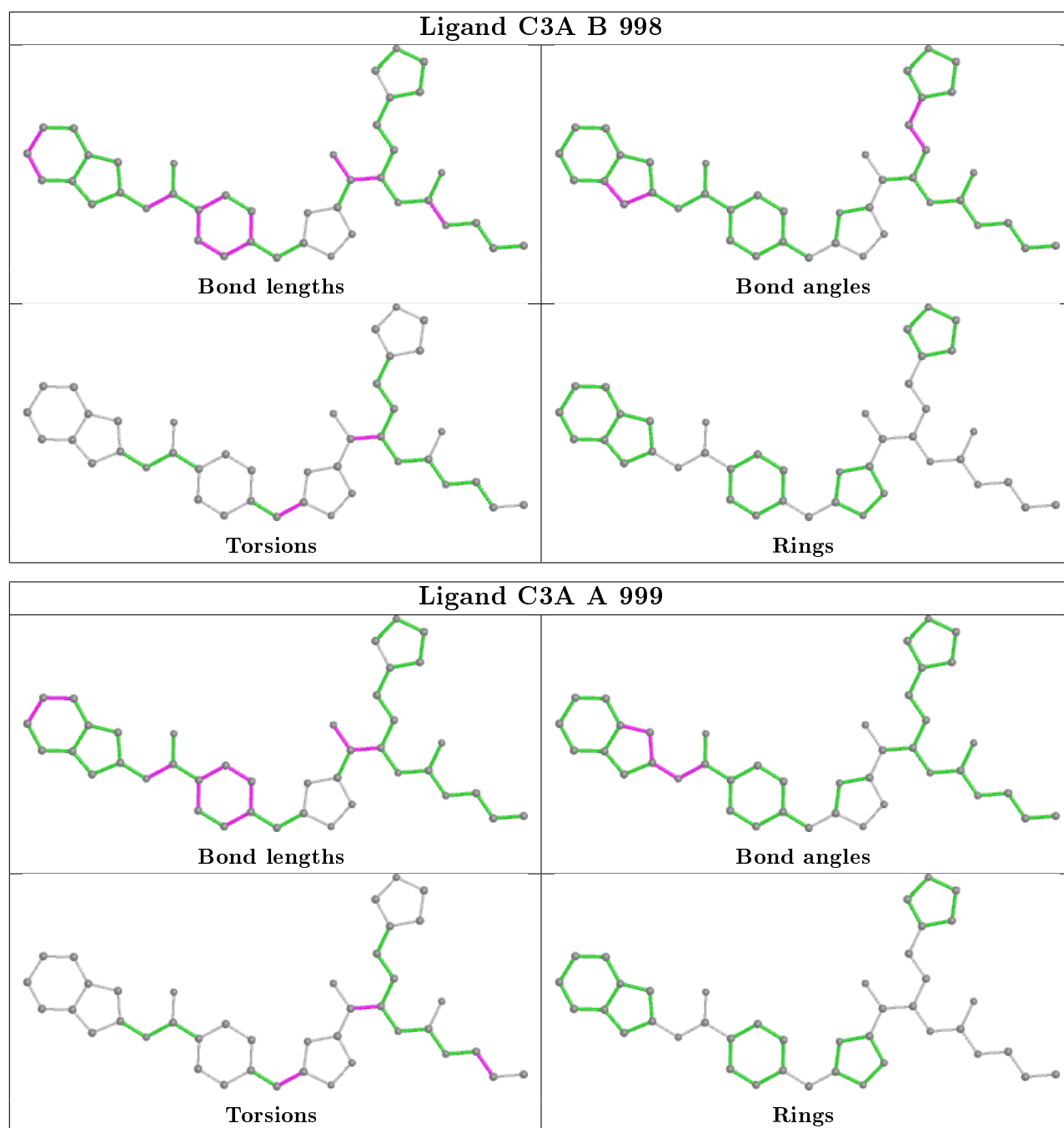
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

## Ligand C3A D 996



## Ligand C3A C 997





## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	243/245 (99%)	-0.67	0 100 100	15, 23, 41, 55	0
1	B	243/245 (99%)	-0.50	0 100 100	16, 28, 50, 61	0
1	C	243/245 (99%)	-0.67	0 100 100	15, 24, 41, 56	0
1	D	243/245 (99%)	-0.50	0 100 100	17, 28, 49, 60	0
All	All	972/980 (99%)	-0.58	0 100 100	15, 26, 46, 61	0

There are no RSRZ outliers to report.

### 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

### 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

### 6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists the number of atoms with occupancy less than 0.9.

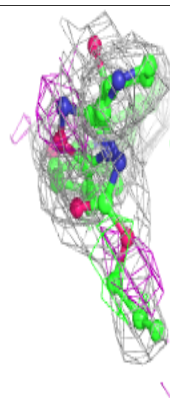
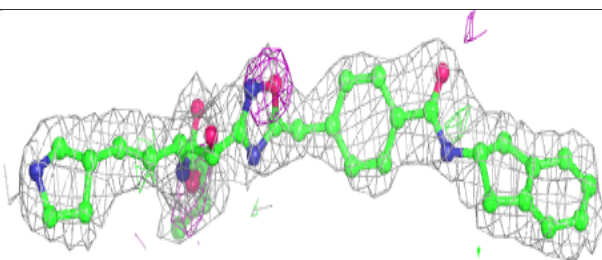
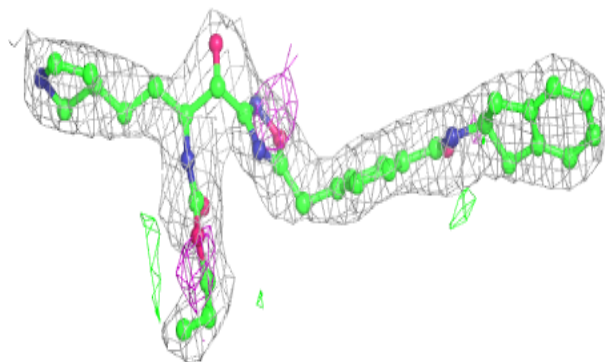
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	C3A	A	999	41/41	0.93	0.16	21,21,21,21	0
2	C3A	C	997	41/41	0.94	0.15	21,21,21,21	0
2	C3A	B	998	41/41	0.95	0.15	21,21,21,21	0
2	C3A	D	996	41/41	0.95	0.13	21,21,21,21	0



The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

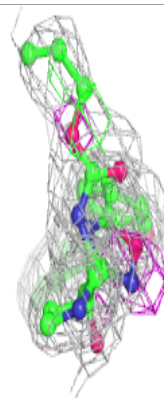
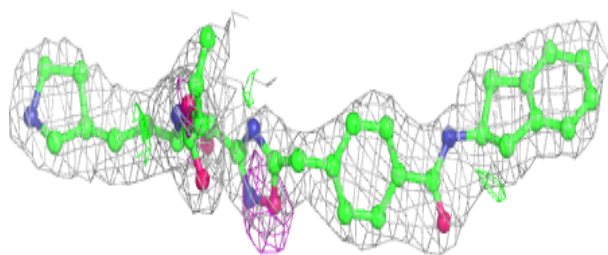
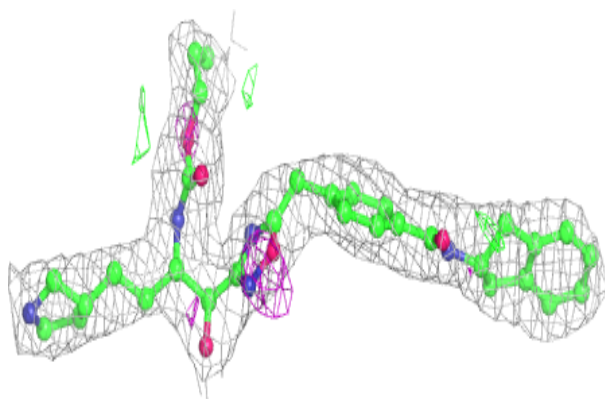
**Electron density around C3A A 999:**

2mF<sub>o</sub>-DF<sub>c</sub> (at 0.7 rmsd) in gray  
mF<sub>o</sub>-DF<sub>c</sub> (at 3 rmsd) in purple (negative)  
and green (positive)

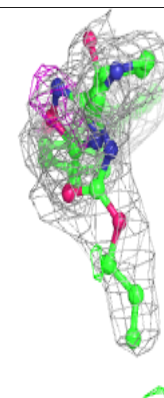
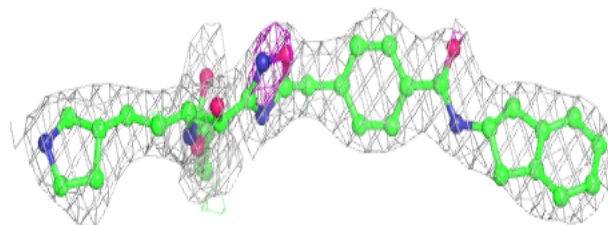
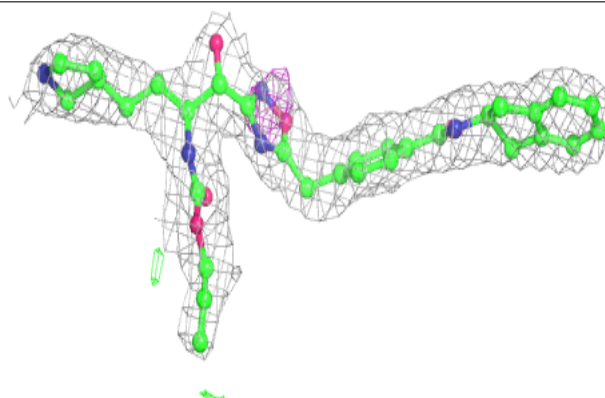


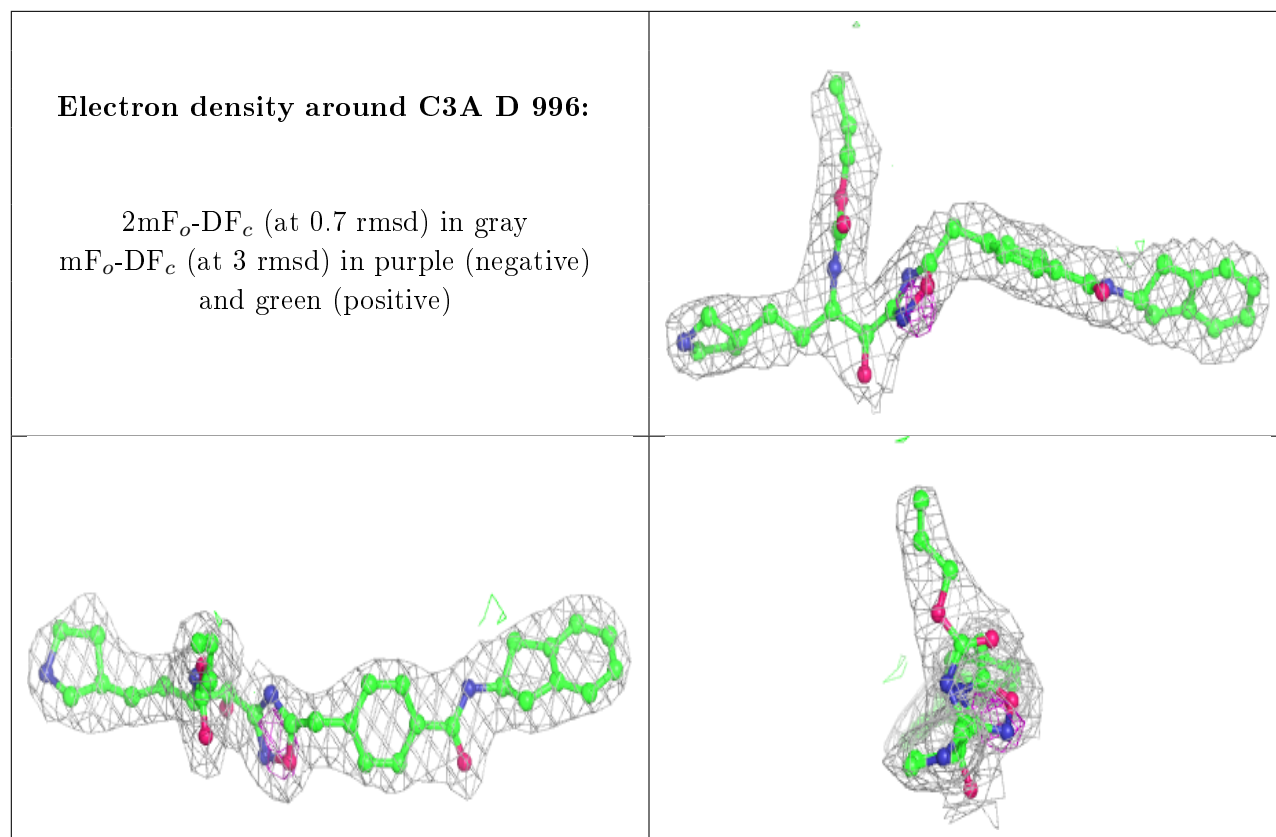
**Electron density around C3A C 997:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around C3A B 998:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





## 6.5 Other polymers [i](#)

There are no such residues in this entry.