



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 09:44 PM GMT

PDB ID : 1QFW  
Title : TERNARY COMPLEX OF HUMAN CHORIONIC GONADOTROPIN  
WITH FV ANTI ALPHA SUBUNIT AND FV ANTI BETA SUBUNIT  
Authors : Tegoni, M.; Spinelli, S.; Cambillau, C.  
Deposited on : 1999-04-15  
Resolution : 3.50 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  
<http://wwpdb.org/validation/2016/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 : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

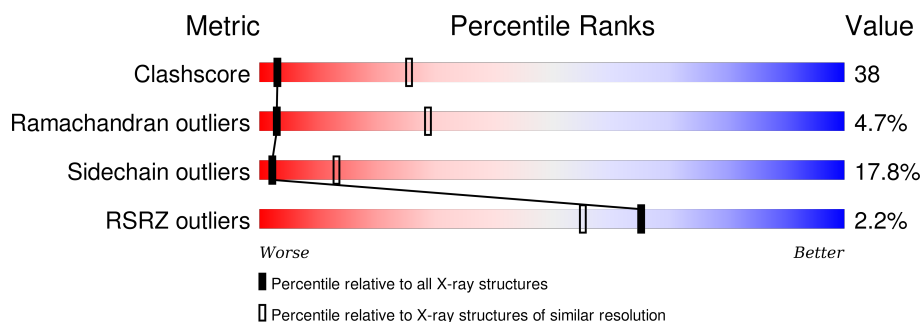
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.50 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))
Clashscore	102246	1157 (3.60-3.40)
Ramachandran outliers	100387	1120 (3.60-3.40)
Sidechain outliers	100360	1121 (3.60-3.40)
RSRZ outliers	91569	1058 (3.60-3.40)

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. 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	92	<div> <div></div> <div>51% 30% 13% 5%</div> </div>
2	B	145	<div> <div>5%</div> <div>38% 24% 8% 6% 23%</div> </div>
3	L	114	<div> <div></div> <div>35% 48% 14% ..</div> </div>
4	H	117	<div> <div>3%</div> <div>36% 40% 14% 6% .</div> </div>
5	M	108	<div> <div></div> <div>50% 42% 6% .</div> </div>
6	I	122	<div> <div></div> <div>46% 45% 8% .</div> </div>

## 2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 4938 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 GONADOTROPIN ALPHA SUBUNIT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	87	Total	C	N	O	S	19	0	0
			666	412	115	126	13			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	4	THR	VAL	CONFLICT	UNP P01215

- Molecule 2 is a protein called GONADOTROPHIN BETA SUBUNIT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	111	Total	C	N	O	S	9	0	1
			810	501	148	148	13			

- Molecule 3 is a protein called ANTIBODY (ANTI ALPHA SUBUNIT) (LIGHT CHAIN).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	L	112	Total	C	N	O	S	0	0	1
			843	531	136	172	4			

- Molecule 4 is a protein called ANTIBODY (ANTI ALPHA SUBUNIT) (HEAVY CHAIN).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	H	112	Total	C	N	O	S	0	0	1
			829	518	139	168	4			

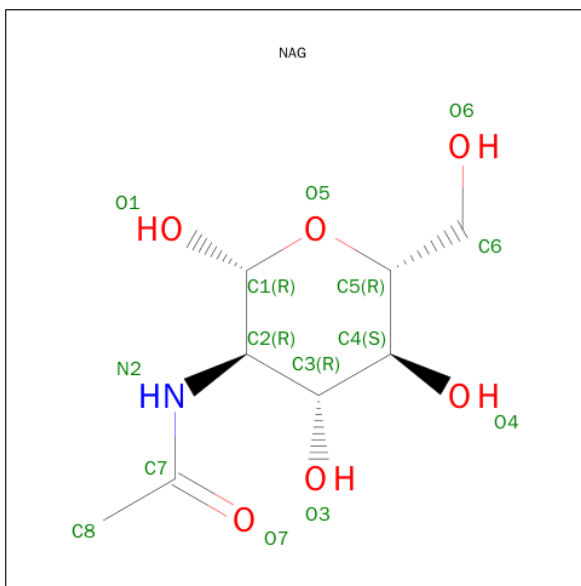
- Molecule 5 is a protein called ANTIBODY (ANTI BETA SUBUNIT) (LIGHT CHAIN).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	M	108	Total	C	N	O	S	19	0	1
			825	517	136	168	4			

- Molecule 6 is a protein called ANTIBODY (ANTI BETA SUBUNIT) (HEAVY CHAIN).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	I	121	Total	C	N	O	S	20	0	1
			937	593	163	177	4			

- Molecule 7 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula:  $C_8H_{15}NO_6$ ).

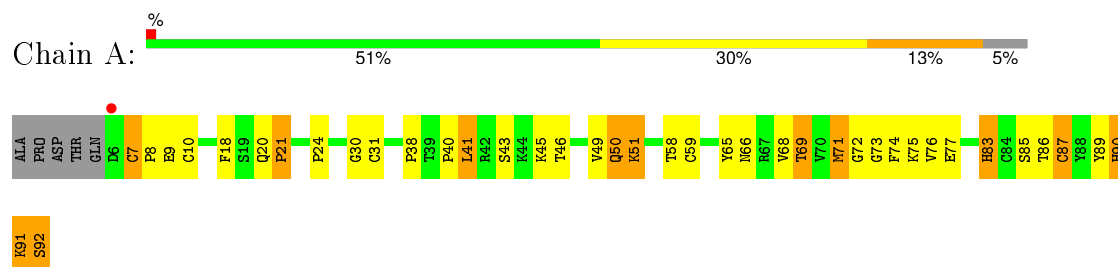


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	A	1	Total	C	N	O	0	0
			14	8	1	5		
7	A	1	Total	C	N	O	0	0
			14	8	1	5		

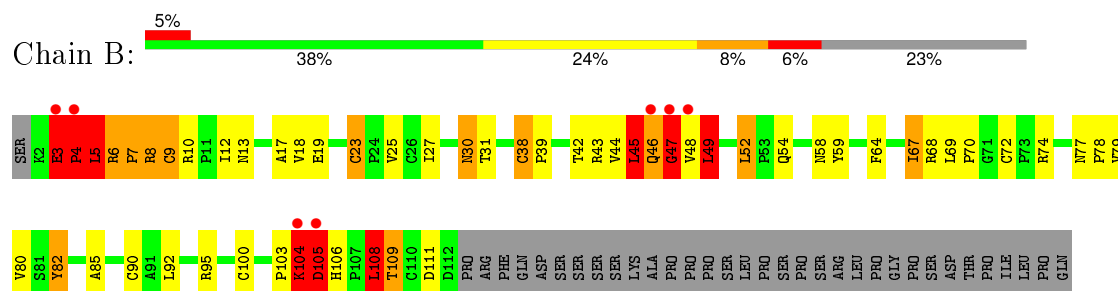
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

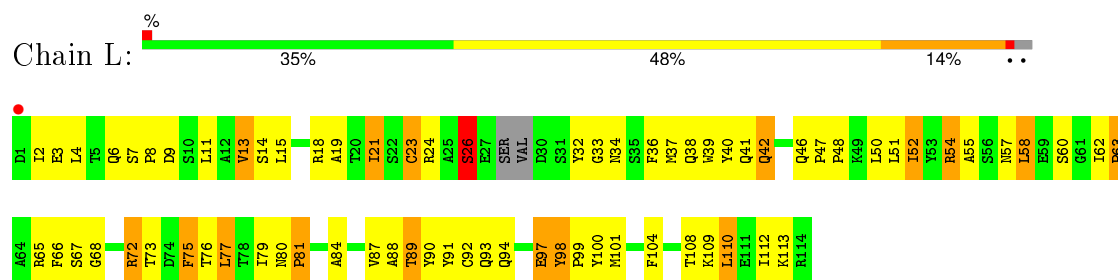
#### • Molecule 1: GONADOTROPIN ALPHA SUBUNIT



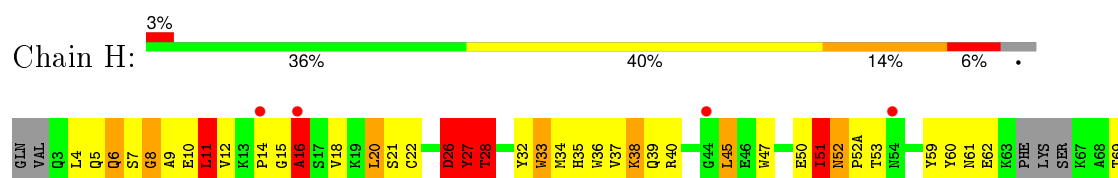
#### • Molecule 2: GONADOTROPHIN BETA SUBUNIT

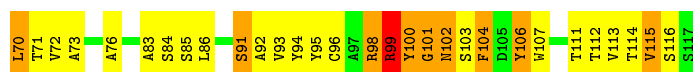


#### • Molecule 3: ANTIBODY (ANTI ALPHA SUBUNIT) (LIGHT CHAIN)



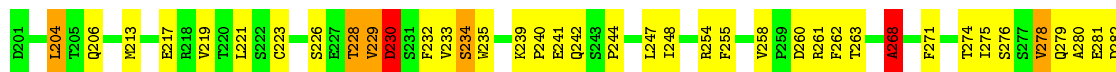
#### • Molecule 4: ANTIBODY (ANTI ALPHA SUBUNIT) (HEAVY CHAIN)





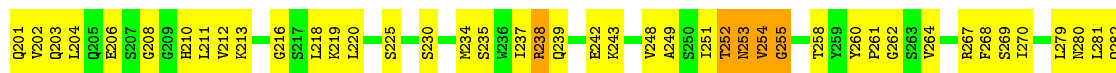
• Molecule 5: ANTIBODY (ANTI BETA SUBUNIT) (LIGHT CHAIN)

Chain M: 50% 42% 6%



• Molecule 6: ANTIBODY (ANTI BETA SUBUNIT) (HEAVY CHAIN)

Chain I: % 46% 45% 8%



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	104.90Å 104.90Å 150.10Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	9.00 – 3.50 21.50 – 3.48	Depositor EDS
% Data completeness (in resolution range)	79.0 (9.00-3.50) 95.9 (21.50-3.48)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.17	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.49 (at 3.45Å)	Xtriage
Refinement program	X-PLOR 3.843	Depositor
R, $R_{free}$	0.260 , 0.310 0.321 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	DCC
Wilson B-factor (Å <sup>2</sup> )	66.0	Xtriage
Anisotropy	0.093	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.24 , 25.1	EDS
Estimated twinning fraction	0.015 for -h,-k,l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 12091 reflections	Xtriage
$F_o, F_c$ correlation	0.74	EDS
Total number of atoms	4938	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.36% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: NAG

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	0.87	3/682 (0.4%)	0.99	4/922 (0.4%)
2	B	1.55	15/826 (1.8%)	1.96	24/1132 (2.1%)
3	L	0.76	1/862 (0.1%)	1.76	7/1172 (0.6%)
4	H	1.01	3/849 (0.4%)	1.47	19/1160 (1.6%)
5	M	0.80	1/844 (0.1%)	1.24	4/1142 (0.4%)
6	I	1.00	3/961 (0.3%)	0.99	6/1305 (0.5%)
All	All	1.03	26/5024 (0.5%)	1.45	64/6833 (0.9%)

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
2	B	0	3
3	L	0	3
4	H	0	3
5	M	0	3
All	All	0	12

All (26) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	6	ARG	C-N	20.18	1.72	1.34
6	I	252	THR	C-O	-16.64	0.91	1.23
5	M	230	ASP	C-N	-16.25	0.96	1.34
6	I	255	GLY	CA-C	-14.89	1.28	1.51
2	B	4	PRO	N-CA	14.37	1.71	1.47
2	B	4	PRO	C-N	-13.02	1.04	1.34
2	B	4	PRO	N-CD	-12.97	1.29	1.47
3	L	26	SER	C-O	-10.90	1.02	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	5	LEU	C-N	-10.43	1.10	1.34
1	A	92	SER	C-O	-9.73	1.04	1.23
4	H	100	TYR	C-N	-8.59	1.17	1.33
2	B	3	GLU	C-O	-8.34	1.07	1.23
2	B	3	GLU	CA-C	7.42	1.72	1.52
4	H	26	ASP	C-N	-7.35	1.17	1.34
2	B	4	PRO	CA-C	7.23	1.67	1.52
6	I	298	ARG	C-N	6.78	1.49	1.34
1	A	9	GLU	CG-CD	6.62	1.61	1.51
2	B	3	GLU	CB-CG	6.47	1.64	1.52
2	B	3	GLU	CG-CD	6.22	1.61	1.51
2	B	3	GLU	CA-CB	6.01	1.67	1.53
2	B	9	CYS	CB-SG	-5.79	1.72	1.81
2	B	9	CYS	CA-C	-5.61	1.38	1.52
2	B	47	GLY	C-O	-5.54	1.14	1.23
1	A	21	PRO	CA-CB	-5.30	1.43	1.53
4	H	62	GLU	CG-CD	5.22	1.59	1.51
2	B	4	PRO	CA-CB	-5.01	1.43	1.53

All (64) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	L	26	SER	O-C-N	-48.59	44.96	122.70
2	B	5	LEU	O-C-N	-37.01	63.48	122.70
5	M	230	ASP	O-C-N	-19.57	91.39	122.70
5	M	230	ASP	C-N-CA	17.96	166.61	121.70
5	M	230	ASP	CA-C-N	14.78	149.72	117.20
2	B	5	LEU	CA-C-N	13.64	147.20	117.20
4	H	28	THR	N-CA-CB	-13.28	85.07	110.30
2	B	10	ARG	NE-CZ-NH1	12.22	126.41	120.30
2	B	4	PRO	N-CA-C	11.31	141.51	112.10
4	H	102	ASN	CA-CB-CG	10.78	137.11	113.40
2	B	49	LEU	CA-CB-CG	9.57	137.32	115.30
4	H	26	ASP	CB-CA-C	-9.50	91.40	110.40
2	B	10	ARG	NE-CZ-NH2	-9.20	115.70	120.30
2	B	9	CYS	CA-C-N	-8.69	98.09	117.20
2	B	9	CYS	O-C-N	8.23	135.86	122.70
4	H	26	ASP	N-CA-CB	8.07	125.12	110.60
2	B	6	ARG	O-C-N	7.98	136.27	121.10
6	I	255	GLY	CA-C-O	7.95	134.90	120.60
2	B	9	CYS	CB-CA-C	7.85	126.11	110.40
4	H	100	TYR	O-C-N	-7.50	110.44	123.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	H	16	ALA	N-CA-C	7.34	130.81	111.00
2	B	9	CYS	C-N-CA	7.28	139.91	121.70
6	I	255	GLY	N-CA-C	7.26	131.24	113.10
2	B	4	PRO	O-C-N	-6.97	111.55	122.70
3	L	72	ARG	NE-CZ-NH1	6.97	123.79	120.30
2	B	8	ARG	O-C-N	6.80	133.58	122.70
4	H	27	TYR	CB-CG-CD1	-6.80	116.92	121.00
3	L	55	ALA	O-C-N	-6.77	111.87	122.70
3	L	54	ARG	NE-CZ-NH2	6.67	123.64	120.30
2	B	8	ARG	CA-C-N	-6.62	102.64	117.20
4	H	26	ASP	CB-CG-OD1	6.61	124.25	118.30
2	B	3	GLU	O-C-N	-6.60	108.55	121.10
2	B	3	GLU	CB-CA-C	6.59	123.58	110.40
4	H	26	ASP	O-C-N	-6.52	112.27	122.70
6	I	255	GLY	C-N-CA	-6.52	105.40	121.70
3	L	26	SER	CA-C-O	6.49	133.72	120.10
1	A	92	SER	CA-C-O	6.46	133.67	120.10
2	B	3	GLU	N-CA-C	-6.37	93.79	111.00
2	B	104	LYS	N-CA-C	6.36	128.17	111.00
4	H	91	SER	N-CA-C	6.34	128.13	111.00
2	B	45	LEU	C-N-CA	-6.29	105.97	121.70
1	A	91	LYS	CB-CA-C	-6.24	97.92	110.40
4	H	8	GLY	N-CA-C	6.24	128.71	113.10
4	H	28	THR	CB-CA-C	6.12	128.12	111.60
2	B	108	LEU	N-CA-C	6.10	127.47	111.00
3	L	99	PRO	N-CA-C	-5.93	96.67	112.10
6	I	255	GLY	CA-C-N	-5.89	104.25	117.20
4	H	102	ASN	N-CA-CB	-5.85	100.07	110.60
2	B	6	ARG	CA-C-O	-5.76	107.99	120.10
4	H	99	ARG	NE-CZ-NH2	5.74	123.17	120.30
4	H	27	TYR	C-N-CA	5.66	135.84	121.70
2	B	5	LEU	C-N-CA	5.65	135.82	121.70
3	L	58	LEU	CA-CB-CG	5.64	128.28	115.30
4	H	26	ASP	C-N-CA	5.61	135.73	121.70
2	B	3	GLU	OE1-CD-OE2	-5.52	116.68	123.30
2	B	45	LEU	N-CA-C	5.45	125.71	111.00
1	A	92	SER	CB-CA-C	-5.42	99.80	110.10
6	I	254	VAL	CA-C-N	-5.40	105.39	116.20
4	H	51	ILE	CA-C-N	-5.35	105.42	117.20
6	I	281	LEU	CA-CB-CG	5.35	127.61	115.30
4	H	34	MET	CG-SD-CE	5.35	108.75	100.20
4	H	26	ASP	CA-C-N	5.31	128.89	117.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	21	PRO	CA-CB-CG	5.28	114.84	104.80
5	M	268	ALA	N-CA-C	5.27	125.23	111.00

There are no chirality outliers.

All (12) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	47	GLY	Mainchain
2	B	5	LEU	Mainchain
2	B	7	PRO	Mainchain
4	H	106	TYR	Sidechain
4	H	11	LEU	Mainchain
4	H	26	ASP	Peptide
3	L	26	SER	Mainchain
3	L	72	ARG	Sidechain
3	L	98	TYR	Sidechain
5	M	229	VAL	Mainchain
5	M	230	ASP	Mainchain,Peptide

## 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	666	0	626	54	1
2	B	810	0	800	66	0
3	L	843	0	782	76	0
4	H	829	0	738	116	0
5	M	825	0	780	41	0
6	I	937	0	898	53	0
7	A	28	0	26	3	0
All	All	4938	0	4650	363	1

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 38.

All (363) 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)
2:B:6:ARG:C	2:B:7:PRO:N	1.72	1.40
2:B:4:PRO:N	2:B:4:PRO:CA	1.71	1.26
1:A:92:SER:HB2	2:B:92:LEU:HD23	1.27	1.15
4:H:101:GLY:O	4:H:102:ASN:HB2	1.34	1.15
4:H:51:ILE:O	4:H:52(A):PRO:HD3	1.50	1.10
1:A:92:SER:CB	2:B:92:LEU:HD23	1.90	1.01
1:A:21:PRO:CB	4:H:53:THR:HG21	1.90	1.01
2:B:46:GLN:HE21	2:B:46:GLN:H	1.06	1.01
4:H:11:LEU:O	4:H:12:VAL:HG23	1.60	0.99
1:A:21:PRO:HB2	4:H:53:THR:HG21	1.44	0.97
6:I:252:THR:O	6:I:255:GLY:HA3	1.65	0.96
4:H:20:LEU:HD21	4:H:111:THR:HG21	1.48	0.95
2:B:27:ILE:HG12	2:B:69:LEU:HD21	1.50	0.94
2:B:27:ILE:HG12	2:B:69:LEU:CD2	1.98	0.93
4:H:10:GLU:HA	4:H:113:VAL:HA	1.52	0.91
2:B:38:CYS:SG	2:B:39:PRO:HD2	2.11	0.91
4:H:37:VAL:HG21	4:H:107:TRP:HZ3	1.37	0.90
2:B:68:ARG:O	2:B:70:PRO:HD3	1.73	0.89
1:A:69:THR:HG23	1:A:75:LYS:HB3	1.54	0.88
2:B:46:GLN:HE21	2:B:46:GLN:N	1.70	0.87
2:B:23:CYS:HB3	2:B:72:CYS:HA	1.54	0.87
4:H:10:GLU:CA	4:H:113:VAL:HA	2.05	0.87
3:L:15:LEU:HD21	3:L:112:ILE:HD12	1.57	0.86
4:H:38:LYS:HD3	4:H:94:TYR:HE1	1.37	0.86
4:H:91:SER:HB3	4:H:114:THR:HA	1.59	0.83
1:A:10:CYS:HA	1:A:30:GLY:HA3	1.60	0.83
4:H:38:LYS:HD3	4:H:94:TYR:CE1	2.15	0.81
4:H:37:VAL:HG21	4:H:107:TRP:CZ3	2.14	0.81
4:H:4:LEU:HD11	4:H:22:CYS:SG	2.19	0.81
1:A:20:GLN:OE1	4:H:33:TRP:CD1	2.34	0.81
2:B:3:GLU:HG2	2:B:4:PRO:HD3	1.63	0.80
3:L:40:TYR:HE2	4:H:107:TRP:HZ2	1.32	0.78
1:A:21:PRO:HB3	4:H:53:THR:HG21	1.64	0.77
2:B:47:GLY:O	2:B:49:LEU:N	2.17	0.77
4:H:52:ASN:OD1	4:H:53:THR:OG1	2.01	0.77
3:L:7:SER:HB3	3:L:8:PRO:HD3	1.66	0.76
4:H:104:PHE:N	4:H:104:PHE:HD1	1.84	0.76
3:L:80:ASN:HB2	3:L:81:PRO:HD3	1.67	0.76
6:I:237:ILE:HD11	6:I:309:PHE:CD2	2.19	0.76
3:L:19:ALA:HB3	3:L:79:ILE:HB	1.68	0.76
1:A:90:HIS:CD2	1:A:90:HIS:H	2.03	0.75
1:A:76:VAL:HB	2:B:44:VAL:HG12	1.66	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:52:ILE:HB	3:L:57:ASN:O	1.86	0.75
1:A:76:VAL:HG12	2:B:45:LEU:HD13	1.68	0.75
5:M:206:GLN:HE21	5:M:299:GLY:HA3	1.53	0.74
4:H:35:HIS:HE1	4:H:99:ARG:HB2	1.51	0.74
6:I:237:ILE:HD11	6:I:309:PHE:CE2	2.22	0.73
6:I:304:GLN:HB3	6:I:306:TYR:CZ	2.23	0.73
6:I:206:GLU:OE2	6:I:313:GLY:HA3	1.87	0.73
1:A:18:PHE:CE1	2:B:45:LEU:CD2	2.71	0.72
3:L:6:GLN:OE1	3:L:108:THR:HG23	1.89	0.72
4:H:104:PHE:N	4:H:104:PHE:CD1	2.57	0.72
1:A:18:PHE:CE1	2:B:45:LEU:HD22	2.25	0.71
3:L:41:GLN:HG3	3:L:90:TYR:HE1	1.56	0.71
3:L:52:ILE:HG22	3:L:58:LEU:HB3	1.70	0.71
3:L:75:PHE:H	3:L:75:PHE:HD1	1.38	0.71
6:I:230:SER:O	6:I:253:ASN:HB2	1.91	0.71
6:I:203:GLN:HB2	6:I:225:SER:OG	1.91	0.71
2:B:104:LYS:NZ	2:B:105:ASP:H	1.88	0.71
2:B:17:ALA:HB2	2:B:109:THR:HA	1.71	0.71
1:A:21:PRO:HB3	4:H:53:THR:CG2	2.20	0.70
3:L:37:MET:HE2	3:L:75:PHE:CE1	2.24	0.70
4:H:18:VAL:HG13	4:H:86:LEU:HG	1.72	0.70
1:A:40:PRO:HG2	1:A:43:SER:OG	1.90	0.70
1:A:21:PRO:CB	4:H:53:THR:CG2	2.69	0.69
4:H:47:TRP:HZ2	4:H:50:GLU:HG3	1.58	0.69
5:M:290:GLN:HG2	5:M:292:TYR:H	1.58	0.69
4:H:39:GLN:HB2	4:H:45:LEU:HB3	1.74	0.68
1:A:20:GLN:CD	1:A:21:PRO:HD2	2.14	0.68
6:I:291:THR:HG22	6:I:320:VAL:HG23	1.76	0.68
3:L:8:PRO:HG2	3:L:21:ILE:HA	1.76	0.68
4:H:14:PRO:HA	4:H:86:LEU:O	1.93	0.68
3:L:11:LEU:HD22	3:L:21:ILE:HG23	1.75	0.68
6:I:237:ILE:HD12	6:I:312:TRP:CH2	2.29	0.68
2:B:4:PRO:N	2:B:4:PRO:HA	2.03	0.67
3:L:84:ALA:O	3:L:87:VAL:HG23	1.95	0.67
6:I:252:THR:O	6:I:255:GLY:CA	2.40	0.67
4:H:10:GLU:HB2	4:H:114:THR:H	1.60	0.67
6:I:304:GLN:HB3	6:I:306:TYR:CE1	2.30	0.67
3:L:52:ILE:HG22	3:L:58:LEU:CB	2.24	0.67
4:H:60:TYR:OH	4:H:70:LEU:HD22	1.94	0.67
2:B:3:GLU:CG	2:B:4:PRO:HD3	2.25	0.66
2:B:74:ARG:HH11	6:I:253:ASN:HD21	1.42	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:75:PHE:CD1	3:L:75:PHE:N	2.62	0.66
4:H:35:HIS:HB3	4:H:50:GLU:HA	1.78	0.66
5:M:230:ASP:O	5:M:230:ASP:OD1	2.14	0.66
1:A:92:SER:HB2	2:B:92:LEU:CD2	2.16	0.65
3:L:6:GLN:HE22	3:L:91:TYR:HA	1.60	0.65
3:L:80:ASN:CB	3:L:81:PRO:HD3	2.27	0.65
4:H:20:LEU:HD12	4:H:36:TRP:HZ3	1.62	0.65
3:L:14:SER:HB3	3:L:113:LYS:HB2	1.78	0.65
4:H:51:ILE:O	4:H:52(A):PRO:CD	2.36	0.65
1:A:89:TYR:HB2	2:B:54:GLN:HG3	1.78	0.65
4:H:99:ARG:HA	4:H:103:SER:O	1.97	0.64
3:L:3:GLU:H	3:L:26:SER:HB3	1.62	0.64
2:B:43:ARG:NH1	2:B:52:LEU:HB2	2.11	0.64
5:M:230:ASP:O	5:M:230:ASP:CG	2.36	0.64
6:I:234:MET:HB3	6:I:279:LEU:HD22	1.80	0.64
3:L:40:TYR:CE1	3:L:50:LEU:HD13	2.33	0.64
3:L:41:GLN:HG3	3:L:90:TYR:CE1	2.33	0.63
4:H:32:TYR:O	4:H:52(A):PRO:HG2	1.99	0.63
4:H:51:ILE:HD12	4:H:72:VAL:HG21	1.80	0.63
4:H:11:LEU:O	4:H:12:VAL:CG2	2.41	0.63
4:H:6:GLN:HG3	4:H:96:CYS:SG	2.39	0.63
5:M:254:ARG:HH12	5:M:263:THR:HA	1.64	0.62
4:H:16:ALA:O	4:H:84:SER:HA	1.99	0.62
4:H:8:GLY:HA3	4:H:111:THR:OG1	2.00	0.62
5:M:241:GLU:O	5:M:242:GLN:HG2	1.99	0.62
6:I:288:SER:O	6:I:291:THR:HG23	1.99	0.62
1:A:49:VAL:O	1:A:51:LYS:HG2	1.99	0.62
4:H:10:GLU:CB	4:H:113:VAL:HA	2.30	0.62
3:L:52:ILE:HA	3:L:58:LEU:HA	1.82	0.62
3:L:40:TYR:HE2	4:H:107:TRP:CZ2	2.17	0.61
3:L:89:THR:HB	3:L:91:TYR:HE1	1.65	0.61
2:B:45:LEU:O	2:B:45:LEU:HD12	2.00	0.61
6:I:253:ASN:N	6:I:253:ASN:HD22	1.99	0.61
4:H:8:GLY:HA3	4:H:111:THR:HG23	1.81	0.61
4:H:20:LEU:HD12	4:H:36:TRP:CZ3	2.36	0.61
3:L:6:GLN:NE2	3:L:92:CYS:H	1.98	0.60
1:A:92:SER:OG	2:B:92:LEU:HD23	2.00	0.60
3:L:14:SER:CB	3:L:113:LYS:HB2	2.32	0.60
2:B:74:ARG:NH1	6:I:253:ASN:HD21	2.00	0.60
3:L:23:CYS:SG	3:L:37:MET:HE1	2.42	0.60
3:L:89:THR:HA	3:L:109:LYS:HA	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:H:71:THR:HG22	4:H:72:VAL:N	2.17	0.59
3:L:93:GLN:HG2	3:L:94:GLN:N	2.17	0.59
5:M:254:ARG:HH11	5:M:254:ARG:HG3	1.67	0.59
3:L:40:TYR:HE1	3:L:50:LEU:HD13	1.67	0.59
1:A:20:GLN:NE2	4:H:32:TYR:HA	2.17	0.59
6:I:249:ALA:HB1	6:I:270:ILE:HD12	1.84	0.59
4:H:93:VAL:HA	4:H:111:THR:O	2.03	0.59
1:A:76:VAL:HB	2:B:44:VAL:CG1	2.31	0.59
1:A:20:GLN:NE2	1:A:21:PRO:HD2	2.18	0.59
4:H:26:ASP:O	4:H:27:TYR:HB2	2.02	0.59
6:I:258:THR:HG22	6:I:260:TYR:CE1	2.37	0.59
4:H:9:ALA:O	4:H:113:VAL:HG22	2.03	0.58
6:I:248:VAL:HA	6:I:261:PRO:HG3	1.83	0.58
3:L:89:THR:HB	3:L:91:TYR:CE1	2.38	0.58
2:B:74:ARG:NH1	6:I:253:ASN:ND2	2.51	0.58
4:H:4:LEU:CD1	4:H:22:CYS:SG	2.90	0.58
2:B:74:ARG:HH21	6:I:305:PRO:HB3	1.69	0.58
3:L:11:LEU:HG	3:L:13:VAL:HG12	1.86	0.58
1:A:20:GLN:OE1	4:H:33:TRP:HD1	1.82	0.57
6:I:212:VAL:HG11	6:I:218:LEU:HG	1.86	0.57
5:M:229:VAL:O	5:M:268:ALA:O	2.22	0.57
2:B:74:ARG:HE	6:I:305:PRO:HG3	1.70	0.57
2:B:18:VAL:HG22	2:B:80:VAL:HG11	1.87	0.56
3:L:14:SER:HB3	3:L:113:LYS:HD3	1.87	0.56
2:B:104:LYS:HZ1	2:B:105:ASP:H	1.53	0.56
1:A:77:GLU:O	2:B:44:VAL:HG13	2.06	0.55
5:M:233:VAL:HG21	5:M:271:PHE:CE1	2.40	0.55
2:B:27:ILE:CG1	2:B:69:LEU:CD2	2.80	0.55
4:H:115:VAL:HG13	4:H:116:SER:N	2.20	0.55
5:M:234:SER:O	5:M:288:CYS:HA	2.06	0.55
1:A:8:PRO:O	1:A:31:CYS:HB2	2.07	0.54
2:B:27:ILE:CG1	2:B:69:LEU:HD21	2.31	0.54
4:H:99:ARG:CG	4:H:99:ARG:HH11	2.20	0.54
5:M:261:ARG:NH2	5:M:282:ASP:OD2	2.40	0.54
3:L:68:GLY:HA2	3:L:76:THR:O	2.08	0.54
1:A:41:LEU:HD22	1:A:41:LEU:O	2.07	0.54
1:A:18:PHE:CE1	2:B:45:LEU:HD23	2.41	0.54
6:I:219:LYS:HB2	6:I:282:GLN:OE1	2.08	0.54
5:M:306:ILE:O	5:M:307:LYS:HB2	2.08	0.54
1:A:21:PRO:HB2	4:H:53:THR:CG2	2.30	0.54
2:B:104:LYS:HZ1	2:B:105:ASP:CB	2.20	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:M:286:TYR:O	5:M:301:GLY:HA2	2.08	0.53
2:B:82:TYR:N	2:B:82:TYR:CD1	2.76	0.53
4:H:8:GLY:HA3	4:H:111:THR:CB	2.37	0.53
4:H:35:HIS:CE1	4:H:99:ARG:HB2	2.39	0.53
4:H:20:LEU:HD13	4:H:21:SER:H	1.74	0.53
4:H:6:GLN:HA	4:H:22:CYS:HA	1.91	0.53
1:A:10:CYS:CA	1:A:30:GLY:HA3	2.35	0.53
3:L:19:ALA:CB	3:L:79:ILE:HB	2.36	0.53
2:B:59:TYR:CE1	2:B:85:ALA:HB2	2.44	0.53
4:H:51:ILE:HD13	4:H:52(A):PRO:N	2.24	0.53
4:H:99:ARG:HG3	4:H:99:ARG:NH1	2.24	0.52
4:H:10:GLU:HB3	4:H:112:THR:O	2.10	0.52
4:H:6:GLN:NE2	4:H:95:TYR:HA	2.24	0.52
5:M:261:ARG:HH22	5:M:282:ASP:CG	2.13	0.52
5:M:213:MET:CE	5:M:219:VAL:HG12	2.39	0.52
6:I:252:THR:O	6:I:255:GLY:N	2.42	0.52
6:I:212:VAL:HG23	6:I:320:VAL:HG22	1.91	0.52
5:M:223:CYS:HB2	5:M:235:TRP:CH2	2.44	0.52
3:L:76:THR:HG22	3:L:77:LEU:N	2.25	0.52
5:M:228:THR:HG23	5:M:228:THR:O	2.10	0.52
4:H:51:ILE:CD1	4:H:52(A):PRO:HA	2.40	0.52
3:L:21:ILE:HB	3:L:108:THR:HG21	1.92	0.52
3:L:37:MET:SD	3:L:93:GLN:O	2.67	0.52
3:L:3:GLU:H	3:L:26:SER:CB	2.23	0.52
4:H:91:SER:CA	4:H:113:VAL:O	2.58	0.52
2:B:47:GLY:O	2:B:48:VAL:C	2.42	0.52
2:B:23:CYS:HB3	2:B:72:CYS:CA	2.30	0.51
5:M:254:ARG:NH1	5:M:262:PHE:O	2.44	0.51
3:L:62:ILE:HG23	3:L:63:PRO:HD2	1.92	0.51
4:H:8:GLY:HA3	4:H:111:THR:CG2	2.39	0.51
5:M:204:LEU:HD21	5:M:290:GLN:HB3	1.91	0.51
6:I:204:LEU:HD23	6:I:311:VAL:O	2.11	0.51
4:H:32:TYR:HH	4:H:100:TYR:HE1	1.58	0.51
4:H:10:GLU:O	4:H:114:THR:N	2.44	0.51
3:L:84:ALA:HA	3:L:112:ILE:HD13	1.92	0.51
5:M:233:VAL:HG21	5:M:271:PHE:CD1	2.46	0.51
6:I:251:ILE:HD11	6:I:255:GLY:O	2.10	0.50
4:H:6:GLN:HG3	4:H:96:CYS:HB3	1.93	0.50
4:H:99:ARG:HG3	4:H:99:ARG:HH11	1.76	0.50
2:B:8:ARG:HG3	2:B:9:CYS:H	1.75	0.50
4:H:91:SER:HA	4:H:113:VAL:O	2.10	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:H:18:VAL:HG22	4:H:86:LEU:HD11	1.93	0.50
3:L:51:LEU:HD12	3:L:77:LEU:HD11	1.94	0.50
1:A:90:HIS:O	2:B:92:LEU:HG	2.11	0.50
4:H:15:GLY:O	4:H:86:LEU:HD23	2.12	0.50
5:M:206:GLN:NE2	5:M:299:GLY:HA3	2.23	0.49
1:A:83:HIS:O	1:A:83:HIS:ND1	2.45	0.49
3:L:23:CYS:SG	3:L:37:MET:CE	3.00	0.49
1:A:58:THR:C	1:A:59:CYS:SG	2.90	0.49
2:B:104:LYS:NZ	2:B:105:ASP:N	2.58	0.49
2:B:19:GLU:HA	2:B:69:LEU:HD11	1.94	0.49
6:I:212:VAL:CG1	6:I:218:LEU:HG	2.42	0.49
3:L:91:TYR:CD1	3:L:91:TYR:N	2.81	0.49
6:I:267:ARG:NH1	6:I:286:LEU:HA	2.27	0.49
6:I:310:ASP:O	6:I:312:TRP:CD1	2.66	0.49
4:H:50:GLU:HB2	4:H:59:TYR:HB2	1.95	0.48
2:B:69:LEU:O	2:B:72:CYS:SG	2.72	0.48
5:M:290:GLN:NE2	5:M:297:THR:OG1	2.46	0.48
4:H:84:SER:O	4:H:85:SER:HB3	2.14	0.48
2:B:18:VAL:CG2	2:B:80:VAL:HG11	2.44	0.48
2:B:6:ARG:CA	2:B:7:PRO:N	2.72	0.48
7:A:93:NAG:O7	7:A:93:NAG:H3	2.13	0.48
4:H:71:THR:HG22	4:H:72:VAL:H	1.78	0.47
5:M:232:PHE:O	5:M:290:GLN:HA	2.14	0.47
4:H:27:TYR:HD1	4:H:27:TYR:HA	1.38	0.47
4:H:60:TYR:CZ	4:H:70:LEU:HD22	2.49	0.47
4:H:8:GLY:CA	4:H:111:THR:HG23	2.44	0.47
4:H:18:VAL:CG2	4:H:83:ALA:HB3	2.44	0.47
3:L:39:TRP:CZ3	3:L:92:CYS:HB3	2.49	0.47
3:L:19:ALA:HB3	3:L:79:ILE:CB	2.41	0.47
3:L:98:TYR:C	3:L:100:TYR:H	2.18	0.47
4:H:32:TYR:OH	4:H:100:TYR:HE1	1.98	0.47
2:B:74:ARG:HD3	6:I:252:THR:CG2	2.45	0.47
1:A:10:CYS:HA	1:A:30:GLY:CA	2.39	0.47
4:H:98:ARG:O	4:H:104:PHE:HA	2.14	0.47
4:H:18:VAL:CG1	4:H:86:LEU:HG	2.42	0.47
4:H:115:VAL:CG1	4:H:116:SER:N	2.77	0.47
3:L:24:ARG:HA	3:L:73:THR:O	2.15	0.46
4:H:94:TYR:N	4:H:111:THR:O	2.48	0.46
4:H:18:VAL:HG22	4:H:83:ALA:HB3	1.98	0.46
3:L:52:ILE:HG22	3:L:58:LEU:HB2	1.98	0.46
3:L:34:ASN:HD22	3:L:34:ASN:N	2.14	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:7:SER:CB	3:L:8:PRO:HD3	2.40	0.46
1:A:87:CYS:SG	2:B:54:GLN:NE2	2.89	0.46
3:L:36:PHE:CD2	4:H:102:ASN:OD1	2.70	0.45
3:L:6:GLN:HA	3:L:23:CYS:HA	1.97	0.45
3:L:101:MET:HA	4:H:47:TRP:CZ3	2.50	0.45
1:A:72:GLY:O	1:A:74:PHE:N	2.50	0.45
1:A:65:TYR:O	1:A:65:TYR:CD2	2.69	0.45
1:A:68:VAL:HG21	7:A:94:NAG:H83	1.98	0.45
6:I:242:GLU:O	6:I:243:LYS:HG2	2.16	0.45
6:I:251:ILE:HB	6:I:270:ILE:HD13	1.98	0.45
4:H:10:GLU:HB3	4:H:112:THR:HG23	1.97	0.45
4:H:40:ARG:HA	4:H:92:ALA:CB	2.46	0.45
2:B:30:ASN:H	2:B:30:ASN:HD22	1.63	0.45
1:A:8:PRO:O	1:A:31:CYS:N	2.50	0.45
3:L:42:GLN:O	3:L:88:ALA:HB1	2.17	0.45
1:A:18:PHE:HE1	2:B:45:LEU:CD2	2.25	0.45
4:H:37:VAL:O	4:H:37:VAL:HG23	2.17	0.45
1:A:90:HIS:HD2	1:A:90:HIS:H	1.60	0.45
6:I:251:ILE:HD12	6:I:251:ILE:HA	1.84	0.45
6:I:260:TYR:CE2	6:I:268:PHE:O	2.69	0.45
4:H:73:ALA:HB3	4:H:76:ALA:HB3	1.97	0.45
3:L:54:ARG:HG3	3:L:54:ARG:NH1	2.32	0.45
6:I:264:VAL:HG13	6:I:268:PHE:CG	2.51	0.45
3:L:112:ILE:O	3:L:112:ILE:HG13	2.15	0.45
4:H:96:CYS:O	4:H:96:CYS:SG	2.75	0.45
6:I:253:ASN:H	6:I:253:ASN:HD22	1.65	0.44
4:H:98:ARG:HH21	4:H:99:ARG:C	2.20	0.44
3:L:98:TYR:C	3:L:100:TYR:N	2.68	0.44
4:H:51:ILE:HD13	4:H:52(A):PRO:CA	2.48	0.44
3:L:37:MET:HE2	3:L:75:PHE:CD1	2.52	0.44
4:H:51:ILE:HD12	4:H:72:VAL:CG2	2.47	0.44
3:L:40:TYR:CE2	4:H:107:TRP:CZ2	3.02	0.44
1:A:24:PRO:O	7:A:94:NAG:O7	2.34	0.44
3:L:89:THR:HG23	3:L:109:LYS:HB2	1.98	0.44
3:L:90:TYR:HE2	3:L:110:LEU:HD12	1.82	0.44
5:M:248:ILE:HG12	5:M:254:ARG:HA	2.00	0.44
1:A:7:CYS:HA	1:A:8:PRO:HD2	1.58	0.44
5:M:239:LYS:HG3	5:M:240:PRO:HD2	1.98	0.44
2:B:67:ILE:O	2:B:80:VAL:HG12	2.18	0.44
5:M:304:LEU:HA	5:M:304:LEU:HD23	1.67	0.44
1:A:20:GLN:CG	1:A:21:PRO:HD2	2.48	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:M:247:LEU:O	5:M:248:ILE:HG13	2.18	0.44
6:I:253:ASN:C	6:I:255:GLY:N	2.66	0.44
5:M:290:GLN:HE22	5:M:297:THR:H	1.66	0.44
4:H:100:TYR:CG	4:H:100:TYR:O	2.71	0.43
1:A:20:GLN:HE21	4:H:32:TYR:HA	1.83	0.43
4:H:6:GLN:HG3	4:H:96:CYS:CB	2.47	0.43
5:M:261:ARG:NH2	5:M:282:ASP:CG	2.72	0.43
4:H:6:GLN:HB2	4:H:96:CYS:SG	2.58	0.43
6:I:264:VAL:CG1	6:I:268:PHE:HB2	2.48	0.43
2:B:3:GLU:CB	2:B:4:PRO:CD	2.96	0.43
5:M:217:GLU:O	5:M:278:VAL:HG12	2.18	0.43
2:B:67:ILE:N	2:B:82:TYR:HE1	2.17	0.43
2:B:3:GLU:CB	2:B:4:PRO:HD3	2.49	0.43
6:I:239:GLN:O	6:I:292:ALA:HB1	2.19	0.43
4:H:26:ASP:O	4:H:27:TYR:CB	2.67	0.43
5:M:213:MET:HG3	5:M:219:VAL:HG11	2.00	0.43
5:M:219:VAL:HG22	5:M:275:ILE:HB	1.99	0.43
4:H:52:ASN:C	4:H:53:THR:H	2.22	0.43
3:L:8:PRO:O	3:L:108:THR:HG22	2.19	0.43
5:M:206:GLN:OE1	5:M:301:GLY:HA2	2.18	0.43
4:H:86:LEU:N	4:H:86:LEU:HD22	2.34	0.43
6:I:258:THR:CG2	6:I:260:TYR:CE1	3.02	0.43
2:B:27:ILE:CD1	2:B:69:LEU:HD23	2.49	0.42
5:M:291:THR:HA	5:M:296:TYR:CD1	2.54	0.42
4:H:9:ALA:O	4:H:112:THR:O	2.38	0.42
3:L:2:ILE:HG22	3:L:4:LEU:HD22	2.01	0.42
5:M:221:LEU:HD22	5:M:302:THR:HG21	2.01	0.42
4:H:100:TYR:CD2	4:H:100:TYR:O	2.72	0.42
5:M:254:ARG:HG3	5:M:254:ARG:NH1	2.33	0.42
6:I:208:GLY:O	6:I:220:LEU:HG	2.20	0.42
2:B:38:CYS:CB	2:B:39:PRO:HD2	2.49	0.42
3:L:80:ASN:CB	3:L:81:PRO:CD	2.94	0.42
1:A:83:HIS:CE1	1:A:85:SER:HG	2.38	0.42
6:I:293:LEU:HD13	6:I:295:PHE:CZ	2.54	0.42
4:H:99:ARG:CG	4:H:99:ARG:NH1	2.81	0.42
6:I:238:ARG:HA	6:I:293:LEU:O	2.19	0.42
1:A:20:GLN:CD	4:H:33:TRP:CD1	2.91	0.42
1:A:43:SER:O	1:A:45:LYS:N	2.53	0.42
5:M:255:PHE:O	5:M:258:VAL:HG23	2.20	0.42
4:H:103:SER:C	4:H:104:PHE:HD1	2.23	0.42
4:H:45:LEU:H	4:H:45:LEU:HD22	1.85	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:104:PHE:CE1	4:H:45:LEU:HD23	2.54	0.42
6:I:301:THR:O	6:I:302:ALA:C	2.59	0.41
2:B:74:ARG:HD3	6:I:252:THR:HG23	2.02	0.41
3:L:87:VAL:O	3:L:88:ALA:HB2	2.20	0.41
2:B:44:VAL:O	2:B:44:VAL:HG12	2.20	0.41
5:M:278:VAL:HG23	5:M:282:ASP:OD2	2.20	0.41
3:L:46:GLN:HA	3:L:47:PRO:HD3	1.96	0.41
3:L:36:PHE:CE2	4:H:102:ASN:OD1	2.73	0.41
3:L:38:GLN:HB2	3:L:93:GLN:HB3	2.01	0.41
5:M:283:PHE:HE2	5:M:305:GLU:O	2.02	0.41
1:A:20:GLN:OE1	4:H:33:TRP:NE1	2.52	0.41
3:L:58:LEU:HD23	3:L:66:PHE:O	2.21	0.41
4:H:71:THR:CG2	4:H:72:VAL:N	2.83	0.41
6:I:235:SER:HB2	6:I:299:GLN:HE21	1.86	0.41
3:L:6:GLN:HG2	3:L:23:CYS:HB2	2.03	0.41
3:L:34:ASN:O	3:L:97:GLU:HB2	2.20	0.41
2:B:64:PHE:CE2	5:M:228:THR:HB	2.56	0.41
1:A:68:VAL:O	1:A:75:LYS:HB2	2.21	0.41
6:I:216:GLY:O	6:I:284:SER:O	2.38	0.41
6:I:264:VAL:HG12	6:I:268:PHE:HB2	2.03	0.41
4:H:51:ILE:HD13	4:H:52(A):PRO:HA	2.04	0.40
3:L:42:GLN:HA	3:L:48:PRO:HA	2.03	0.40
4:H:45:LEU:N	4:H:45:LEU:HD13	2.36	0.40
6:I:286:LEU:O	6:I:286:LEU:HD12	2.21	0.40
1:A:38:PRO:HD3	2:B:59:TYR:OH	2.21	0.40
2:B:12:ILE:HG13	2:B:13:ASN:N	2.36	0.40
1:A:20:GLN:HA	4:H:33:TRP:HE1	1.86	0.40
3:L:6:GLN:HE21	3:L:92:CYS:H	1.66	0.40
2:B:77:ASN:HA	2:B:78:PRO:HD3	1.86	0.40
6:I:299:GLN:OE1	6:I:305:PRO:O	2.40	0.40
4:H:47:TRP:CZ2	4:H:50:GLU:HG3	2.46	0.40
1:A:43:SER:O	1:A:46:THR:HG22	2.21	0.40
6:I:253:ASN:N	6:I:253:ASN:ND2	2.67	0.40
5:M:290:GLN:NE2	5:M:297:THR:H	2.20	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:46:THR:O	1:A:46:THR:O[6_765]	1.97	0.23

## 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	85/92 (92%)	73 (86%)	8 (9%)	4 (5%)	3	29
2	B	109/145 (75%)	84 (77%)	18 (16%)	7 (6%)	2	20
3	L	108/114 (95%)	81 (75%)	22 (20%)	5 (5%)	3	29
4	H	108/117 (92%)	80 (74%)	22 (20%)	6 (6%)	2	24
5	M	106/108 (98%)	91 (86%)	12 (11%)	3 (3%)	6	43
6	I	119/122 (98%)	90 (76%)	24 (20%)	5 (4%)	3	32
All	All	635/698 (91%)	499 (79%)	106 (17%)	30 (5%)	3	29

All (30) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	73	GLY
2	B	4	PRO
4	H	16	ALA
4	H	33	TRP
5	M	268	ALA
6	I	311	VAL
1	A	50	GLN
2	B	3	GLU
2	B	79	VAL
2	B	105	ASP
3	L	65	ARG
4	H	27	TYR
4	H	28	THR
5	M	260	ASP
5	M	280	ALA
6	I	300	GLY
1	A	51	LYS
1	A	71	MET
2	B	49	LEU
2	B	108	LEU

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Mol	Chain	Res	Type
3	L	26	SER
3	L	81	PRO
4	H	52	ASN
6	I	262	GLY
3	L	33	GLY
4	H	101	GLY
3	L	63	PRO
2	B	103	PRO
6	I	202	VAL
6	I	254	VAL

### 5.3.2 Protein sidechains [i](#)

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	78/83 (94%)	67 (86%)	11 (14%)	4	24
2	B	93/129 (72%)	70 (75%)	23 (25%)	1	5
3	L	89/98 (91%)	73 (82%)	16 (18%)	2	12
4	H	82/96 (85%)	63 (77%)	19 (23%)	1	5
5	M	93/94 (99%)	80 (86%)	13 (14%)	4	24
6	I	99/101 (98%)	86 (87%)	13 (13%)	5	27
All	All	534/601 (89%)	439 (82%)	95 (18%)	2	13

All (95) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	CYS
1	A	41	LEU
1	A	50	GLN
1	A	66	ASN
1	A	69	THR
1	A	71	MET
1	A	83	HIS
1	A	86	THR

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Mol	Chain	Res	Type
1	A	87	CYS
1	A	90	HIS
1	A	91	LYS
2	B	5	LEU
2	B	23	CYS
2	B	25	VAL
2	B	30	ASN
2	B	31	THR
2	B	38	CYS
2	B	42	THR
2	B	45	LEU
2	B	46	GLN
2	B	49	LEU
2	B	52	LEU
2	B	58	ASN
2	B	67	ILE
2	B	82	TYR
2	B	90	CYS
2	B	95	ARG
2	B	100	CYS
2	B	104	LYS
2	B	105	ASP
2	B	106	HIS
2	B	108	LEU
2	B	109	THR
2	B	111	ASP
3	L	9	ASP
3	L	13	VAL
3	L	18	ARG
3	L	21	ILE
3	L	23	CYS
3	L	26	SER
3	L	32	TYR
3	L	42	GLN
3	L	52	ILE
3	L	60	SER
3	L	67	SER
3	L	75	PHE
3	L	77	LEU
3	L	89	THR
3	L	97	GLU
3	L	110	LEU

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Mol	Chain	Res	Type
4	H	5	GLN
4	H	6	GLN
4	H	7	SER
4	H	11	LEU
4	H	20	LEU
4	H	26	ASP
4	H	27	TYR
4	H	28	THR
4	H	38	LYS
4	H	45	LEU
4	H	51	ILE
4	H	61	ASN
4	H	69	THR
4	H	70	LEU
4	H	98	ARG
4	H	99	ARG
4	H	104	PHE
4	H	106	TYR
4	H	115	VAL
5	M	204	LEU
5	M	226	SER
5	M	228	THR
5	M	230	ASP
5	M	234	SER
5	M	244	PRO
5	M	274	THR
5	M	276	SER
5	M	278	VAL
5	M	279	GLN
5	M	281	GLU
5	M	290	GLN
5	M	306	ILE
6	I	201	GLN
6	I	210	HIS
6	I	211	LEU
6	I	213	LYS
6	I	238	ARG
6	I	253	ASN
6	I	269	SER
6	I	280	ASN
6	I	286	LEU
6	I	287	ARG

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Mol	Chain	Res	Type
6	I	293	LEU
6	I	298	ARG
6	I	310	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	20	GLN
1	A	27	GLN
1	A	90	HIS
2	B	30	ASN
2	B	46	GLN
3	L	6	GLN
3	L	34	ASN
3	L	38	GLN
3	L	41	GLN
3	L	57	ASN
4	H	5	GLN
4	H	35	HIS
4	H	82	GLN
5	M	206	GLN
5	M	238	GLN
5	M	290	GLN
6	I	253	ASN
6	I	299	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry

2 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	$\# Z  > 2$
7	NAG	A	93	1	14,14,15	0.61	0	15,19,21	0.96	1 (6%)
7	NAG	A	94	1	14,14,15	0.38	0	15,19,21	1.08	1 (6%)

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
7	NAG	A	93	1	-	0/6/23/26	0/1/1/1
7	NAG	A	94	1	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	A	94	NAG	C4-C3-C2	-2.82	106.85	111.23
7	A	93	NAG	C4-C3-C2	-2.40	107.50	111.23

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	A	93	NAG	1	0
7	A	94	NAG	2	0

## 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	87/92 (94%)	-0.04	1 (1%) 82 73	17, 17, 17, 17	5 (5%)
2	B	111/145 (76%)	0.20	7 (6%) 23 18	17, 17, 17, 20	1 (0%)
3	L	112/114 (98%)	0.22	1 (0%) 85 78	20, 44, 44, 44	0
4	H	112/117 (95%)	0.32	4 (3%) 46 37	15, 44, 44, 44	0
5	M	108/108 (100%)	-0.08	0 100 100	15, 15, 15, 15	6 (5%)
6	I	120/122 (98%)	0.06	1 (0%) 87 80	15, 15, 15, 15	4 (3%)
All	All	650/698 (93%)	0.12	14 (2%) 65 55	15, 17, 44, 44	16 (2%)

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	105	ASP	3.7
2	B	47	GLY	3.3
3	L	1	ASP	3.1
2	B	48	VAL	3.0
2	B	4	PRO	2.9
4	H	16	ALA	2.7
2	B	104	LYS	2.4
2	B	3	GLU	2.2
4	H	14	PRO	2.1
6	I	321	SER	2.0
2	B	46	GLN	2.0
4	H	44	GLY	2.0
1	A	6	ASP	2.0
4	H	54	ASN	2.0

### 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 [i](#)

There are no carbohydrates in this entry.

### 6.4 Ligands [i](#)

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
7	NAG	A	93	14/15	0.87	0.30	0.33	34,34,34,34	0
7	NAG	A	94	14/15	0.82	0.26	-0.01	32,32,32,32	0

### 6.5 Other polymers [i](#)

There are no such residues in this entry.