



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:54 PM UTC

PDB ID : 3G6A / pdb_00003g6a
Title : Crystal structure of anti-IL-13 antibody CNTO607
Authors : Teplyakov, A.; Obmolova, G.; Gilliland, G.L.
Deposited on : 2009-02-06
Resolution : 2.10 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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

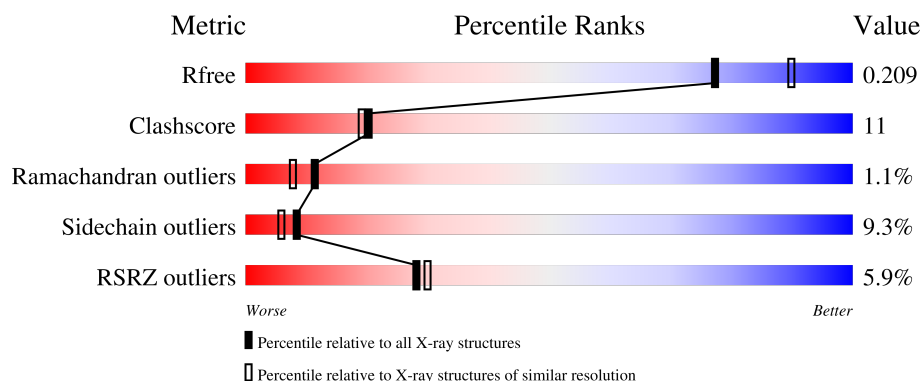
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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}	180053	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 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	213	10% 62% 31% 6%
1	L	213	4% 74% 21%
2	B	229	7% 72% 20%
2	H	229	3% 85% 10%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6738 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 CNTO607 Fab Light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	210	Total	C	N	O	S	0	0	0
			1568	978	260	325	5			
1	A	210	Total	C	N	O	S	0	0	0
			1568	978	260	325	5			

- Molecule 2 is a protein called CNTO607 Fab Heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	222	Total	C	N	O	S	0	0	0
			1666	1053	281	326	6			
2	B	222	Total	C	N	O	S	0	0	0
			1666	1053	281	326	6			

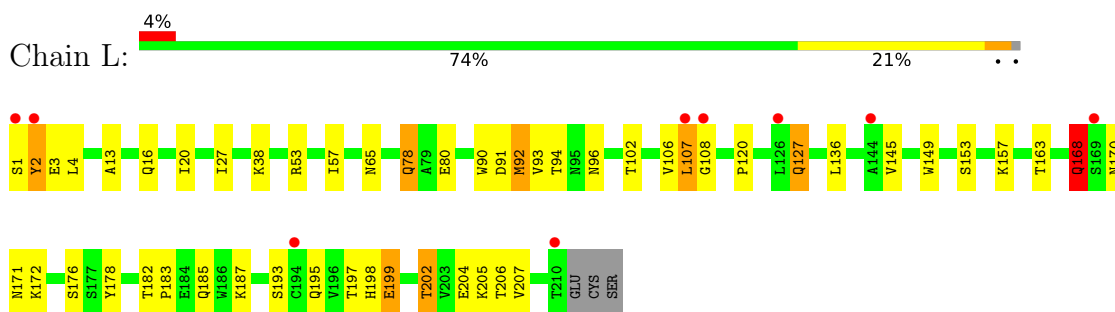
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	L	57	Total	O	0	0
			57	57		
3	H	101	Total	O	0	0
			101	101		
3	A	37	Total	O	0	0
			37	37		
3	B	75	Total	O	0	0
			75	75		

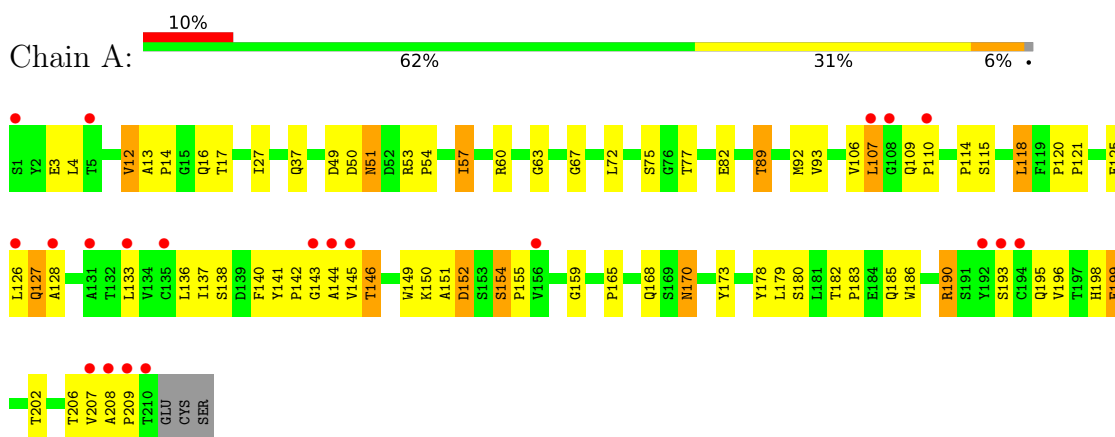
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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.

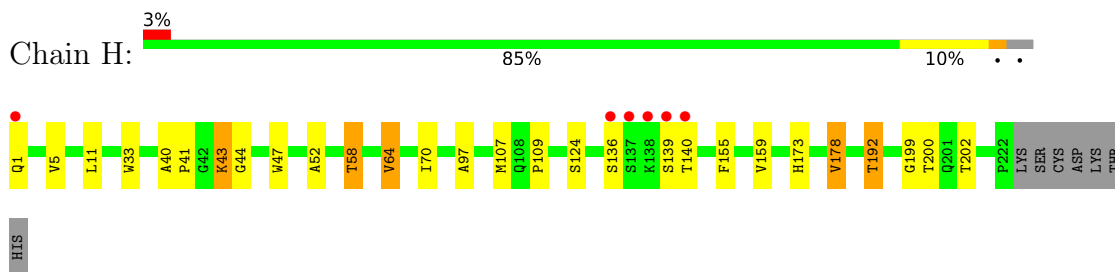
- Molecule 1: CNTO607 Fab Light chain



- Molecule 1: CNTO607 Fab Light chain

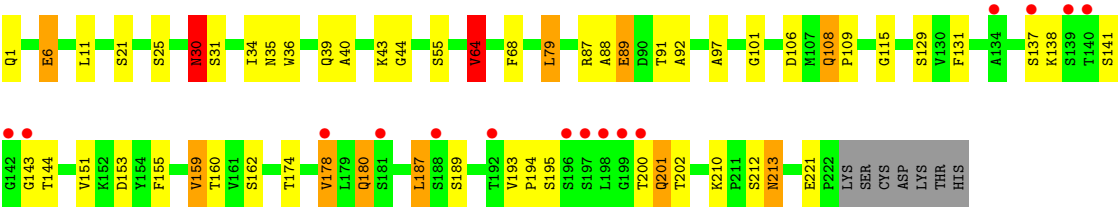


- Molecule 2: CNTO607 Fab Heavy chain



- Molecule 2: CNTO607 Fab Heavy chain





4 Data and refinement statistics

Property	Value	Source
Space group	P 42	Depositor
Cell constants a, b, c, α , β , γ	108.05Å 108.05Å 78.56Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.00 – 2.10 15.00 – 2.10	Depositor EDS
% Data completeness (in resolution range)	94.7 (15.00-2.10) 94.5 (15.00-2.10)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$	-	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.208 , 0.283 0.211 , 0.209	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	(Not available)	Xtriage
Anisotropy	(Not available)	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 56.8	EDS
L-test for twinning ¹	$\langle L \rangle =$ (Not available), $\langle L^2 \rangle =$ (Not available)	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6738	wwPDB-VP
Average B, all atoms (Å ²)	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *(Not available)*

¹Theoretical 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

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	0/1607	1.09	5/2200 (0.2%)
1	L	0.95	0/1607	1.07	1/2200 (0.0%)
2	B	1.04	1/1709 (0.1%)	1.21	7/2332 (0.3%)
2	H	1.12	1/1709 (0.1%)	1.18	3/2332 (0.1%)
All	All	1.00	2/6632 (0.0%)	1.14	16/9064 (0.2%)

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	L	0	1
2	H	0	1
All	All	0	2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	5	VAL	CA-CB	5.14	1.60	1.53
2	B	115	GLY	N-CA	5.05	1.50	1.45

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	64	VAL	N-CA-CB	-10.07	100.34	112.33
2	B	108	GLN	CA-C-N	9.13	129.16	119.76
2	B	108	GLN	C-N-CA	9.13	129.16	119.76
1	A	208	ALA	CA-C-N	7.23	128.88	119.84
1	A	208	ALA	C-N-CA	7.23	128.88	119.84
2	B	64	VAL	N-CA-CB	-7.07	103.92	112.33
1	A	154	SER	CA-C-N	6.85	126.83	119.78
1	A	154	SER	C-N-CA	6.85	126.83	119.78

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	212	SER	N-CA-C	-5.91	105.65	112.92
1	L	2	TYR	N-CA-C	5.89	119.00	107.57
2	B	6	GLU	CA-CB-CG	5.40	124.91	114.10
2	B	213	ASN	N-CA-C	5.15	118.55	111.54
2	H	178	VAL	CB-CA-C	-5.13	104.55	110.91
2	H	200	THR	N-CA-CB	5.07	117.34	110.59
2	B	35	ASN	CB-CA-C	-5.03	99.42	109.33
1	A	138	SER	N-CA-C	5.01	116.80	109.24

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	H	199	GLY	Peptide
1	L	1	SER	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	1568	0	1499	54	0
1	L	1568	0	1499	40	0
2	B	1666	0	1618	38	0
2	H	1666	0	1618	25	0
3	A	37	0	0	1	0
3	B	75	0	0	0	0
3	H	101	0	0	3	0
3	L	57	0	0	2	0
All	All	6738	0	6234	144	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 11.

All (144) 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:L:20:ILE:HG21	1:L:102:THR:HG21	1.34	1.08
2:H:40:ALA:H	2:H:43:LYS:HE2	1.05	1.06
1:L:20:ILE:CG2	1:L:102:THR:HG21	1.87	1.05
2:H:40:ALA:N	2:H:43:LYS:HE2	1.89	0.87
1:A:193:SER:CB	1:A:206:THR:HG22	2.06	0.86
2:H:43:LYS:HG2	2:H:44:GLY:N	1.92	0.84
1:L:120:PRO:HB3	1:L:207:VAL:HG21	1.60	0.83
1:L:178:TYR:CE1	2:H:178:VAL:HG21	2.16	0.81
1:A:118:LEU:HD13	1:A:207:VAL:HG13	1.62	0.81
1:A:182:THR:HG23	1:A:185:GLN:H	1.47	0.79
1:A:193:SER:HB3	1:A:206:THR:HG22	1.64	0.79
1:A:178:TYR:CE1	2:B:178:VAL:HG11	2.18	0.78
1:L:182:THR:H	1:L:185:GLN:HE21	1.31	0.78
2:B:43:LYS:HE3	2:B:44:GLY:O	1.84	0.77
1:L:193:SER:HB3	1:L:206:THR:HG22	1.68	0.75
2:B:143:GLY:O	2:B:195:SER:N	2.14	0.74
2:H:58:THR:HG21	3:H:244:HOH:O	1.90	0.72
1:L:20:ILE:HG21	1:L:102:THR:CG2	2.19	0.70
1:L:168:GLN:HB2	1:L:172:LYS:O	1.92	0.70
1:A:16:GLN:O	1:A:77:THR:HG23	1.92	0.69
1:A:178:TYR:CD1	2:B:178:VAL:HG11	2.28	0.69
1:L:20:ILE:HG23	1:L:102:THR:HG21	1.73	0.68
1:L:178:TYR:CD1	2:H:178:VAL:HG21	2.29	0.68
2:H:43:LYS:HE3	2:H:44:GLY:O	1.93	0.68
2:B:201:GLN:HE21	2:B:201:GLN:HA	1.59	0.67
1:A:168:GLN:HE21	1:A:170:ASN:HD21	1.42	0.67
1:A:12:VAL:HG13	1:A:16:GLN:HB2	1.77	0.67
1:A:193:SER:HB2	1:A:206:THR:HG22	1.76	0.65
1:L:198:HIS:O	1:L:199:GLU:C	2.41	0.64
1:L:178:TYR:CE1	2:H:178:VAL:CG2	2.80	0.64
2:B:43:LYS:HG2	2:B:44:GLY:H	1.61	0.64
1:L:199:GLU:O	1:L:199:GLU:HG3	1.98	0.63
2:H:139:SER:O	2:H:140:THR:HB	1.99	0.63
2:B:87:ARG:HG2	2:B:89:GLU:HG2	1.79	0.63
1:A:51:ASN:C	1:A:51:ASN:HD22	2.07	0.62
2:H:40:ALA:HB3	2:H:43:LYS:HD3	1.82	0.61
1:A:17:THR:HG22	1:A:75:SER:HA	1.82	0.61
1:A:127:GLN:HG2	1:A:127:GLN:O	2.01	0.60
2:B:40:ALA:H	2:B:43:LYS:NZ	2.00	0.60
1:A:106:VAL:O	1:A:107:LEU:HG	2.03	0.59
1:L:106:VAL:O	1:L:107:LEU:C	2.46	0.59
1:A:107:LEU:HB2	1:A:109:GLN:HG2	1.85	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:91:ASP:OD1	1:L:94:THR:OG1	2.22	0.58
1:A:37:GLN:HE22	2:B:39:GLN:HE22	1.52	0.58
1:L:92:MET:HG2	3:L:242:HOH:O	2.04	0.57
1:L:193:SER:CB	1:L:206:THR:HG22	2.33	0.57
1:L:163:THR:HG23	2:H:178:VAL:HG22	1.87	0.57
2:B:144:THR:HA	2:B:193:VAL:O	2.05	0.57
2:B:141:SER:HB3	2:B:144:THR:OG1	2.06	0.56
2:H:192:THR:HG23	3:H:263:HOH:O	2.06	0.56
2:B:143:GLY:O	2:B:194:PRO:HA	2.06	0.55
1:L:127:GLN:O	1:L:127:GLN:HG3	2.07	0.55
1:A:182:THR:OG1	1:A:183:PRO:HD2	2.07	0.54
2:B:151:VAL:HG11	2:B:159:VAL:HG11	1.88	0.54
1:L:4:LEU:HD22	1:L:27:ILE:HD12	1.89	0.54
1:L:78:GLN:HE21	1:L:80:GLU:H	1.54	0.54
1:L:178:TYR:HE1	2:H:178:VAL:CG2	2.21	0.53
1:A:190:ARG:HB2	3:A:263:HOH:O	2.08	0.53
1:L:195:GLN:HG2	1:L:204:GLU:HG3	1.90	0.53
1:A:178:TYR:CE1	2:B:178:VAL:CG1	2.92	0.53
1:A:198:HIS:O	1:A:199:GLU:C	2.52	0.53
1:A:114:PRO:HB3	1:A:140:PHE:CD2	2.44	0.52
2:H:43:LYS:HG2	2:H:44:GLY:H	1.69	0.52
1:A:137:ILE:HG22	1:A:140:PHE:CE2	2.45	0.51
2:H:33:TRP:CE3	2:H:52:ALA:HA	2.46	0.51
2:B:187:LEU:HD12	2:B:187:LEU:C	2.36	0.50
2:B:91:THR:O	2:B:92:ALA:HB2	2.12	0.50
2:B:87:ARG:CG	2:B:89:GLU:HG2	2.42	0.50
1:A:182:THR:H	1:A:185:GLN:HE21	1.60	0.49
1:L:27:ILE:HG23	1:L:65:ASN:HD21	1.77	0.49
1:A:126:LEU:C	1:A:128:ALA:H	2.20	0.49
1:A:110:PRO:HD2	1:A:141:TYR:CE1	2.48	0.49
1:L:38:LYS:NZ	1:L:80:GLU:HG2	2.28	0.48
1:L:120:PRO:HB3	1:L:207:VAL:CG2	2.36	0.48
1:L:197:THR:OG1	1:L:202:THR:OG1	2.29	0.48
1:A:140:PHE:CE1	1:A:173:TYR:HB2	2.49	0.48
2:B:153:ASP:OD1	2:B:180:GLN:NE2	2.46	0.48
2:B:178:VAL:CG2	2:B:180:GLN:HG2	2.43	0.48
2:B:174:THR:HA	2:B:189:SER:HA	1.96	0.47
1:L:90:TRP:CD1	2:H:107:MET:HE3	2.50	0.47
1:A:168:GLN:NE2	1:A:170:ASN:HD21	2.11	0.47
1:L:53:ARG:HD2	1:L:57:ILE:O	2.15	0.47
1:A:133:LEU:HD13	1:A:179:LEU:HD23	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:160:THR:HG21	2:B:210:LYS:NZ	2.29	0.47
1:L:90:TRP:NE1	2:H:107:MET:HE3	2.30	0.46
1:A:133:LEU:HD22	1:A:149:TRP:HZ3	1.80	0.46
1:L:13:ALA:HB3	1:L:16:GLN:HG3	1.97	0.46
1:L:183:PRO:O	1:L:187:LYS:HG2	2.16	0.46
1:A:13:ALA:H	1:A:16:GLN:NE2	2.13	0.46
1:A:127:GLN:O	1:A:127:GLN:CG	2.64	0.46
2:H:11:LEU:HD11	2:H:155:PHE:HZ	1.81	0.45
1:A:53:ARG:HD3	1:A:57:ILE:HG22	1.98	0.45
2:B:11:LEU:HD13	2:B:155:PHE:HZ	1.80	0.45
1:L:168:GLN:NE2	1:L:170:ASN:OD1	2.45	0.45
1:A:110:PRO:HD2	1:A:141:TYR:HE1	1.81	0.45
1:A:137:ILE:HG22	1:A:140:PHE:HE2	1.80	0.45
2:H:58:THR:OG1	2:H:70:ILE:CG2	2.65	0.44
1:A:125:GLU:HG3	2:B:131:PHE:CE1	2.52	0.44
1:A:142:PRO:O	1:A:198:HIS:NE2	2.50	0.44
1:A:151:ALA:O	1:A:152:ASP:C	2.61	0.44
2:B:201:GLN:HE21	2:B:201:GLN:CA	2.27	0.44
2:H:43:LYS:CG	2:H:44:GLY:N	2.72	0.44
1:A:14:PRO:HD3	1:A:107:LEU:HB3	2.00	0.44
2:B:64:VAL:HG13	2:B:68:PHE:HB2	2.00	0.44
1:A:186:TRP:CZ2	1:A:209:PRO:HA	2.52	0.44
1:A:154:SER:HA	1:A:155:PRO:HD3	1.76	0.43
2:H:43:LYS:HG3	3:H:308:HOH:O	2.17	0.43
2:B:108:GLN:HA	2:B:109:PRO:HD3	1.75	0.43
2:B:160:THR:HG21	2:B:210:LYS:HZ1	1.84	0.43
1:A:146:THR:O	1:A:196:VAL:HA	2.19	0.43
1:L:78:GLN:HG3	1:L:80:GLU:H	1.84	0.43
1:A:49:ASP:O	1:A:50:ASP:HB2	2.19	0.43
1:A:159:GLY:O	1:A:179:LEU:HA	2.19	0.43
2:B:43:LYS:HB3	2:B:43:LYS:HE2	1.52	0.42
1:A:51:ASN:HB3	1:A:63:GLY:O	2.19	0.42
1:A:53:ARG:CD	1:A:57:ILE:HG22	2.49	0.42
1:A:143:GLY:O	1:A:144:ALA:C	2.63	0.42
1:L:106:VAL:O	1:L:108:GLY:N	2.53	0.42
1:A:4:LEU:HD11	1:A:89:THR:HB	2.00	0.42
2:B:40:ALA:H	2:B:43:LYS:HZ3	1.68	0.42
2:B:97:ALA:HB1	2:B:109:PRO:HB3	2.01	0.42
1:L:93:VAL:HG12	1:L:94:THR:HG23	2.01	0.42
1:L:53:ARG:HG3	1:L:57:ILE:HB	2.01	0.42
1:A:53:ARG:HA	1:A:54:PRO:HD3	1.93	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:140:PHE:CD1	1:A:173:TYR:HB2	2.54	0.42
2:B:36:TRP:NE1	2:B:79:LEU:HD13	2.34	0.42
1:L:170:ASN:O	1:L:171:ASN:HB2	2.20	0.41
1:A:152:ASP:C	1:A:154:SER:H	2.28	0.41
1:A:51:ASN:C	1:A:51:ASN:ND2	2.76	0.41
1:A:82:GLU:HG2	1:A:106:VAL:HG23	2.02	0.41
1:L:149:TRP:HA	1:L:193:SER:O	2.21	0.41
2:H:97:ALA:HB1	2:H:109:PRO:HB3	2.03	0.41
2:B:87:ARG:CD	2:B:89:GLU:HG2	2.50	0.41
3:L:225:HOH:O	2:H:173:HIS:HD2	2.03	0.41
1:A:114:PRO:HB3	1:A:140:PHE:HD2	1.84	0.41
2:B:178:VAL:HG23	2:B:180:GLN:HG2	2.03	0.41
2:B:201:GLN:HE21	2:B:202:THR:H	1.69	0.41
2:B:30:ASN:HD22	2:B:30:ASN:HA	1.61	0.41
1:A:120:PRO:HB2	1:A:121:PRO:HD2	2.01	0.40
2:B:101:GLY:N	2:B:106:ASP:HB3	2.36	0.40
2:B:87:ARG:O	2:B:88:ALA:C	2.63	0.40
1:L:96:ASN:HD22	2:H:47:TRP:CD1	2.40	0.40
2:H:40:ALA:O	2:H:41:PRO:C	2.62	0.40
2:B:68:PHE:N	2:B:68:PHE:CD1	2.90	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	208/213 (98%)	190 (91%)	13 (6%)	5 (2%)	4	2
1	L	208/213 (98%)	196 (94%)	9 (4%)	3 (1%)	9	5
2	B	220/229 (96%)	209 (95%)	10 (4%)	1 (0%)	24	22
2	H	220/229 (96%)	211 (96%)	9 (4%)	0	100	100
All	All	856/884 (97%)	806 (94%)	41 (5%)	9 (1%)	11	8

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	199	GLU
1	L	107	LEU
1	L	199	GLU
1	A	127	GLN
1	A	152	ASP
1	L	168	GLN
1	A	165	PRO
2	B	30	ASN
1	A	67	GLY

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	176/179 (98%)	154 (88%)	22 (12%)	4	2
1	L	176/179 (98%)	163 (93%)	13 (7%)	13	10
2	B	185/192 (96%)	162 (88%)	23 (12%)	4	2
2	H	185/192 (96%)	176 (95%)	9 (5%)	22	22
All	All	722/742 (97%)	655 (91%)	67 (9%)	8	6

All (67) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	2	TYR
1	L	3	GLU
1	L	78	GLN
1	L	92	MET
1	L	127	GLN
1	L	136	LEU
1	L	145	VAL
1	L	153	SER
1	L	157	LYS
1	L	168	GLN
1	L	176	SER

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Mol	Chain	Res	Type
1	L	202	THR
1	L	205	LYS
2	H	1	GLN
2	H	43	LYS
2	H	58	THR
2	H	64	VAL
2	H	124	SER
2	H	136	SER
2	H	159	VAL
2	H	192	THR
2	H	202	THR
1	A	3	GLU
1	A	12	VAL
1	A	27	ILE
1	A	51	ASN
1	A	57	ILE
1	A	60	ARG
1	A	72	LEU
1	A	89	THR
1	A	92	MET
1	A	93	VAL
1	A	107	LEU
1	A	115	SER
1	A	118	LEU
1	A	136	LEU
1	A	145	VAL
1	A	146	THR
1	A	150	LYS
1	A	170	ASN
1	A	180	SER
1	A	190	ARG
1	A	195	GLN
1	A	202	THR
2	B	1	GLN
2	B	6	GLU
2	B	21	SER
2	B	25	SER
2	B	30	ASN
2	B	31	SER
2	B	34	ILE
2	B	55	SER
2	B	64	VAL

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Mol	Chain	Res	Type
2	B	79	LEU
2	B	89	GLU
2	B	129	SER
2	B	137	SER
2	B	138	LYS
2	B	159	VAL
2	B	162	SER
2	B	178	VAL
2	B	180	GLN
2	B	187	LEU
2	B	200	THR
2	B	201	GLN
2	B	213	ASN
2	B	221	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	37	GLN
1	L	65	ASN
1	L	78	GLN
1	L	185	GLN
2	H	39	GLN
2	H	82	GLN
2	H	108	GLN
1	A	16	GLN
1	A	41	GLN
1	A	51	ASN
1	A	127	GLN
1	A	170	ASN
1	A	185	GLN
1	A	195	GLN
2	B	1	GLN
2	B	30	ASN
2	B	39	GLN
2	B	108	GLN
2	B	180	GLN
2	B	201	GLN
2	B	208	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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, 95th 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 ²)	Q<0.9
1	A	210/213 (98%)	0.73	21 (10%) 12 13	37, 65, 95, 109	0
1	L	210/213 (98%)	0.32	9 (4%) 40 41	33, 52, 88, 108	0
2	B	222/229 (96%)	0.29	15 (6%) 23 25	32, 46, 83, 132	0
2	H	222/229 (96%)	-0.13	6 (2%) 56 59	29, 41, 61, 125	0
All	All	864/884 (97%)	0.30	51 (5%) 28 30	29, 50, 88, 132	0

All (51) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	H	139	SER	5.0
1	A	108	GLY	4.0
2	B	197	SER	4.0
2	B	143	GLY	3.8
2	B	142	GLY	3.8
2	B	137	SER	3.7
2	B	139	SER	3.6
2	B	200	THR	3.3
2	H	137	SER	3.3
1	A	143	GLY	3.3
1	A	194	CYS	3.2
1	A	210	THR	3.2
1	A	192	TYR	3.2
1	L	2	TYR	3.1
1	L	108	GLY	3.1
2	B	178	VAL	3.1
1	A	110	PRO	3.0
2	B	134	ALA	3.0
1	A	156	VAL	2.9
1	L	144	ALA	2.7
1	A	131	ALA	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	193	SER	2.7
1	A	133	LEU	2.6
2	B	140	THR	2.6
1	L	169	SER	2.5
2	B	198	LEU	2.5
1	A	126	LEU	2.5
1	A	144	ALA	2.5
1	A	107	LEU	2.4
1	A	5	THR	2.4
2	B	188	SER	2.4
1	A	209	PRO	2.4
2	H	1	GLN	2.3
2	B	199	GLY	2.3
1	A	135	CYS	2.3
1	A	145	VAL	2.3
1	L	107	LEU	2.3
2	B	192	THR	2.2
1	A	208	ALA	2.2
1	A	207	VAL	2.2
1	A	128	ALA	2.2
1	L	1	SER	2.2
1	A	1	SER	2.1
2	B	181	SER	2.1
1	L	210	THR	2.1
2	H	140	THR	2.1
2	B	196	SER	2.1
1	L	126	LEU	2.1
2	H	138	LYS	2.1
2	H	136	SER	2.0
1	L	194	CYS	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 ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.