



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 06:51 am GMT

PDB ID : 3G6A
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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.9-1692
EDS	:	trunk28620
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
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)	:	recalc28949

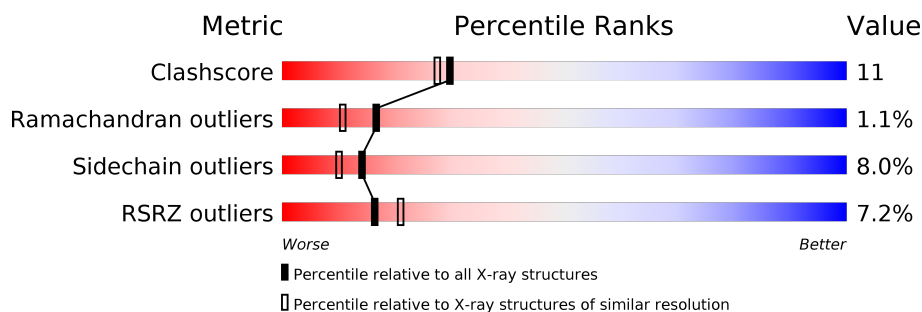
1 Overall quality at a glance

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))
Clashscore	112137	4788 (2.10-2.10)
Ramachandran outliers	110173	4740 (2.10-2.10)
Sidechain outliers	110143	4741 (2.10-2.10)
RSRZ outliers	101464	4275 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 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	213	<div> <div>14%</div> <div>66%</div> <div>29%</div> <div>..</div> </div>
1	L	213	<div> <div>8%</div> <div>76%</div> <div>20%</div> <div>..</div> </div>
2	B	229	<div> <div>4%</div> <div>73%</div> <div>20%</div> <div>...</div> </div>
2	H	229	<div> <div>3%</div> <div>85%</div> <div>10%</div> <div>..</div> </div>

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			

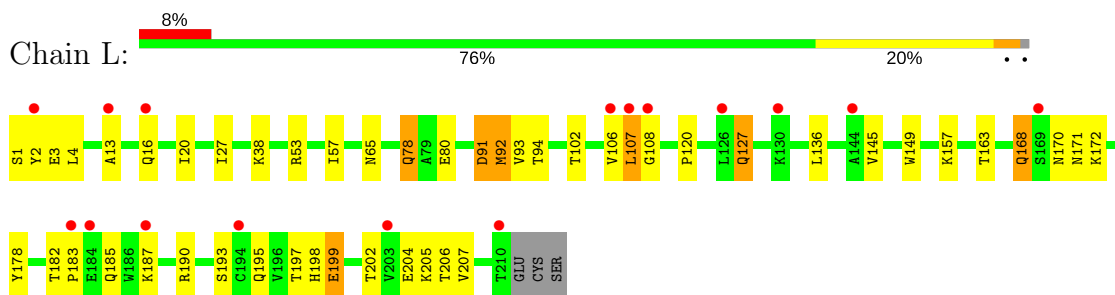
- Molecule 3 is water.

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

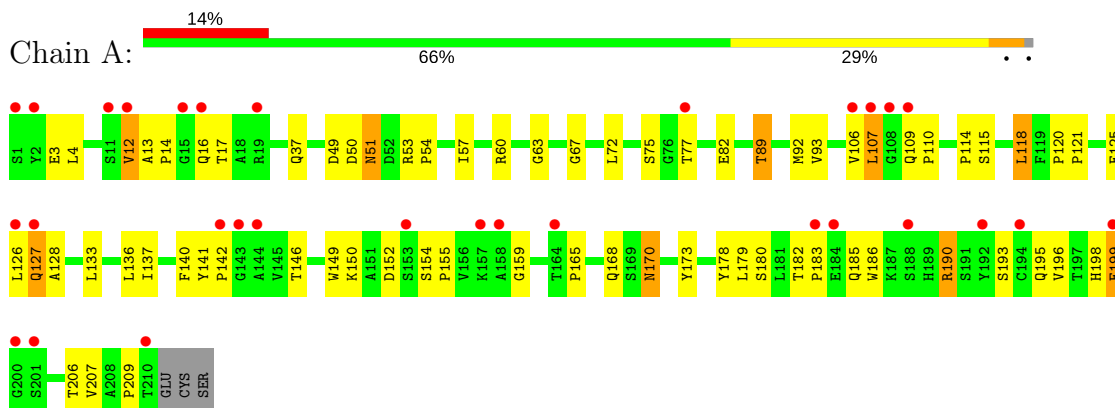
3 Residue-property plots [i](#)

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.

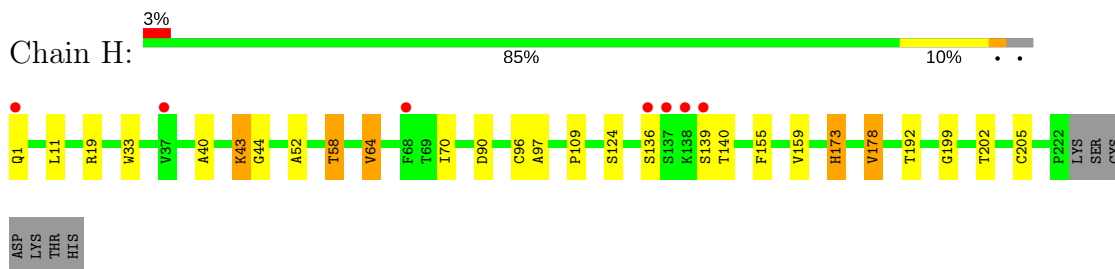
- Molecule 1: CNTO607 Fab Light chain



- Molecule 1: CNTO607 Fab Light chain

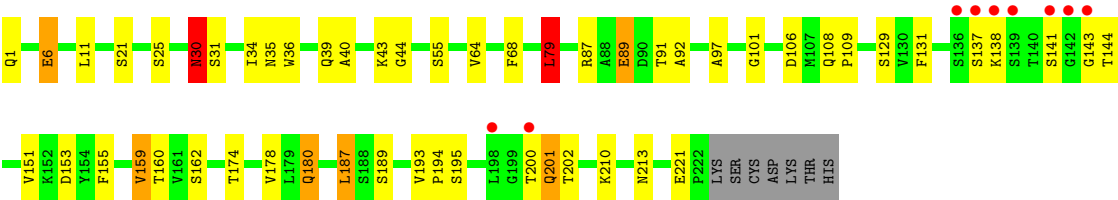


- Molecule 2: CNTO607 Fab Heavy chain



- 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 36.92 – 2.00	Depositor EDS
% Data completeness (in resolution range)	94.7 (15.00-2.10) 93.7 (36.92-2.00)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.63 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.208 , 0.283 0.211 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	35.7	Xtriage
Anisotropy	0.237	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 57.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.031 for h,-k,-l	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: *The largest off-origin peak in the Patterson function is 14.87% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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	0.69	0/1607	0.78	1/2200 (0.0%)
1	L	0.73	0/1607	0.80	0/2200
2	B	0.84	0/1709	0.93	3/2332 (0.1%)
2	H	0.92	3/1709 (0.2%)	0.90	3/2332 (0.1%)
All	All	0.80	3/6632 (0.0%)	0.85	7/9064 (0.1%)

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 (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	205	CYS	CB-SG	-7.34	1.69	1.82
2	H	96	CYS	CB-SG	-6.54	1.71	1.82
2	H	178	VAL	CB-CG1	5.16	1.63	1.52

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	79	LEU	CA-CB-CG	6.99	131.38	115.30
1	A	107	LEU	CA-CB-CG	5.94	128.96	115.30
2	H	90	ASP	CB-CG-OD1	5.61	123.34	118.30
2	B	35	ASN	CB-CA-C	-5.49	99.42	110.40
2	H	64	VAL	CG1-CB-CG2	5.41	119.56	110.90
2	B	6	GLU	CA-CB-CG	5.23	124.91	113.40
2	H	64	VAL	N-CA-CB	-5.07	100.34	111.50

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	50	0
1	L	1568	0	1499	38	0
2	B	1666	0	1618	37	0
2	H	1666	0	1618	21	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	136	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 (136) 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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:HA	2:B:201:GLN:HE21	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:178:TYR:CE1	2:H:178:VAL:CG2	2.80	0.64
1:L:199:GLU:HG3	1:L:199:GLU:O	1.98	0.64
2:B:43:LYS:HG2	2:B:44:GLY:H	1.61	0.64
2:B:87:ARG:HG2	2:B:89:GLU:HG2	1.79	0.63
2:H:139:SER:O	2:H:140:THR:HB	1.99	0.63
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:O	1:A:127:GLN:HG2	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:198:HIS:O	1:L:199:GLU:C	2.41	0.59
1:A:107:LEU:HB2	1:A:109:GLN:HG2	1.85	0.59
1:L:91:ASP:OD1	1:L:94:THR:OG1	2.22	0.58
1:A:51:ASN:C	1:A:51:ASN:HD22	2.07	0.58
1:A:37:GLN:HE22	2:B:39:GLN:HE22	1.52	0.58
1:L:193:SER:CB	1:L:206:THR:HG22	2.33	0.57
1:L:92:MET:HG2	3:L:242:HOH:O	2.04	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:151:VAL:HG11	2:B:159:VAL:HG11	1.88	0.54
1:L:106:VAL:O	1:L:107:LEU:C	2.46	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: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: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:110:PRO:HD2	1:A:141:TYR:CE1	2.48	0.49
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:L:38:LYS:NZ	1:L:80:GLU:HG2	2.28	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
1:A:198:HIS:O	1:A:199:GLU:C	2.52	0.48
2:B:174:THR:HA	2:B:189:SER:HA	1.96	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
2:B:160:THR:HG21	2:B:210:LYS:NZ	2.29	0.47
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
2:B:187:LEU:C	2:B:187:LEU:HD12	2.36	0.46
1:A:127:GLN:CG	1:A:127:GLN:O	2.64	0.46
1:A:13:ALA:H	1:A:16:GLN:NE2	2.13	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:A:110:PRO:HD2	1:A:141:TYR:HE1	1.81	0.45
1:L:168:GLN:NE2	1:L:170:ASN:OD1	2.45	0.45
1:A:137:ILE:HG22	1:A:140:PHE:HE2	1.80	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
2:B:201:GLN:CA	2:B:201:GLN:HE21	2.27	0.44
1:A:126:LEU:C	1:A:128:ALA:H	2.20	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:B:108:GLN:HA	2:B:109:PRO:HD3	1.75	0.43
2:H:43:LYS:HG3	3:H:308:HOH:O	2.17	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:159:GLY:O	1:A:179:LEU:HA	2.19	0.43
1:A:49:ASP:O	1:A:50:ASP:HB2	2.19	0.43
2:B:43:LYS:HE2	2:B:43:LYS:HB3	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:4:LEU:HD11	1:A:89:THR:HB	2.00	0.42
1:L:106:VAL:O	1:L:108:GLY:N	2.53	0.42
2:B:97:ALA:HB1	2:B:109:PRO:HB3	2.01	0.42
2:B:40:ALA:H	2:B:43:LYS:HZ3	1.68	0.42
1:L:93:VAL:HG12	1:L:94:THR:HG23	2.01	0.42
1:A:140:PHE:CD1	1:A:173:TYR:HB2	2.54	0.42
1:A:53:ARG:HA	1:A:54:PRO:HD3	1.93	0.42
2:B:36:TRP:NE1	2:B:79:LEU:HD13	2.34	0.42
1:L:53:ARG:HG3	1:L:57:ILE:HB	2.01	0.42
1:L:170:ASN:O	1:L:171:ASN:HB2	2.20	0.41
1:A:82:GLU:HG2	1:A:106:VAL:HG23	2.02	0.41
2:B:160:THR:HG21	2:B:210:LYS:HZ1	1.86	0.41
2:B:87:ARG:CD	2:B:89:GLU:HG2	2.50	0.41
2:H:97:ALA:HB1	2:H:109:PRO:HB3	2.03	0.41
1:L:149:TRP:HA	1:L:193:SER:O	2.21	0.41
1:A:114:PRO:HB3	1:A:140:PHE:HD2	1.84	0.41
3:L:225:HOH:O	2:H:173:HIS:HD2	2.03	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
1:L:38:LYS:HZ1	1:L:80:GLU:HG2	1.86	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:101:GLY:N	2:B:106:ASP:HB3	2.36	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 [i](#)

5.3.1 Protein backbone [i](#)

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%)	7	2
1	L	208/213 (98%)	196 (94%)	9 (4%)	3 (1%)	13	7
2	B	220/229 (96%)	209 (95%)	10 (4%)	1 (0%)	32	28
2	H	220/229 (96%)	211 (96%)	9 (4%)	0	100	100
All	All	856/884 (97%)	806 (94%)	41 (5%)	9 (1%)	17	11

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%)	160 (91%)	16 (9%)	11	7
1	L	176/179 (98%)	165 (94%)	11 (6%)	21	17
2	B	185/192 (96%)	164 (89%)	21 (11%)	7	3
2	H	185/192 (96%)	175 (95%)	10 (5%)	26	23
All	All	722/742 (97%)	664 (92%)	58 (8%)	14	10

All (58) 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	91	ASP
1	L	92	MET
1	L	127	GLN
1	L	136	LEU
1	L	145	VAL
1	L	157	LYS
1	L	190	ARG
1	L	205	LYS
2	H	1	GLN
2	H	19	ARG
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	173	HIS
2	H	202	THR
1	A	3	GLU
1	A	12	VAL
1	A	51	ASN

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Mol	Chain	Res	Type
1	A	60	ARG
1	A	72	LEU
1	A	89	THR
1	A	92	MET
1	A	93	VAL
1	A	115	SER
1	A	118	LEU
1	A	136	LEU
1	A	150	LYS
1	A	170	ASN
1	A	180	SER
1	A	190	ARG
1	A	195	GLN
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	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	180	GLN
2	B	187	LEU
2	B	200	THR
2	B	201	GLN
2	B	213	ASN
2	B	221	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	65	ASN
1	L	78	GLN
1	L	185	GLN
2	H	57	ASN

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Mol	Chain	Res	Type
2	H	82	GLN
1	A	16	GLN
1	A	37	GLN
1	A	41	GLN
1	A	51	ASN
1	A	78	GLN
1	A	127	GLN
1	A	170	ASN
1	A	185	GLN
2	B	1	GLN
2	B	30	ASN
2	B	82	GLN
2	B	173	HIS
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 carbohydrates 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 ⓘ

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%)	0.92	30 (14%) 3 4	37, 65, 95, 109	0
1	L	210/213 (98%)	0.61	16 (7%) 15 19	33, 52, 88, 108	0
2	B	222/229 (96%)	0.38	9 (4%) 38 45	32, 46, 83, 132	0
2	H	222/229 (96%)	0.16	7 (3%) 48 55	29, 41, 61, 125	0
All	All	864/884 (97%)	0.51	62 (7%) 16 21	29, 50, 88, 132	0

All (62) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	H	139	SER	8.4
2	B	139	SER	8.2
1	L	108	GLY	7.3
1	L	210	THR	6.3
1	A	210	THR	6.0
1	A	1	SER	5.8
2	B	142	GLY	5.4
1	A	109	GLN	5.0
1	L	107	LEU	5.0
1	A	201	SER	4.9
1	A	108	GLY	4.7
2	B	138	LYS	4.5
2	B	137	SER	4.4
1	L	126	LEU	4.4
1	A	142	PRO	4.2
1	A	106	VAL	4.1
1	A	2	TYR	4.1
1	L	169	SER	3.9
1	L	2	TYR	3.9
1	L	184	GLU	3.8
1	A	126	LEU	3.8

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Mol	Chain	Res	Type	RSRZ
2	B	198	LEU	3.8
2	H	138	LYS	3.5
2	H	137	SER	3.4
1	A	158	ALA	3.3
2	B	136	SER	3.3
1	A	144	ALA	3.3
1	A	194	CYS	3.2
2	B	143	GLY	3.1
2	H	1	GLN	3.0
2	B	141	SER	3.0
1	L	106	VAL	3.0
1	A	164	THR	3.0
1	A	192	TYR	3.0
1	A	107	LEU	2.9
1	A	19	ARG	2.8
1	L	16	GLN	2.8
1	L	130	LYS	2.8
1	A	77	THR	2.8
1	L	187	LYS	2.7
1	A	16	GLN	2.6
1	A	183	PRO	2.6
1	A	153	SER	2.6
1	A	188	SER	2.6
1	A	143	GLY	2.6
1	L	203	VAL	2.6
2	B	200	THR	2.5
1	A	157	LYS	2.5
1	A	199	GLU	2.5
1	A	15	GLY	2.5
1	A	200	GLY	2.5
1	L	183	PRO	2.3
1	A	184	GLU	2.3
1	A	127	GLN	2.3
2	H	37	VAL	2.3
1	A	11	SER	2.3
2	H	136	SER	2.3
1	A	12	VAL	2.2
1	L	13	ALA	2.2
2	H	68	PHE	2.1
1	L	144	ALA	2.1
1	L	194	CYS	2.0

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

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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.