



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 12:08 PM EDT

PDB ID : 3QBQ / pdb_00003qbq
Title : Crystal structure of extracellular domains of mouse RANK-RANKL complex
Authors : Ta, H.M.; Nguyen, G.T.T.; Jin, H.M.; Choi, J.K.; Park, H.; Kim, N.S.; Hwang, H.Y.; Kim, K.K.
Deposited on : 2011-01-13
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

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

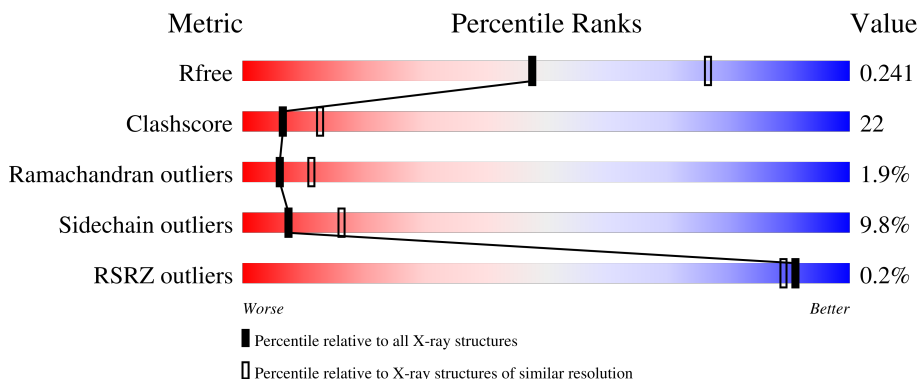
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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}	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (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 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	160	
1	C	160	
2	B	170	
2	D	170	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5114 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 Tumor necrosis factor ligand superfamily member 11.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	155	Total	C	N	O	S	0	0	0
			1224	783	209	228	4			
1	C	155	Total	C	N	O	S	0	0	0
			1224	783	209	228	4			

- Molecule 2 is a protein called Tumor necrosis factor receptor superfamily member 11A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	164	Total	C	N	O	S	0	0	0
			1242	757	220	245	20			
2	D	164	Total	C	N	O	S	0	0	0
			1242	757	220	245	20			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	45	Total	O	0	0
			45	45		
3	B	45	Total	O	0	0
			45	45		
3	C	47	Total	O	0	0
			47	47		
3	D	45	Total	O	0	0
			45	45		



4 Data and refinement statistics

Property	Value	Source
Space group	P 3	Depositor
Cell constants a, b, c, α , β , γ	120.86Å 120.86Å 93.42Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	45.66 – 2.50 45.66 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.6 (45.66-2.50) 99.6 (45.66-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.15 (at 2.51Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.227 , 0.249 0.222 , 0.241	Depositor DCC
R_{free} test set	2626 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	52.7	Xtriage
Anisotropy	0.598	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 42.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.487 for -h,-k,l 0.041 for h,-h-k,-l 0.041 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5114	wwPDB-VP
Average B, all atoms (Å ²)	62.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.08% 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	0.57	0/1259	1.12	9/1707 (0.5%)
1	C	0.64	0/1259	1.15	11/1707 (0.6%)
2	B	0.50	0/1272	1.06	5/1729 (0.3%)
2	D	0.49	0/1272	1.06	8/1729 (0.5%)
All	All	0.55	0/5062	1.10	33/6872 (0.5%)

There are no bond length outliers.

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	248	ILE	CA-C-N	9.08	131.19	119.84
1	C	248	ILE	C-N-CA	9.08	131.19	119.84
1	A	248	ILE	CA-C-N	9.07	131.18	119.84
1	A	248	ILE	C-N-CA	9.07	131.18	119.84
2	D	82	ASN	N-CA-C	8.40	121.11	109.18
2	B	82	ASN	N-CA-C	7.94	120.45	109.18
1	C	230	VAL	N-CA-C	-7.09	101.11	107.56
1	A	230	VAL	N-CA-C	-6.84	101.34	107.56
1	A	242	VAL	N-CA-C	6.57	118.05	108.53
1	C	188	HIS	N-CA-C	-6.55	103.94	114.09
2	D	168	ASP	N-CA-C	6.52	119.92	110.28
2	D	80	THR	N-CA-C	6.50	118.03	108.86
1	C	189	ASP	N-CA-C	6.35	122.95	114.12
1	A	260	THR	N-CA-C	-6.31	100.23	110.20
2	D	37	GLN	N-CA-C	6.29	118.22	111.36
1	C	254	LEU	N-CA-C	-6.05	105.93	113.55
1	C	242	VAL	N-CA-C	6.02	118.04	108.89
2	B	102	VAL	N-CA-C	-5.96	106.69	111.81
1	C	260	THR	N-CA-C	-5.86	100.95	110.20
2	D	62	THR	CA-C-N	-5.72	113.86	119.64
2	D	62	THR	C-N-CA	-5.72	113.86	119.64
1	A	293	SER	N-CA-C	5.72	117.31	111.14
2	B	78	LEU	N-CA-C	-5.69	98.24	108.48

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	293	SER	N-CA-C	5.38	117.23	111.36
1	A	269	PHE	N-CA-C	5.36	118.19	109.24
2	D	83	GLU	N-CA-C	-5.24	105.17	112.30
2	B	81	TRP	N-CA-C	-5.19	101.96	109.59
2	D	176	CYS	N-CA-C	5.18	117.60	111.33
1	C	274	ILE	CB-CA-C	-5.14	102.46	110.69
1	A	176	SER	N-CA-C	5.09	119.59	113.17
1	C	269	PHE	N-CA-C	5.04	117.67	109.24
2	B	168	ASP	N-CA-C	5.04	117.73	110.28
1	A	249	PRO	N-CA-C	5.03	122.84	112.47

There are no chirality outliers.

There are no planarity outliers.

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	1224	0	1178	39	0
1	C	1224	0	1178	46	0
2	B	1242	0	1138	64	0
2	D	1242	0	1138	64	0
3	A	45	0	0	4	0
3	B	45	0	0	1	0
3	C	47	0	0	3	0
3	D	45	0	0	1	0
All	All	5114	0	4632	207	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 22.

All (207) 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:D:106:ASN:ND2	2:D:109:ALA:H	1.63	0.94
1:A:245:SER:HB2	1:A:248:ILE:HB	1.54	0.87

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:242:VAL:CG1	1:A:251:SER:HB3	2.05	0.86
2:B:58:SER:HB2	2:B:70:LEU:HD12	1.59	0.85
2:B:106:ASN:ND2	2:B:109:ALA:H	1.76	0.83
1:C:225:GLU:OE1	2:D:130:ARG:NH2	2.13	0.82
2:D:41:TYR:HB3	2:D:50:ARG:HE	1.45	0.81
2:D:145:GLN:HE21	2:D:145:GLN:HA	1.47	0.79
2:D:131:ASN:HB3	2:D:150:THR:HG22	1.64	0.79
1:A:204:LYS:NZ	1:A:244:THR:HG21	1.99	0.78
1:A:225:GLU:CD	2:B:130:ARG:HH22	1.92	0.77
2:D:51:CYS:O	2:D:81:TRP:HA	1.86	0.76
2:D:186:GLN:HE21	2:D:187:GLY:H	1.35	0.74
1:A:247:LYS:C	1:A:248:ILE:HD12	2.12	0.73
2:B:106:ASN:HD22	2:B:106:ASN:C	1.97	0.70
2:D:145:GLN:HB2	2:D:148:LYS:HB2	1.73	0.70
2:B:37:GLN:C	2:B:39:ARG:H	2.00	0.69
1:C:312:VAL:HG22	1:C:313:GLN:HG3	1.74	0.69
2:B:51:CYS:O	2:B:81:TRP:HA	1.92	0.69
2:B:132:THR:HG22	2:B:149:ASP:OD2	1.93	0.69
2:D:131:ASN:HB3	2:D:150:THR:CG2	2.23	0.68
2:B:106:ASN:ND2	2:B:108:THR:H	1.91	0.68
2:B:185:HIS:HB2	2:B:194:VAL:HB	1.76	0.68
2:B:131:ASN:HB3	2:B:150:THR:HG23	1.75	0.67
1:A:270:HIS:HD2	3:A:97:HOH:O	1.77	0.67
2:D:122:ASN:ND2	2:D:125:CYS:H	1.93	0.67
1:A:170:ASN:HD22	1:A:170:ASN:C	2.02	0.67
1:A:245:SER:HB3	1:A:248:ILE:HD13	1.76	0.67
2:B:50:ARG:HG2	2:B:50:ARG:HH11	1.59	0.67
1:C:225:GLU:HG3	1:C:268:GLU:HA	1.75	0.67
2:D:35:CYS:HB2	2:D:42:GLU:OE2	1.93	0.67
1:A:210:ASP:HB3	1:A:315:ILE:HD12	1.76	0.66
1:C:192:TRP:O	1:C:194:LYS:HE2	1.96	0.66
2:D:106:ASN:ND2	2:D:108:THR:H	1.93	0.66
2:D:106:ASN:HD21	2:D:109:ALA:H	1.42	0.65
2:D:106:ASN:C	2:D:106:ASN:HD22	2.03	0.65
2:D:57:LEU:HD12	2:D:57:LEU:O	1.97	0.64
1:C:248:ILE:N	1:C:248:ILE:HD12	2.12	0.64
1:C:170:ASN:C	1:C:170:ASN:HD22	2.02	0.64
2:B:55:LYS:HB2	2:B:70:LEU:O	1.97	0.64
2:B:130:ARG:HH21	2:B:130:ARG:HG3	1.63	0.64
1:C:223:HIS:HD2	3:C:54:HOH:O	1.80	0.64
1:C:242:VAL:CG1	1:C:251:SER:HB2	2.28	0.64

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:145:GLN:HB2	2:B:148:LYS:HB2	1.79	0.64
2:D:106:ASN:HD22	2:D:109:ALA:H	1.43	0.63
1:C:268:GLU:HG2	2:D:130:ARG:HB2	1.79	0.63
2:D:41:TYR:O	2:D:42:GLU:HG3	1.99	0.63
2:D:122:ASN:HD22	2:D:125:CYS:H	1.46	0.63
2:B:57:LEU:HD12	2:B:57:LEU:O	1.99	0.63
2:D:106:ASN:ND2	2:D:109:ALA:N	2.44	0.63
2:D:55:LYS:HB2	2:D:70:LEU:O	1.99	0.63
1:C:245:SER:C	1:C:247:LYS:H	2.07	0.62
2:B:145:GLN:HE21	2:B:145:GLN:HA	1.65	0.62
2:D:37:GLN:C	2:D:39:ARG:H	2.06	0.62
1:A:206:ARG:HD3	1:A:246:ILE:HD12	1.81	0.62
2:B:36:THR:O	2:B:40:HIS:HB2	2.00	0.61
2:B:41:TYR:HB3	2:B:50:ARG:HE	1.64	0.61
1:A:242:VAL:HG13	1:A:251:SER:HB3	1.81	0.61
2:B:37:GLN:O	2:B:39:ARG:N	2.33	0.61
2:B:38:GLU:O	2:B:39:ARG:HB2	2.01	0.60
2:B:106:ASN:HD22	2:B:109:ALA:H	1.49	0.60
1:A:244:THR:HB	1:A:251:SER:OG	2.00	0.60
1:A:268:GLU:HG2	2:B:130:ARG:HB2	1.84	0.60
2:D:53:PRO:HG3	2:D:80:THR:HA	1.83	0.60
1:A:223:HIS:HD2	3:A:97:HOH:O	1.85	0.59
1:A:242:VAL:HG11	1:A:251:SER:HB3	1.81	0.59
1:C:245:SER:O	1:C:247:LYS:N	2.35	0.59
2:B:156:LEU:CD2	2:B:156:LEU:H	2.15	0.59
2:D:156:LEU:H	2:D:156:LEU:CD2	2.16	0.59
1:A:245:SER:CB	1:A:248:ILE:HD13	2.34	0.58
1:A:164:PHE:C	1:A:164:PHE:CD1	2.83	0.57
2:D:120:HIS:CD2	2:D:149:ASP:HB2	2.40	0.56
1:A:245:SER:C	1:A:247:LYS:H	2.12	0.56
2:D:57:LEU:HD23	2:D:83:GLU:HA	1.87	0.56
1:A:204:LYS:HZ2	1:A:244:THR:HG21	1.70	0.56
2:D:50:ARG:HH11	2:D:50:ARG:HG2	1.71	0.56
2:B:58:SER:CB	2:B:70:LEU:HD12	2.34	0.56
2:D:145:GLN:HE21	2:D:145:GLN:CA	2.14	0.56
1:A:225:GLU:OE2	2:B:130:ARG:NH2	2.38	0.55
2:D:150:THR:HG21	2:D:166:SER:HB2	1.89	0.55
1:A:268:GLU:CG	2:B:130:ARG:HB2	2.37	0.54
2:B:91:LYS:O	2:B:111:ARG:HD2	2.07	0.54
2:B:57:LEU:O	2:B:57:LEU:CD1	2.56	0.54
2:D:57:LEU:HD12	2:D:57:LEU:C	2.33	0.54

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:91:LYS:HE3	3:B:238:HOH:O	2.07	0.54
2:D:132:THR:CG2	2:D:149:ASP:OD2	2.57	0.54
1:A:232:THR:O	1:A:232:THR:HG22	2.07	0.53
1:C:164:PHE:C	1:C:164:PHE:CD1	2.87	0.53
2:D:79:ASP:OD2	2:D:110:PRO:HD3	2.09	0.53
2:B:50:ARG:HG2	2:B:50:ARG:NH1	2.22	0.53
2:D:122:ASN:HB3	2:D:125:CYS:SG	2.49	0.52
1:C:242:VAL:HG22	1:C:253:ASN:HD22	1.74	0.52
2:D:106:ASN:HD21	2:D:109:ALA:N	2.02	0.52
1:C:210:ASP:HB3	1:C:315:ILE:HD12	1.91	0.52
1:C:230:VAL:HG11	1:C:263:TRP:O	2.09	0.52
1:C:270:HIS:HD2	3:C:54:HOH:O	1.91	0.52
1:C:243:LYS:HB2	1:C:254:LEU:HD11	1.92	0.52
2:D:119:TYR:CD2	2:D:130:ARG:HA	2.44	0.51
2:D:134:CYS:HB2	3:D:211:HOH:O	2.10	0.51
2:D:99:LEU:HD13	2:D:128:CYS:SG	2.50	0.51
2:D:37:GLN:O	2:D:39:ARG:N	2.43	0.51
1:C:232:THR:CG2	1:C:232:THR:O	2.59	0.50
2:D:55:LYS:HG3	2:D:69:CYS:HB3	1.92	0.50
1:C:279:PHE:CE1	1:C:312:VAL:HG11	2.46	0.50
1:A:207:VAL:O	1:A:284:ALA:HA	2.12	0.50
2:B:58:SER:HB2	2:B:70:LEU:CD1	2.36	0.50
2:B:99:LEU:HD13	2:B:128:CYS:SG	2.52	0.50
2:B:106:ASN:ND2	2:B:106:ASN:C	2.68	0.50
2:B:106:ASN:HD22	2:B:108:THR:H	1.60	0.49
2:D:38:GLU:O	2:D:39:ARG:HB2	2.10	0.49
2:D:103:ASP:HB3	2:D:112:ARG:HB2	1.92	0.49
1:C:279:PHE:HE1	1:C:312:VAL:HG11	1.76	0.49
1:A:236:GLN:HG2	1:A:238:MET:HG2	1.93	0.49
2:B:66:ASP:CG	2:B:67:SER:H	2.20	0.49
1:C:248:ILE:N	1:C:248:ILE:CD1	2.76	0.49
1:A:170:ASN:C	1:A:170:ASN:ND2	2.71	0.49
1:A:245:SER:C	1:A:247:LYS:N	2.71	0.49
2:B:160:PHE:CZ	2:B:192:ASP:HB2	2.48	0.49
1:C:236:GLN:HE21	1:C:260:THR:HG23	1.78	0.49
2:D:55:LYS:HB2	2:D:70:LEU:C	2.38	0.48
2:B:51:CYS:SG	2:B:57:LEU:HB3	2.52	0.48
1:C:206:ARG:HD3	1:C:246:ILE:HD12	1.96	0.48
1:C:206:ARG:HH11	1:C:206:ARG:HG2	1.78	0.48
2:D:50:ARG:HG2	2:D:50:ARG:NH1	2.28	0.48
2:B:52:GLU:HB3	2:B:81:TRP:CZ3	2.48	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:245:SER:O	1:A:247:LYS:N	2.46	0.48
2:B:46:ARG:NH1	2:B:62:THR:O	2.46	0.48
2:B:55:LYS:HB3	2:B:71:PRO:HA	1.95	0.48
2:B:156:LEU:H	2:B:156:LEU:HD23	1.78	0.48
2:D:52:GLU:HB3	2:D:81:TRP:CH2	2.48	0.48
1:A:170:ASN:ND2	1:A:172:ALA:H	2.11	0.47
1:A:223:HIS:HE1	3:A:106:HOH:O	1.96	0.47
2:D:37:GLN:C	2:D:39:ARG:N	2.72	0.47
2:D:186:GLN:HG3	2:D:187:GLY:N	2.30	0.47
2:B:150:THR:HG21	2:B:166:SER:O	2.14	0.47
1:A:170:ASN:HD22	1:A:172:ALA:H	1.60	0.47
1:C:170:ASN:C	1:C:170:ASN:ND2	2.72	0.47
2:B:52:GLU:HB3	2:B:81:TRP:CH2	2.49	0.47
2:D:160:PHE:HA	2:D:189:THR:O	2.15	0.47
2:D:91:LYS:O	2:D:111:ARG:HD2	2.15	0.46
2:B:106:ASN:ND2	2:B:109:ALA:N	2.55	0.46
2:B:132:THR:CG2	2:B:149:ASP:OD2	2.61	0.46
2:B:132:THR:HG22	2:B:149:ASP:CG	2.41	0.46
1:C:223:HIS:HE1	3:C:127:HOH:O	1.99	0.46
2:B:186:GLN:HE21	2:B:187:GLY:H	1.62	0.46
1:A:242:VAL:HG22	1:A:253:ASN:ND2	2.31	0.46
2:B:122:ASN:HB3	2:B:125:CYS:SG	2.55	0.46
2:B:103:ASP:HB3	2:B:112:ARG:HB2	1.98	0.45
2:D:53:PRO:HG3	2:D:80:THR:CA	2.44	0.45
2:D:36:THR:O	2:D:40:HIS:HB2	2.16	0.45
1:C:207:VAL:O	1:C:284:ALA:HA	2.16	0.45
2:D:132:THR:HG23	2:D:149:ASP:OD2	2.16	0.45
1:C:306:TYR:CD1	1:C:306:TYR:C	2.95	0.45
2:D:52:GLU:HB3	2:D:81:TRP:CZ3	2.51	0.45
2:D:106:ASN:HD22	2:D:108:THR:H	1.63	0.45
1:A:218:ASN:O	1:A:219:ILE:HD13	2.17	0.45
2:B:186:GLN:HG3	2:B:187:GLY:N	2.31	0.44
2:B:91:LYS:HE2	2:B:111:ARG:O	2.17	0.44
1:C:244:THR:HG23	1:C:287:GLU:HB2	2.00	0.44
1:C:246:ILE:HG12	1:C:246:ILE:O	2.17	0.44
2:B:122:ASN:ND2	2:B:125:CYS:H	2.16	0.44
1:C:244:THR:HB	1:C:251:SER:HB3	1.99	0.44
1:C:245:SER:HA	1:C:286:GLU:HA	1.99	0.44
2:B:182:LEU:HD23	2:B:198:SER:CB	2.48	0.44
2:B:37:GLN:C	2:B:39:ARG:N	2.67	0.43
2:B:85:ASP:O	2:B:86:LYS:HD2	2.17	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:78:LEU:HD23	2:D:78:LEU:HA	1.80	0.43
1:A:221:PHE:HE1	1:A:274:ILE:HD11	1.83	0.43
1:C:263:TRP:CE2	1:C:270:HIS:HB3	2.53	0.43
2:B:52:GLU:CG	2:B:55:LYS:HD3	2.48	0.43
1:C:186:TRP:CD1	1:C:203:GLY:HA2	2.54	0.43
1:C:215:LEU:O	1:C:277:GLY:HA2	2.18	0.43
2:D:43:HIS:CD2	2:D:44:LEU:HG	2.54	0.43
1:A:264:SER:O	1:A:265:GLY:O	2.37	0.42
1:C:262:ASN:HD22	1:C:262:ASN:C	2.27	0.42
2:D:39:ARG:HG3	2:D:81:TRP:CZ3	2.54	0.42
1:C:204:LYS:NZ	1:C:244:THR:HG21	2.34	0.42
2:B:145:GLN:HE21	2:B:145:GLN:CA	2.27	0.42
2:D:159:PHE:CE2	2:D:172:PRO:HB3	2.55	0.42
2:D:184:ALA:HB2	2:D:196:SER:HB2	2.00	0.42
2:B:50:ARG:CD	2:B:83:GLU:HG2	2.50	0.42
1:C:183:LEU:HD11	1:C:298:LEU:HD11	2.00	0.42
1:A:183:LEU:HD23	1:A:183:LEU:HA	1.93	0.42
2:B:42:GLU:OE2	2:B:47:CYS:HA	2.19	0.42
2:D:106:ASN:ND2	2:D:106:ASN:C	2.74	0.42
1:A:247:LYS:O	1:A:248:ILE:HD12	2.20	0.42
2:B:106:ASN:HD21	2:B:109:ALA:H	1.62	0.42
2:D:55:LYS:HB3	2:D:71:PRO:HA	2.01	0.41
1:C:162:GLN:HA	1:C:163:PRO:HD3	1.95	0.41
1:C:245:SER:C	1:C:247:LYS:N	2.74	0.41
1:C:245:SER:OG	1:C:248:ILE:HD13	2.19	0.41
2:B:130:ARG:HG3	2:B:130:ARG:NH2	2.33	0.41
1:C:230:VAL:HG13	1:C:264:SER:HA	2.01	0.41
1:C:242:VAL:HG22	1:C:253:ASN:ND2	2.34	0.41
2:D:185:HIS:HB2	2:D:194:VAL:CG2	2.50	0.41
1:A:211:GLY:HA2	1:A:315:ILE:HG13	2.02	0.41
2:B:44:LEU:C	2:B:46:ARG:H	2.29	0.41
1:C:225:GLU:CG	1:C:268:GLU:HA	2.46	0.41
2:B:160:PHE:HA	2:B:189:THR:O	2.21	0.41
2:D:156:LEU:H	2:D:156:LEU:HD23	1.84	0.41
2:B:43:HIS:CD2	2:B:44:LEU:HG	2.56	0.41
1:C:186:TRP:HD1	1:C:203:GLY:HA2	1.86	0.41
1:C:212:PHE:HB3	1:C:279:PHE:CZ	2.56	0.41
1:A:294:ASN:HB3	3:A:107:HOH:O	2.21	0.40
2:D:185:HIS:HB2	2:D:194:VAL:HB	2.02	0.40
1:A:215:LEU:O	1:A:277:GLY:HA2	2.21	0.40
2:D:122:ASN:HD21	2:D:124:ASP:HB2	1.86	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	153/160 (96%)	140 (92%)	9 (6%)	4 (3%)	4	7
1	C	153/160 (96%)	146 (95%)	5 (3%)	2 (1%)	9	18
2	B	162/170 (95%)	152 (94%)	7 (4%)	3 (2%)	6	11
2	D	162/170 (95%)	152 (94%)	7 (4%)	3 (2%)	6	11
All	All	630/660 (96%)	590 (94%)	28 (4%)	12 (2%)	6	11

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	249	PRO
2	B	36	THR
2	B	38	GLU
2	B	39	ARG
1	C	246	ILE
1	C	249	PRO
2	D	39	ARG
1	A	246	ILE
1	A	265	GLY
2	D	38	GLU
2	D	44	LEU
1	A	210	ASP

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	134/138 (97%)	125 (93%)	9 (7%)	15	31
1	C	134/138 (97%)	121 (90%)	13 (10%)	8	17
2	B	142/150 (95%)	127 (89%)	15 (11%)	6	14
2	D	142/150 (95%)	125 (88%)	17 (12%)	5	10
All	All	552/576 (96%)	498 (90%)	54 (10%)	7	16

All (54) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	170	ASN
1	A	232	THR
1	A	235	LEU
1	A	244	THR
1	A	249	PRO
1	A	266	ASN
1	A	274	ILE
1	A	281	LYS
1	A	312	VAL
2	B	36	THR
2	B	57	LEU
2	B	86	LYS
2	B	88	LEU
2	B	91	LYS
2	B	99	LEU
2	B	106	ASN
2	B	132	THR
2	B	144	LEU
2	B	145	GLN
2	B	151	VAL
2	B	156	LEU
2	B	179	LEU
2	B	191	SER
2	B	193	VAL
1	C	170	ASN
1	C	194	LYS
1	C	199	THR
1	C	230	VAL
1	C	232	THR
1	C	235	LEU
1	C	244	THR

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	C	248	ILE
1	C	249	PRO
1	C	262	ASN
1	C	267	SER
1	C	297	LEU
1	C	312	VAL
2	D	35	CYS
2	D	36	THR
2	D	47	CYS
2	D	57	LEU
2	D	79	ASP
2	D	88	LEU
2	D	99	LEU
2	D	106	ASN
2	D	132	THR
2	D	144	LEU
2	D	145	GLN
2	D	150	THR
2	D	151	VAL
2	D	156	LEU
2	D	157	LEU
2	D	179	LEU
2	D	193	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	170	ASN
1	A	218	ASN
1	A	236	GLN
1	A	253	ASN
1	A	270	HIS
1	A	275	ASN
2	B	43	HIS
2	B	106	ASN
2	B	122	ASN
2	B	145	GLN
2	B	186	GLN
1	C	170	ASN
1	C	202	ASN
1	C	218	ASN
1	C	236	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	C	253	ASN
1	C	262	ASN
1	C	270	HIS
1	C	275	ASN
1	C	313	GLN
2	D	43	HIS
2	D	106	ASN
2	D	122	ASN
2	D	131	ASN
2	D	145	GLN
2	D	186	GLN

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	155/160 (96%)	-1.06	0 100 100	34, 52, 84, 101	0
1	C	155/160 (96%)	-1.02	0 100 100	36, 52, 85, 100	0
2	B	164/170 (96%)	-0.98	1 (0%) 85 83	43, 64, 107, 126	0
2	D	164/170 (96%)	-0.94	0 100 100	44, 64, 108, 128	0
All	All	638/660 (96%)	-1.00	1 (0%) 91 89	34, 59, 95, 128	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	36	THR	2.1

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 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.