



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 11:55 AM GMT

PDB ID : 3QBQ
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
<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
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

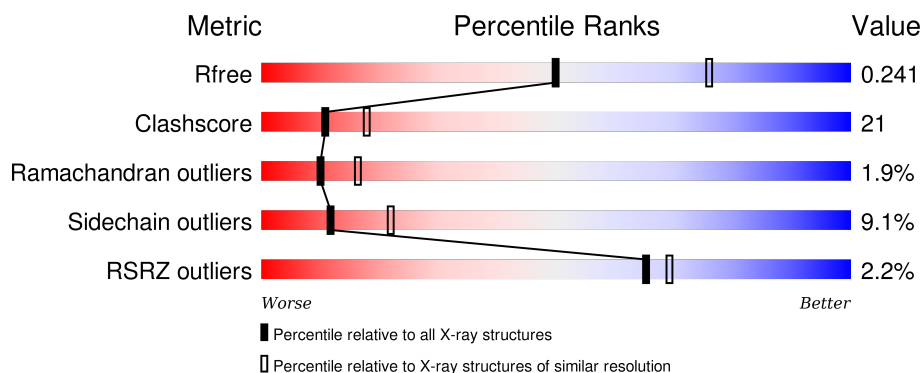
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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}	91344	3553 (2.50-2.50)
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (2.50-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 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	160	<div> <div>3%</div> <div> <div></div> <div>71%</div> <div>21%</div> <div>• •</div> </div> </div>
1	C	160	<div> <div>2%</div> <div> <div></div> <div>65%</div> <div>27%</div> <div>• • •</div> </div> </div>
2	B	170	<div> <div>%</div> <div> <div></div> <div>61%</div> <div>31%</div> <div>5% • •</div> </div> </div>
2	D	170	<div> <div>2%</div> <div> <div></div> <div>58%</div> <div>31%</div> <div>6% • •</div> </div> </div>

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			

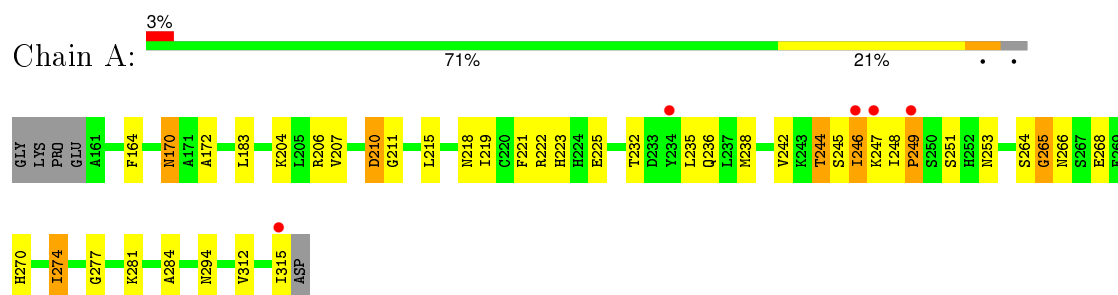
- Molecule 3 is water.

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		

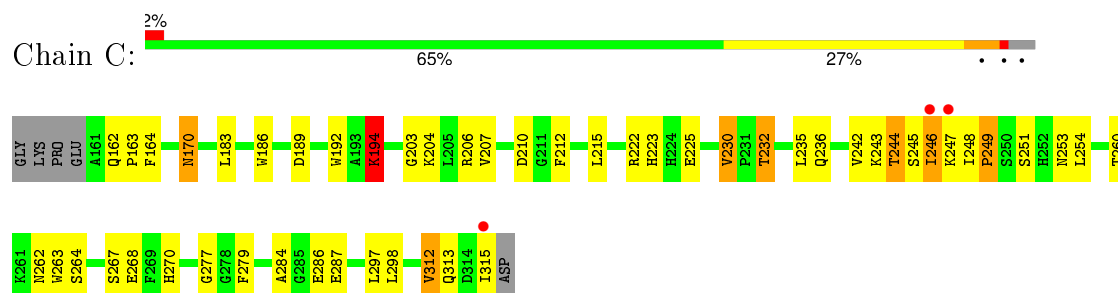
3 Residue-property plots

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.

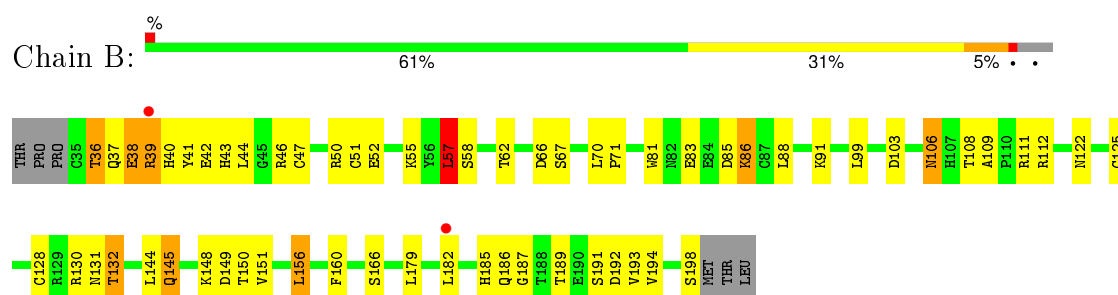
- Molecule 1: Tumor necrosis factor ligand superfamily member 11



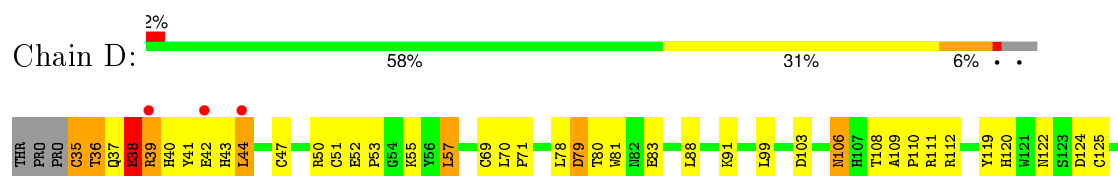
- Molecule 1: Tumor necrosis factor ligand superfamily member 11



- Molecule 2: Tumor necrosis factor receptor superfamily member 11A



- Molecule 2: Tumor necrosis factor receptor superfamily member 11A





4 Data and refinement statistics

Property	Value	Source
Space group	P 3	Depositor
Cell constants a, b, c, α , β , γ	120.86 Å 120.86 Å 93.42 Å 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 2.51 Å)	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%)	DCC
Wilson B-factor (Å ²)	52.7	Xtriage
Anisotropy	0.598	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 44.3	EDS
Estimated twinning fraction	0.487 for -h,-k,l 0.041 for h,-h-k,-l 0.041 for -k,-h,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 52575 reflections	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.*

¹ Intensities estimated from amplitudes.

² 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 [i](#)

5.1 Standard geometry [i](#)

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	0.50	0/1259	0.73	0/1707
1	C	0.59	2/1259 (0.2%)	0.80	1/1707 (0.1%)
2	B	0.44	0/1272	0.76	1/1729 (0.1%)
2	D	0.42	0/1272	0.75	0/1729
All	All	0.49	2/5062 (0.0%)	0.76	2/6872 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	194	LYS	CE-NZ	6.00	1.64	1.49
1	C	194	LYS	CD-CE	5.34	1.64	1.51

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	194	LYS	CD-CE-NZ	5.77	124.97	111.70
2	B	57	LEU	CA-CB-CG	5.03	126.86	115.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	43	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	1242	0	1138	63	0
2	D	1242	0	1138	63	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	202	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 21.

All (202) 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
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:HA	2:D:145:GLN:HE21	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
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:225:GLU:CD	2:B:130:ARG:HH22	1.92	0.72
2:D:145:GLN:HB2	2:D:148:LYS:HB2	1.73	0.70
1:A:247:LYS:C	1:A:248:ILE:HD12	2.12	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:106:ASN:HD22	2:B:106:ASN:C	1.97	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:245:SER:HB3	1:A:248:ILE:HD13	1.76	0.67
1:C:225:GLU:HG3	1:C:268:GLU:HA	1.75	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:50:ARG:HH11	2:B:50:ARG:HG2	1.59	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
2:D:106:ASN:ND2	2:D:108:THR:H	1.93	0.66
1:C:192:TRP:O	1:C:194:LYS:HE2	1.96	0.66
2:D:106:ASN:HD21	2:D:109:ALA:H	1.42	0.65
2:B:37:GLN:C	2:B:39:ARG:H	2.00	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
2:B:55:LYS:HB2	2:B:70:LEU:O	1.97	0.64
2:B:130:ARG:HG3	2:B:130:ARG:HH21	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
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:D:106:ASN:ND2	2:D:109:ALA:N	2.44	0.63
2:B:57:LEU:HD12	2:B:57:LEU:O	1.99	0.63
2:D:55:LYS:HB2	2:D:70:LEU:O	1.99	0.63
1:A:170:ASN:C	1:A:170:ASN:HD22	2.02	0.62
2:B:145:GLN:HE21	2:B:145:GLN:HA	1.65	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
1:C:170:ASN:C	1:C:170:ASN:HD22	2.02	0.61
2:D:106:ASN:C	2:D:106:ASN:HD22	2.03	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:242:VAL:HG11	1:A:251:SER:HB3	1.81	0.59
1:A:223:HIS:HD2	3:A:97:HOH:O	1.85	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
2:D:37:GLN:C	2:D:39:ARG:H	2.06	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:245:SER:CB	1:A:248:ILE:HD13	2.34	0.58
1:C:245:SER:C	1:C:247:LYS:H	2.07	0.58
2:D:120:HIS:CD2	2:D:149:ASP:HB2	2.40	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:HG2	2:D:50:ARG:HH11	1.71	0.56
2:B:58:SER:CB	2:B:70:LEU:HD12	2.34	0.56
2:D:145:GLN:CA	2:D:145:GLN:HE21	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:CD1	2:B:57:LEU:O	2.56	0.54
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
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
1:A:245:SER:C	1:A:247:LYS:H	2.12	0.53
2:D:122:ASN:HB3	2:D:125:CYS:SG	2.49	0.52
2:D:106:ASN:HD21	2:D:109:ALA:N	2.02	0.52
1:C:242:VAL:HG22	1:C:253:ASN:HD22	1.74	0.52
1:C:230:VAL:HG11	1:C:263:TRP:O	2.09	0.52
1:C:210:ASP:HB3	1:C:315:ILE:HD12	1.91	0.52
1:C:270:HIS:HD2	3:C:54:HOH:O	1.91	0.52
1:A:164:PHE:CD1	1:A:164:PHE:C	2.83	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
2:D:55:LYS:HG3	2:D:69:CYS:HB3	1.92	0.50
1:C:232:THR:O	1:C:232:THR:CG2	2.59	0.50
1:C:279:PHE:CE1	1:C:312:VAL:HG11	2.46	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
1:A:207:VAL:O	1:A:284:ALA:HA	2.12	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:236:GLN:HG2	1:A:238:MET:HG2	1.93	0.49
2:D:57:LEU:C	2:D:57:LEU:HD12	2.33	0.49
1:C:248:ILE:CD1	1:C:248:ILE:N	2.76	0.49
1:C:236:GLN:HE21	1:C:260:THR:HG23	1.78	0.49
2:B:160:PHE:CZ	2:B:192:ASP:HB2	2.48	0.49
2:B:51:CYS:SG	2:B:57:LEU:HB3	2.52	0.48
2:D:50:ARG:HG2	2:D:50:ARG:NH1	2.28	0.48
1:C:206:ARG:HH11	1:C:206:ARG:HG2	1.78	0.48
1:C:206:ARG:HD3	1:C:246:ILE:HD12	1.96	0.48
1:C:164:PHE:C	1:C:164:PHE:CD1	2.87	0.48
2:B:52:GLU:HB3	2:B:81:TRP:CZ3	2.48	0.48
1:A:245:SER:O	1:A:247:LYS:N	2.46	0.48
2:B:55:LYS:HB3	2:B:71:PRO:HA	1.95	0.48
2:B:46:ARG:NH1	2:B:62:THR:O	2.46	0.48
2:D:52:GLU:HB3	2:D:81:TRP:CH2	2.48	0.48
2:B:156:LEU:H	2:B:156:LEU:HD23	1.78	0.48
2:D:186:GLN:HG3	2:D:187:GLY:N	2.30	0.47
1:A:223:HIS:HE1	3:A:106:HOH:O	1.96	0.47
1:A:170:ASN:ND2	1:A:172:ALA:H	2.11	0.47
2:B:150:THR:HG21	2:B:166:SER:O	2.14	0.47
2:B:37:GLN:C	2:B:39:ARG:N	2.67	0.47
1:A:170:ASN:HD22	1:A:172:ALA:H	1.60	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
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:D:53:PRO:HG3	2:D:80:THR:CA	2.44	0.45
2:B:103:ASP:HB3	2:B:112:ARG:HB2	1.98	0.45
2:D:36:THR:O	2:D:40:HIS:HB2	2.16	0.45
2:D:132:THR:HG23	2:D:149:ASP:OD2	2.16	0.45
1:C:207:VAL:O	1:C:284:ALA:HA	2.16	0.45
2:D:106:ASN:HD22	2:D:108:THR:H	1.63	0.45
2:D:52:GLU:HB3	2:D:81:TRP:CZ3	2.51	0.45
1:A:218:ASN:O	1:A:219:ILE:HD13	2.17	0.45
2:B:66:ASP:CG	2:B:67:SER:H	2.20	0.45
2:B:106:ASN:C	2:B:106:ASN:ND2	2.68	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:186:GLN:HG3	2:B:187:GLY:N	2.31	0.44
1:A:245:SER:C	1:A:247:LYS:N	2.71	0.44
1:C:244:THR:HG23	1:C:287:GLU:HB2	2.00	0.44
2:B:91:LYS:HE2	2:B:111:ARG:O	2.17	0.44
1:C:246:ILE:HG12	1:C:246:ILE:O	2.17	0.44
2:D:55:LYS:HB2	2:D:70:LEU:C	2.38	0.44
2:B:122:ASN:ND2	2:B:125:CYS:H	2.16	0.44
1:C:245:SER:HA	1:C:286:GLU:HA	1.99	0.44
1:C:244:THR:HB	1:C:251:SER:HB3	1.99	0.44
2:B:182:LEU:HD23	2:B:198:SER:CB	2.48	0.44
2:B:85:ASP:O	2:B:86:LYS:HD2	2.17	0.43
2:D:78:LEU:HA	2:D:78:LEU:HD23	1.80	0.43
1:C:263:TRP:CE2	1:C:270:HIS:HB3	2.53	0.43
1:A:221:PHE:HE1	1:A:274:ILE:HD11	1.83	0.43
2:B:52:GLU:CG	2:B:55:LYS:HD3	2.48	0.43
1:C:215:LEU:O	1:C:277:GLY:HA2	2.18	0.43
1:C:186:TRP:CD1	1:C:203:GLY:HA2	2.54	0.43
2:D:37:GLN:C	2:D:39:ARG:N	2.72	0.43
2:D:43:HIS:CD2	2:D:44:LEU:HG	2.54	0.43
2:D:39:ARG:HG3	2:D:81:TRP:CZ3	2.54	0.42
1:A:264:SER:O	1:A:265:GLY:O	2.37	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:184:ALA:HB2	2:D:196:SER:HB2	2.00	0.42
2:D:159:PHE:CE2	2:D:172:PRO:HB3	2.55	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
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
1:C:170:ASN:C	1:C:170:ASN:ND2	2.72	0.41
2:D:55:LYS:HB3	2:D:71:PRO:HA	2.01	0.41
2:B:132:THR:HG22	2:B:149:ASP:CG	2.41	0.41
1:C:162:GLN:HA	1:C:163:PRO:HD3	1.95	0.41
1:C:245:SER:OG	1:C:248:ILE:HD13	2.19	0.41
1:A:170:ASN:C	1:A:170:ASN:ND2	2.71	0.41
2:B:130:ARG:NH2	2:B:130:ARG:HG3	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:225:GLU:CG	1:C:268:GLU:HA	2.46	0.41
1:A:211:GLY:HA2	1:A:315:ILE:HG13	2.02	0.41
2:B:160:PHE:HA	2:B:189:THR:O	2.21	0.41
2:D:156:LEU:HD23	2:D:156:LEU:H	1.84	0.41
1:C:212:PHE:HB3	1:C:279:PHE:CZ	2.56	0.41
1:C:186:TRP:HD1	1:C:203:GLY:HA2	1.86	0.41
2:B:43:HIS:CD2	2:B:44:LEU:HG	2.56	0.41
2:D:185:HIS:HB2	2:D:194:VAL:HB	2.02	0.40
1:A:294:ASN:HB3	3:A:107:HOH:O	2.21	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%)	7	10
1	C	153/160 (96%)	146 (95%)	5 (3%)	2 (1%)	15	26
2	B	162/170 (95%)	152 (94%)	7 (4%)	3 (2%)	10	16
2	D	162/170 (95%)	152 (94%)	7 (4%)	3 (2%)	10	16
All	All	630/660 (96%)	590 (94%)	28 (4%)	12 (2%)	10	16

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

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Mol	Chain	Res	Type
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 ⓘ

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%)	20	37
1	C	134/138 (97%)	121 (90%)	13 (10%)	10	19
2	B	142/150 (95%)	130 (92%)	12 (8%)	13	25
2	D	142/150 (95%)	126 (89%)	16 (11%)	7	13
All	All	552/576 (96%)	502 (91%)	50 (9%)	12	22

All (50) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	170	ASN
1	A	222	ARG
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	57	LEU
2	B	86	LYS
2	B	88	LEU
2	B	106	ASN

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Mol	Chain	Res	Type
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	189	ASP
1	C	194	LYS
1	C	222	ARG
1	C	230	VAL
1	C	232	THR
1	C	235	LEU
1	C	244	THR
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	38	GLU
2	D	47	CYS
2	D	57	LEU
2	D	79	ASP
2	D	88	LEU
2	D	106	ASN
2	D	132	THR
2	D	144	LEU
2	D	145	GLN
2	D	150	THR
2	D	156	LEU
2	D	157	LEU
2	D	175	ASN
2	D	179	LEU

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	170	ASN

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Mol	Chain	Res	Type
1	A	253	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	236	GLN
1	C	253	ASN
1	C	262	ASN
1	C	313	GLN
2	D	43	HIS
2	D	106	ASN
2	D	122	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 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 [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%)	0.50	5 (3%) 51 56	34, 52, 84, 101	0
1	C	155/160 (96%)	0.53	3 (1%) 70 73	36, 52, 85, 100	0
2	B	164/170 (96%)	0.33	2 (1%) 81 83	43, 64, 107, 126	0
2	D	164/170 (96%)	0.34	4 (2%) 62 66	44, 64, 108, 128	0
All	All	638/660 (96%)	0.42	14 (2%) 65 69	34, 59, 95, 128	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	39	ARG	5.6
1	C	246	ILE	5.2
1	A	246	ILE	4.1
2	D	44	LEU	3.9
2	B	182	LEU	3.1
1	A	247	LYS	2.9
2	D	185	HIS	2.9
1	A	234	TYR	2.7
1	C	315	ILE	2.6
1	C	247	LYS	2.4
1	A	249	PRO	2.4
2	D	42	GLU	2.4
2	D	39	ARG	2.1
1	A	315	ILE	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.