Full wwPDB X-ray Structure Validation Report

Jan 31, 2016 – 10:26 PM GMT

PDB ID	:	1TNF
Title	:	THE STRUCTURE OF TUMOR NECROSIS FACTOR-ALPHA AT 2.6 ANGSTROMS RESOLUTION. IMPLICATIONS FOR RECEPTOR BINDING
Authors	:	Eck, M.J.; Sprang, S.R.
Deposited on	:	1989-08-25
Resolution	:	2.60 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:

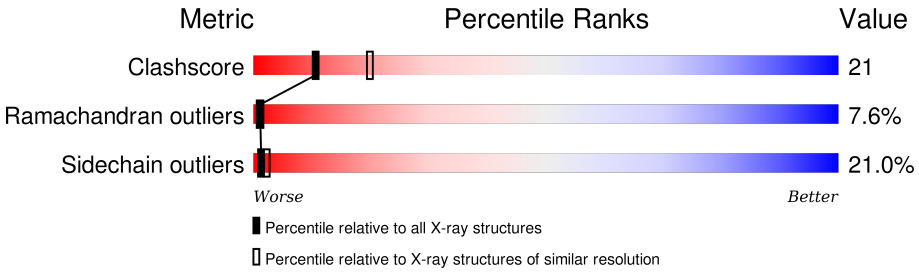
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	2679 (2.60-2.60)
Ramachandran outliers	100387	2635 (2.60-2.60)
Sidechain outliers	100360	2635 (2.60-2.60)

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 <=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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	157	42% 37% 13%
1	B	157	36% 44% 13%
1	C	157	31% 44% 15% 6%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 3552 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-ALPHA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	152	Total	C	N	O	S	0	0	0
			1184	757	204	221	2			
1	B	152	Total	C	N	O	S	0	0	0
			1184	757	204	221	2			
1	C	152	Total	C	N	O	S	0	0	0
			1184	757	204	221	2			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	45	ASP	LEU	CONFLICT	UNP P01375
B	45	ASP	LEU	CONFLICT	UNP P01375
C	45	ASP	LEU	CONFLICT	UNP P01375

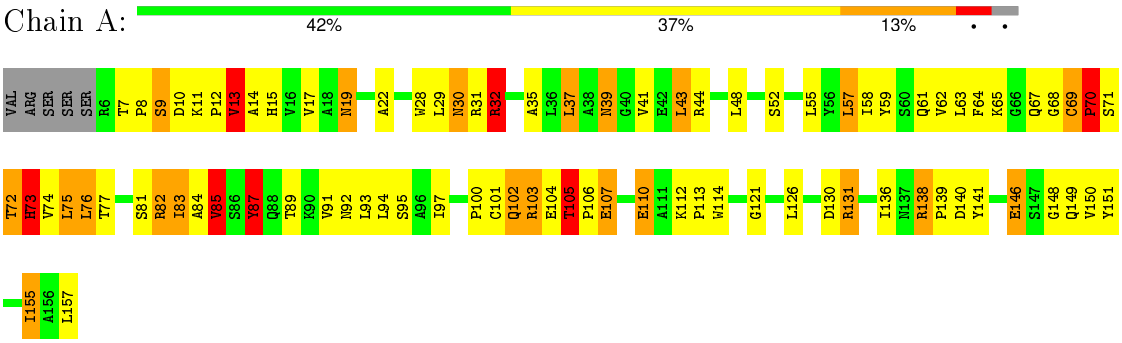
3 Residue-property plots (i)

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

Note EDS was not executed.

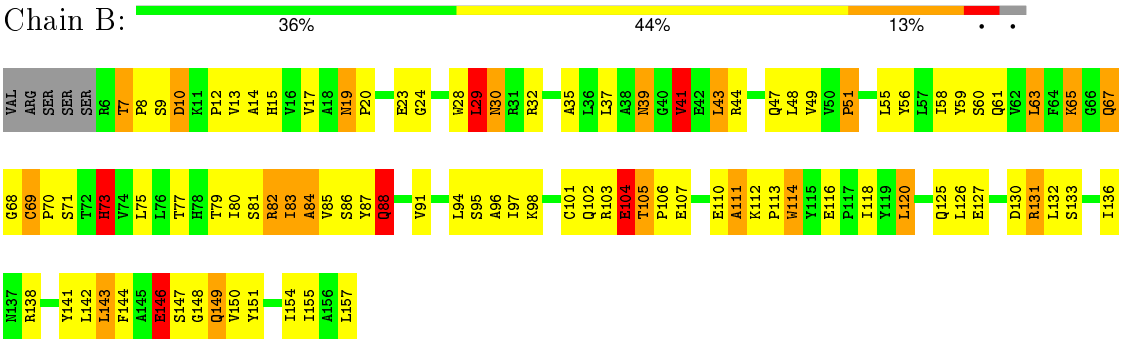
- Molecule 1: TUMOR NECROSIS FACTOR-ALPHA

Chain A:



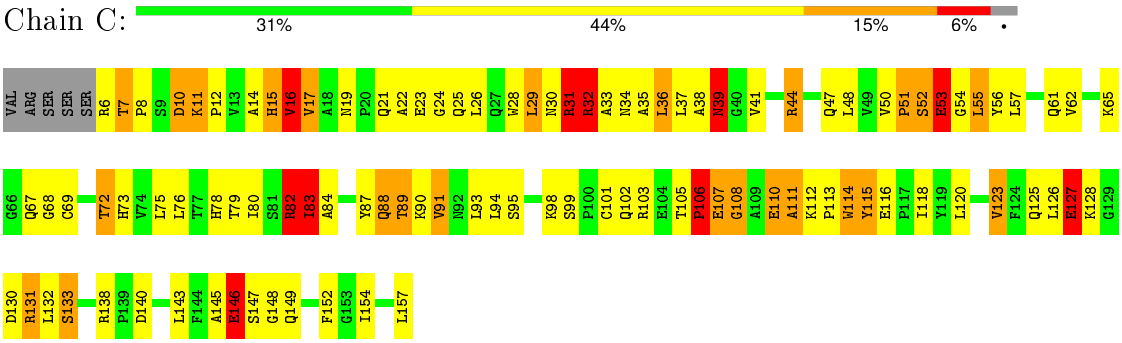
- Molecule 1: TUMOR NECROSIS FACTOR-ALPHA

Chain B:



- Molecule 1: TUMOR NECROSIS FACTOR-ALPHA

Chain C:



4 Data and refinement statistics (i)

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α, β, γ	95.00Å 95.00Å 117.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	(Not available) – 2.60	Depositor
% Data completeness (in resolution range)	(Not available) ((Not available)-2.60)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.230 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3552	wwPDB-VP
Average B, all atoms (Å^2)	25.0	wwPDB-VP

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	0.94	0/1211	2.03	43/1649 (2.6%)
1	B	0.96	1/1211 (0.1%)	1.95	30/1649 (1.8%)
1	C	0.97	0/1211	1.95	43/1649 (2.6%)
All	All	0.96	1/3633 (0.0%)	1.98	116/4947 (2.3%)

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	A	0	2
1	B	0	2
1	C	0	3
All	All	0	7

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	41	VAL	CA-CB	5.03	1.65	1.54

All (116) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	138	ARG	NE-CZ-NH2	16.25	128.43	120.30
1	B	138	ARG	NE-CZ-NH2	11.87	126.24	120.30
1	C	32	ARG	NE-CZ-NH2	11.09	125.84	120.30
1	B	131	ARG	NE-CZ-NH2	10.50	125.55	120.30
1	B	88	GLN	CA-CB-CG	10.07	135.56	113.40
1	B	138	ARG	NE-CZ-NH1	-9.55	115.52	120.30
1	A	28	TRP	CD1-CG-CD2	9.51	113.91	106.30
1	A	138	ARG	NE-CZ-NH1	-9.24	115.68	120.30
1	A	69	CYS	CA-CB-SG	-8.53	98.65	114.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	28	TRP	CE2-CD2-CG	-8.50	100.50	107.30
1	A	44	ARG	NE-CZ-NH2	8.32	124.46	120.30
1	B	114	TRP	CD1-CG-CD2	8.19	112.85	106.30
1	A	82	ARG	NE-CZ-NH1	-8.12	116.24	120.30
1	A	73	HIS	CA-CB-CG	-8.02	99.97	113.60
1	A	82	ARG	NE-CZ-NH2	7.81	124.21	120.30
1	A	114	TRP	CD1-CG-CD2	7.81	112.55	106.30
1	C	146	GLU	N-CA-C	7.77	131.97	111.00
1	B	28	TRP	CD1-CG-CD2	7.77	112.51	106.30
1	C	87	TYR	CB-CG-CD2	-7.58	116.45	121.00
1	C	106	PRO	N-CA-C	7.55	131.73	112.10
1	A	85	VAL	CA-CB-CG2	-7.52	99.63	110.90
1	C	107	GLU	CA-CB-CG	7.47	129.84	113.40
1	C	114	TRP	CD1-CG-CD2	7.45	112.26	106.30
1	B	28	TRP	CE2-CD2-CG	-7.42	101.37	107.30
1	C	114	TRP	CE2-CD2-CG	-7.35	101.42	107.30
1	C	44	ARG	CA-C-N	7.28	133.22	117.20
1	B	114	TRP	CE2-CD2-CG	-7.27	101.48	107.30
1	C	115	TYR	CB-CG-CD1	-7.11	116.74	121.00
1	A	114	TRP	CE2-CD2-CG	-7.10	101.62	107.30
1	C	28	TRP	CG-CD2-CE3	7.04	140.24	133.90
1	A	102	GLN	N-CA-C	-7.03	92.01	111.00
1	A	44	ARG	NE-CZ-NH1	-6.99	116.81	120.30
1	A	103	ARG	NE-CZ-NH2	6.98	123.79	120.30
1	C	87	TYR	N-CA-C	-6.97	92.19	111.00
1	A	37	LEU	N-CA-C	-6.84	92.54	111.00
1	A	87	TYR	CA-CB-CG	6.82	126.36	113.40
1	C	38	ALA	CA-C-N	-6.79	102.26	117.20
1	C	28	TRP	CE2-CD2-CG	-6.78	101.88	107.30
1	C	113	PRO	N-CA-C	6.76	129.67	112.10
1	C	88	GLN	CA-CB-CG	-6.72	98.62	113.40
1	A	32	ARG	NE-CZ-NH2	6.69	123.64	120.30
1	A	32	ARG	NE-CZ-NH1	-6.63	116.98	120.30
1	B	29	LEU	CA-CB-CG	6.59	130.45	115.30
1	C	10	ASP	CA-C-N	-6.58	102.73	117.20
1	C	91	VAL	CA-CB-CG2	-6.54	101.10	110.90
1	B	151	TYR	CB-CG-CD1	-6.52	117.09	121.00
1	B	82	ARG	CA-CB-CG	6.44	127.56	113.40
1	C	44	ARG	NE-CZ-NH2	6.41	123.50	120.30
1	C	38	ALA	O-C-N	6.40	132.94	122.70
1	B	71	SER	CA-C-N	-6.33	103.27	117.20
1	B	111	ALA	N-CA-C	6.31	128.03	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	43	LEU	CA-CB-CG	6.29	129.77	115.30
1	A	101	CYS	CA-CB-SG	-6.28	102.70	114.00
1	C	28	TRP	CD1-CG-CD2	6.26	111.31	106.30
1	C	87	TYR	CA-CB-CG	6.17	125.13	113.40
1	B	7	THR	N-CA-CB	-6.13	98.66	110.30
1	C	82	ARG	NE-CZ-NH2	6.08	123.34	120.30
1	C	28	TRP	CB-CG-CD1	-6.03	119.17	127.00
1	B	143	LEU	CA-CB-CG	6.00	129.09	115.30
1	A	85	VAL	CG1-CB-CG2	5.98	120.47	110.90
1	A	76	LEU	CA-CB-CG	5.92	128.93	115.30
1	C	138	ARG	NE-CZ-NH2	5.92	123.26	120.30
1	C	83	ILE	CA-CB-CG2	-5.85	99.19	110.90
1	C	25	GLN	N-CA-C	5.84	126.77	111.00
1	A	28	TRP	CB-CG-CD1	-5.81	119.45	127.00
1	C	29	LEU	CA-CB-CG	5.81	128.66	115.30
1	A	151	TYR	CB-CG-CD1	-5.78	117.53	121.00
1	C	149	GLN	N-CA-CB	-5.77	100.21	110.60
1	C	83	ILE	CA-CB-CG1	5.76	121.94	111.00
1	B	88	GLN	N-CA-CB	-5.75	100.24	110.60
1	A	148	GLY	O-C-N	5.70	131.81	122.70
1	B	32	ARG	NE-CZ-NH2	5.67	123.14	120.30
1	C	11	LYS	N-CA-C	-5.63	95.81	111.00
1	B	69	CYS	CA-CB-SG	-5.61	103.90	114.00
1	B	87	TYR	CA-CB-CG	5.56	123.97	113.40
1	A	30	ASN	CA-C-N	5.56	129.43	117.20
1	A	74	VAL	CG1-CB-CG2	-5.54	102.04	110.90
1	A	103	ARG	NE-CZ-NH1	-5.53	117.53	120.30
1	A	146	GLU	CA-C-N	5.53	129.35	117.20
1	A	114	TRP	CB-CG-CD1	-5.52	119.83	127.00
1	B	149	GLN	CG-CD-NE2	5.52	129.94	116.70
1	A	28	TRP	CG-CD1-NE1	-5.50	104.60	110.10
1	A	131	ARG	CA-CB-CG	5.42	125.33	113.40
1	C	123	VAL	CA-CB-CG2	-5.42	102.77	110.90
1	A	107	GLU	N-CA-C	5.40	125.59	111.00
1	C	37	LEU	N-CA-C	-5.40	96.41	111.00
1	B	114	TRP	CG-CD1-NE1	-5.40	104.70	110.10
1	A	28	TRP	CG-CD2-CE3	5.34	138.71	133.90
1	C	15	HIS	CA-CB-CG	-5.33	104.54	113.60
1	B	73	HIS	O-C-N	5.32	131.21	122.70
1	C	87	TYR	CB-CG-CD1	5.32	124.19	121.00
1	C	53	GLU	N-CA-C	-5.32	96.65	111.00
1	C	32	ARG	NE-CZ-NH1	-5.31	117.64	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	146	GLU	CA-CB-CG	5.30	125.07	113.40
1	B	28	TRP	CG-CD2-CE3	5.29	138.66	133.90
1	B	56	TYR	CA-CB-CG	5.22	123.33	113.40
1	C	110	GLU	CA-C-N	-5.20	105.76	117.20
1	C	44	ARG	O-C-N	-5.17	114.43	122.70
1	B	73	HIS	CA-C-N	-5.15	105.87	117.20
1	A	114	TRP	CG-CD1-NE1	-5.15	104.95	110.10
1	C	16	VAL	CG1-CB-CG2	5.11	119.07	110.90
1	A	81	SER	CA-CB-OG	5.09	124.94	111.20
1	C	72	THR	CA-CB-OG1	-5.09	98.31	109.00
1	A	114	TRP	CG-CD2-CE3	5.08	138.47	133.90
1	A	43	LEU	CB-CG-CD1	-5.08	102.36	111.00
1	A	68	GLY	CA-C-O	5.07	129.73	120.60
1	B	10	ASP	CB-CG-OD2	5.07	122.86	118.30
1	B	91	VAL	CA-CB-CG1	-5.05	103.32	110.90
1	A	91	VAL	CA-CB-CG2	-5.04	103.33	110.90
1	C	15	HIS	N-CA-C	-5.04	97.39	111.00
1	C	127	GLU	CA-CB-CG	-5.04	102.31	113.40
1	B	39	ASN	CB-CG-ND2	5.03	128.78	116.70
1	C	89	THR	CA-CB-CG2	5.03	119.44	112.40
1	A	13	VAL	CB-CA-C	-5.02	101.87	111.40
1	B	141	TYR	CB-CG-CD1	-5.01	117.99	121.00
1	A	22	ALA	CA-C-O	5.01	130.61	120.10

There are no chirality outliers.

All (7) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	105	THR	Peptide
1	A	7	THR	Peptide
1	B	105	THR	Peptide
1	B	7	THR	Peptide
1	C	11	LYS	Peptide
1	C	146	GLU	Peptide
1	C	7	THR	Peptide

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	1184	0	1181	49	0
1	B	1184	0	1181	67	0
1	C	1184	0	1181	56	0
All	All	3552	0	3543	151	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 (151) 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:A:8:PRO:HG3	1:C:54:GLY:HA2	1.49	0.93
1:B:94:LEU:HA	1:C:148:GLY:HA3	1.60	0.82
1:B:98:LYS:HD2	1:B:116:GLU:HG3	1.65	0.79
1:A:149:GLN:HG3	1:A:150:VAL:H	1.47	0.79
1:C:53:GLU:HG3	1:C:127:GLU:HG3	1.67	0.76
1:B:60:SER:HB2	1:B:80:ILE:HD11	1.69	0.75
1:B:106:PRO:HB2	1:B:111:ALA:HB2	1.69	0.73
1:A:15:HIS:O	1:A:35:ALA:HA	1.88	0.72
1:C:82:ARG:HB2	1:C:93:LEU:HD11	1.71	0.70
1:A:77:THR:HG22	1:A:97:ILE:HG23	1.72	0.70
1:B:77:THR:HG22	1:B:97:ILE:HG23	1.74	0.70
1:C:65:LYS:HD3	1:C:143:LEU:HD12	1.72	0.69
1:C:30:ASN:ND2	1:C:31:ARG:HH22	1.91	0.69
1:A:149:GLN:HG3	1:A:150:VAL:HG12	1.75	0.68
1:A:149:GLN:HG3	1:A:150:VAL:N	2.09	0.67
1:C:47:GLN:OE1	1:C:131:ARG:HB3	1.95	0.67
1:A:146:GLU:O	1:A:149:GLN:HG2	1.97	0.65
1:B:146:GLU:HG3	1:B:149:GLN:HB2	1.79	0.64
1:C:16:VAL:HG12	1:C:152:PHE:HB3	1.79	0.64
1:B:144:PHE:HA	1:B:146:GLU:HB3	1.78	0.63
1:B:103:ARG:HD2	1:B:104:GLU:H	1.62	0.63
1:A:30:ASN:HA	1:A:35:ALA:HB1	1.81	0.63
1:C:19:ASN:OD1	1:C:21:GLN:HB2	1.98	0.62
1:A:87:TYR:HD2	1:A:87:TYR:H	1.47	0.61
1:A:84:ALA:HA	1:A:130:ASP:HB2	1.82	0.60
1:A:62:VAL:HG12	1:A:150:VAL:HG23	1.83	0.59
1:A:69:CYS:O	1:A:105:THR:HA	2.01	0.59
1:B:48:LEU:O	1:B:131:ARG:HA	2.02	0.59
1:B:94:LEU:CA	1:C:148:GLY:HA3	2.33	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:132:LEU:HD13	1:B:154:ILE:HG21	1.86	0.58
1:C:69:CYS:O	1:C:105:THR:HA	2.04	0.58
1:C:14:ALA:HA	1:C:36:LEU:O	2.04	0.57
1:B:65:LYS:HD2	1:B:143:LEU:HD23	1.85	0.57
1:A:67:GLN:HA	1:A:113:PRO:HA	1.85	0.57
1:A:10:ASP:HA	1:A:39:ASN:ND2	2.20	0.57
1:C:80:ILE:HA	1:C:133:SER:O	2.05	0.56
1:A:9:SER:HB3	1:A:11:LYS:HG2	1.88	0.56
1:B:84:ALA:HB3	1:B:88:GLN:HA	1.88	0.56
1:A:155:ILE:HD13	1:C:157:LEU:HD12	1.87	0.56
1:A:106:PRO:HB2	1:A:110:GLU:HA	1.87	0.55
1:A:138:ARG:NH2	1:A:141:TYR:CE1	2.74	0.55
1:A:64:PHE:HA	1:A:141:TYR:O	2.05	0.55
1:A:9:SER:CB	1:A:11:LYS:HG2	2.37	0.55
1:B:49:VAL:HA	1:B:130:ASP:O	2.06	0.55
1:B:96:ALA:HB3	1:B:118:ILE:HG21	1.88	0.54
1:C:30:ASN:O	1:C:31:ARG:HB2	2.07	0.54
1:B:65:LYS:HD2	1:B:143:LEU:HB2	1.90	0.54
1:C:17:VAL:HG23	1:C:29:LEU:HB3	1.89	0.54
1:A:13:VAL:HG12	1:A:155:ILE:HB	1.89	0.54
1:C:19:ASN:HD22	1:C:29:LEU:HD23	1.73	0.54
1:A:58:ILE:O	1:A:121:GLY:HA2	2.08	0.54
1:B:60:SER:CB	1:B:80:ILE:HD11	2.38	0.53
1:A:14:ALA:HB2	1:A:41:VAL:HG11	1.89	0.53
1:A:102:GLN:NE2	1:B:114:TRP:HB3	2.24	0.53
1:C:15:HIS:O	1:C:35:ALA:HA	2.08	0.53
1:A:112:LYS:HG3	1:C:102:GLN:HE22	1.75	0.53
1:A:94:LEU:HA	1:B:148:GLY:HA3	1.90	0.52
1:A:57:LEU:HB2	1:A:157:LEU:HD11	1.92	0.52
1:B:94:LEU:HA	1:C:148:GLY:CA	2.36	0.52
1:C:84:ALA:O	1:C:88:GLN:HA	2.10	0.52
1:B:94:LEU:HB3	1:B:120:LEU:HG	1.90	0.52
1:B:13:VAL:HG23	1:B:155:ILE:HD13	1.90	0.52
1:B:67:GLN:HA	1:B:113:PRO:HA	1.91	0.52
1:C:79:THR:HG23	1:C:95:SER:HB3	1.90	0.52
1:A:73:HIS:CE1	1:B:112:LYS:HB3	2.45	0.52
1:B:73:HIS:HE1	1:B:102:GLN:HA	1.75	0.51
1:B:103:ARG:HD2	1:B:104:GLU:N	2.25	0.51
1:C:65:LYS:HE2	1:C:143:LEU:HA	1.91	0.51
1:B:80:ILE:HA	1:B:133:SER:O	2.11	0.51
1:C:10:ASP:HA	1:C:39:ASN:ND2	2.26	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:106:PRO:CB	1:B:111:ALA:HB2	2.40	0.51
1:A:103:ARG:HD2	1:B:104:GLU:OE1	2.12	0.50
1:A:136:ILE:HD11	1:A:139:PRO:HA	1.92	0.50
1:C:78:HIS:O	1:C:95:SER:HA	2.11	0.50
1:C:7:THR:HA	1:C:8:PRO:O	2.12	0.50
1:B:43:LEU:HA	1:B:47:GLN:O	2.12	0.50
1:C:98:LYS:HD3	1:C:116:GLU:HG3	1.93	0.49
1:A:59:TYR:OH	1:C:123:VAL:HB	2.12	0.49
1:B:14:ALA:HB2	1:B:41:VAL:HG21	1.93	0.49
1:C:55:LEU:HD12	1:C:125:GLN:HB2	1.95	0.49
1:B:146:GLU:HB2	1:B:149:GLN:HB2	1.94	0.49
1:B:47:GLN:HG2	1:B:133:SER:HB3	1.94	0.48
1:C:84:ALA:HA	1:C:130:ASP:OD2	2.12	0.48
1:B:104:GLU:HB3	1:B:106:PRO:HG3	1.95	0.48
1:B:79:THR:HG22	1:B:95:SER:HB3	1.96	0.48
1:C:22:ALA:O	1:C:24:GLY:N	2.47	0.48
1:A:92:ASN:HB3	1:B:147:SER:OG	2.13	0.48
1:C:30:ASN:ND2	1:C:31:ARG:NH2	2.61	0.48
1:C:68:GLY:HA2	1:C:106:PRO:HG2	1.96	0.48
1:B:102:GLN:NE2	1:C:114:TRP:HB3	2.28	0.48
1:A:61:GLN:O	1:A:150:VAL:HA	2.14	0.48
1:A:72:THR:OG1	1:B:112:LYS:HE3	2.14	0.47
1:B:144:PHE:C	1:B:146:GLU:H	2.17	0.47
1:A:84:ALA:HA	1:A:130:ASP:CB	2.43	0.47
1:B:125:GLN:OE1	1:C:6:ARG:N	2.48	0.47
1:A:30:ASN:O	1:A:31:ARG:HG2	2.14	0.47
1:B:146:GLU:HB2	1:B:149:GLN:OE1	2.14	0.47
1:B:63:LEU:HD12	1:B:149:GLN:CD	2.35	0.47
1:B:19:ASN:HB2	1:B:29:LEU:HD13	1.97	0.47
1:B:69:CYS:HA	1:B:70:PRO:HD3	1.72	0.47
1:A:63:LEU:HG	1:A:149:GLN:HB2	1.97	0.47
1:C:17:VAL:HB	1:C:32:ARG:HH11	1.80	0.46
1:C:84:ALA:O	1:C:88:GLN:HG2	2.16	0.46
1:C:94:LEU:HB3	1:C:120:LEU:HD22	1.97	0.46
1:B:73:HIS:CE1	1:B:101:CYS:O	2.69	0.46
1:C:98:LYS:HD2	1:C:118:ILE:HG12	1.97	0.46
1:C:107:GLU:HG3	1:C:108:GLY:H	1.80	0.46
1:C:50:VAL:HG21	1:C:126:LEU:HD13	1.97	0.46
1:B:10:ASP:HA	1:B:39:ASN:OD1	2.16	0.45
1:C:16:VAL:HG22	1:C:29:LEU:O	2.17	0.45
1:A:58:ILE:HD11	1:A:126:LEU:HD11	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:61:GLN:O	1:B:150:VAL:HA	2.17	0.45
1:A:70:PRO:O	1:A:105:THR:HG23	2.17	0.45
1:C:56:TYR:HB2	1:C:126:LEU:HD12	1.98	0.45
1:C:57:LEU:O	1:C:154:ILE:HA	2.17	0.45
1:B:146:GLU:CG	1:B:149:GLN:HB2	2.44	0.45
1:A:92:ASN:HB3	1:B:147:SER:CB	2.46	0.45
1:B:12:PRO:HB3	1:B:51:PRO:HG3	2.00	0.44
1:B:58:ILE:HG23	1:B:154:ILE:HG22	1.99	0.44
1:B:80:ILE:HG13	1:B:120:LEU:HD23	1.99	0.44
1:A:83:ILE:HA	1:A:89:THR:O	2.17	0.44
1:B:47:GLN:OE1	1:B:131:ARG:HD3	2.18	0.43
1:A:95:SER:N	1:B:148:GLY:O	2.51	0.43
1:B:79:THR:HG22	1:B:95:SER:CB	2.48	0.43
1:B:136:ILE:HD13	1:B:142:LEU:HD21	2.00	0.43
1:A:82:ARG:HB2	1:A:93:LEU:HD11	2.01	0.43
1:B:30:ASN:HA	1:B:35:ALA:HB1	2.01	0.43
1:A:8:PRO:HD2	1:C:125:GLN:OE1	2.19	0.43
1:C:32:ARG:NH1	1:C:146:GLU:HB3	2.34	0.42
1:A:12:PRO:HB3	1:A:41:VAL:HG23	2.01	0.42
1:A:30:ASN:CA	1:A:35:ALA:HB1	2.48	0.42
1:B:102:GLN:CD	1:C:114:TRP:HB3	2.40	0.42
1:C:61:GLN:HA	1:C:118:ILE:O	2.20	0.42
1:A:19:ASN:HA	1:A:29:LEU:HD22	2.00	0.42
1:B:157:LEU:HA	1:B:157:LEU:HD23	1.79	0.42
1:C:110:GLU:OE1	1:C:111:ALA:HB3	2.20	0.42
1:B:20:PRO:HA	1:B:144:PHE:CD2	2.54	0.42
1:A:59:TYR:CZ	1:C:123:VAL:HB	2.55	0.41
1:C:52:SER:O	1:C:128:LYS:HB2	2.20	0.41
1:C:82:ARG:NH2	1:C:130:ASP:OD1	2.52	0.41
1:B:19:ASN:HA	1:B:20:PRO:HD3	1.91	0.41
1:C:32:ARG:O	1:C:34:ASN:N	2.54	0.41
1:B:83:ILE:HG23	1:B:131:ARG:HB2	2.03	0.41
1:B:146:GLU:HG2	1:B:147:SER:O	2.21	0.41
1:A:85:VAL:H	1:A:130:ASP:HB3	1.86	0.41
1:B:15:HIS:O	1:B:35:ALA:HA	2.21	0.41
1:C:83:ILE:CG2	1:C:90:LYS:HG2	2.51	0.41
1:B:59:TYR:HA	1:B:120:LEU:O	2.21	0.41
1:B:146:GLU:HB2	1:B:149:GLN:CD	2.42	0.41
1:B:97:ILE:O	1:C:115:TYR:HB3	2.22	0.40
1:C:10:ASP:HA	1:C:39:ASN:HD21	1.85	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	150/157 (96%)	121 (81%)	19 (13%)	10 (7%)	1	1
1	B	150/157 (96%)	115 (77%)	25 (17%)	10 (7%)	1	1
1	C	150/157 (96%)	111 (74%)	25 (17%)	14 (9%)	1	0
All	All	450/471 (96%)	347 (77%)	69 (15%)	34 (8%)	1	1

All (34) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	72	THR
1	A	100	PRO
1	A	104	GLU
1	A	107	GLU
1	B	84	ALA
1	B	88	GLN
1	B	104	GLU
1	C	23	GLU
1	C	53	GLU
1	C	75	LEU
1	C	106	PRO
1	A	70	PRO
1	A	75	LEU
1	B	24	GLY
1	B	37	LEU
1	B	75	LEU
1	C	26	LEU
1	C	31	ARG
1	C	33	ALA
1	C	52	SER
1	C	108	GLY
1	C	111	ALA
1	A	9	SER
1	A	32	ARG

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Mol	Chain	Res	Type
1	A	110	GLU
1	B	8	PRO
1	B	9	SER
1	B	68	GLY
1	C	51	PRO
1	C	145	ALA
1	A	155	ILE
1	C	39	ASN
1	C	147	SER
1	B	51	PRO

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	127/133 (96%)	104 (82%)	23 (18%)	2 3
1	B	127/133 (96%)	101 (80%)	26 (20%)	1 2
1	C	127/133 (96%)	96 (76%)	31 (24%)	1 1
All	All	381/399 (96%)	301 (79%)	80 (21%)	1 2

All (80) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	VAL
1	A	17	VAL
1	A	19	ASN
1	A	32	ARG
1	A	37	LEU
1	A	39	ASN
1	A	43	LEU
1	A	48	LEU
1	A	52	SER
1	A	55	LEU
1	A	57	LEU
1	A	65	LYS

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Mol	Chain	Res	Type
1	A	70	PRO
1	A	71	SER
1	A	73	HIS
1	A	75	LEU
1	A	76	LEU
1	A	83	ILE
1	A	85	VAL
1	A	87	TYR
1	A	105	THR
1	A	131	ARG
1	A	140	ASP
1	B	17	VAL
1	B	19	ASN
1	B	23	GLU
1	B	29	LEU
1	B	30	ASN
1	B	41	VAL
1	B	43	LEU
1	B	44	ARG
1	B	55	LEU
1	B	63	LEU
1	B	65	LYS
1	B	67	GLN
1	B	73	HIS
1	B	81	SER
1	B	82	ARG
1	B	83	ILE
1	B	85	VAL
1	B	86	SER
1	B	104	GLU
1	B	105	THR
1	B	107	GLU
1	B	110	GLU
1	B	120	LEU
1	B	126	LEU
1	B	127	GLU
1	B	146	GLU
1	C	12	PRO
1	C	16	VAL
1	C	17	VAL
1	C	31	ARG
1	C	32	ARG

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Mol	Chain	Res	Type
1	C	36	LEU
1	C	39	ASN
1	C	41	VAL
1	C	44	ARG
1	C	48	LEU
1	C	51	PRO
1	C	55	LEU
1	C	62	VAL
1	C	67	GLN
1	C	72	THR
1	C	73	HIS
1	C	76	LEU
1	C	82	ARG
1	C	83	ILE
1	C	89	THR
1	C	91	VAL
1	C	99	SER
1	C	101	CYS
1	C	103	ARG
1	C	106	PRO
1	C	112	LYS
1	C	127	GLU
1	C	131	ARG
1	C	132	LEU
1	C	133	SER
1	C	140	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	19	ASN
1	A	61	GLN
1	A	78	HIS
1	A	102	GLN
1	B	125	GLN
1	C	30	ASN
1	C	78	HIS

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

There are no ligands in this entry.

5.7 Other polymers

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

EDS was not executed - this section will therefore be empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates

EDS was not executed - this section will therefore be empty.

6.4 Ligands

EDS was not executed - this section will therefore be empty.

6.5 Other polymers

EDS was not executed - this section will therefore be empty.