



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:03 PM GMT

PDB ID : 3VXT  
Title : T36-5 TCR specific for HLA-A24-Nef134-10  
Authors : Shimizu, A.; Fukai, S.; Yamagata, A.; Iwamoto, A.  
Deposited on : 2012-09-20  
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](mailto: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.

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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

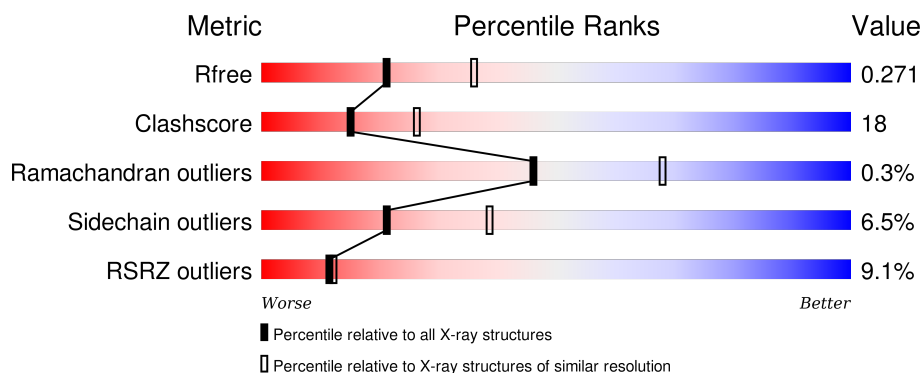
# 1 Overall quality at a glance

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  $\geq 3$ , 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	205	<div> <div>10%</div> <div>59%</div> <div>34%</div> <div>• •</div> </div>
1	C	205	<div> <div>13%</div> <div>62%</div> <div>32%</div> <div>• •</div> </div>
2	B	242	<div> <div>5%</div> <div>66%</div> <div>32%</div> <div>•</div> </div>
2	D	242	<div> <div>9%</div> <div>73%</div> <div>24%</div> <div>•</div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7080 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 T36-5 TCR alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	199	Total	C	N	O	S	0	0	0
			1553	968	257	321	7			
1	A	199	Total	C	N	O	S	0	0	0
			1553	968	257	321	7			

- Molecule 2 is a protein called T36-5 TCR beta chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	241	Total	C	N	O	S	0	0	0
			1933	1217	336	372	8			
2	B	241	Total	C	N	O	S	0	0	0
			1933	1217	336	372	8			

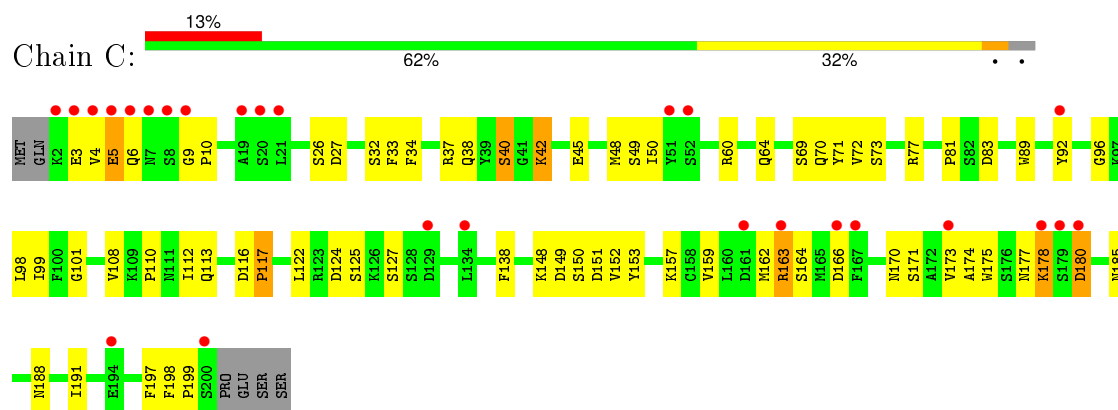
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	C	15	Total	O	0	0
			15	15		
3	D	50	Total	O	0	0
			50	50		
3	A	12	Total	O	0	0
			12	12		
3	B	31	Total	O	0	0
			31	31		

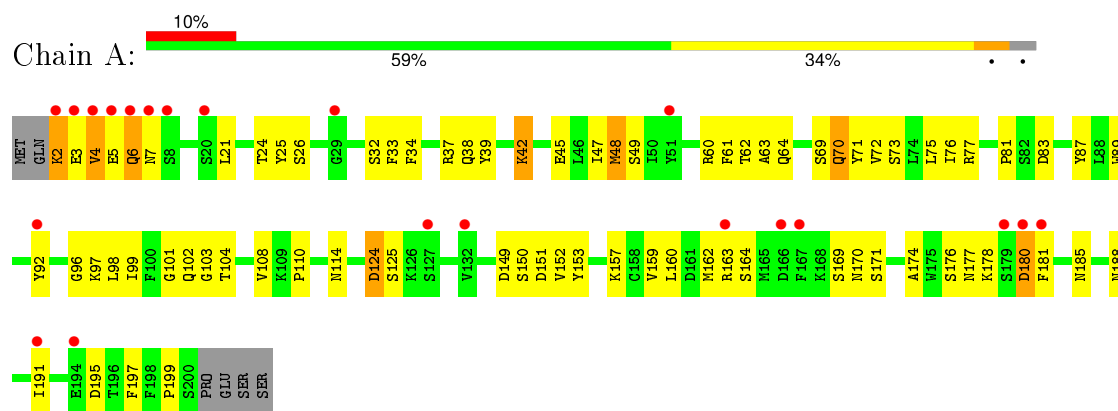
### 3 Residue-property plots

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.

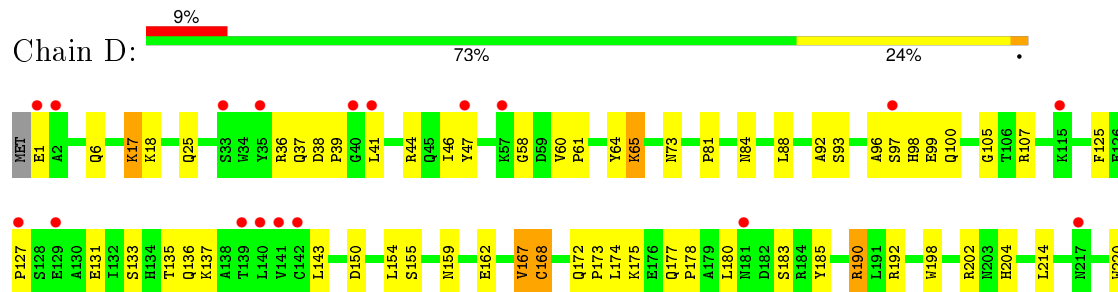
#### • Molecule 1: T36-5 TCR alpha chain

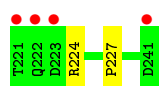


#### • Molecule 1: T36-5 TCR alpha chain

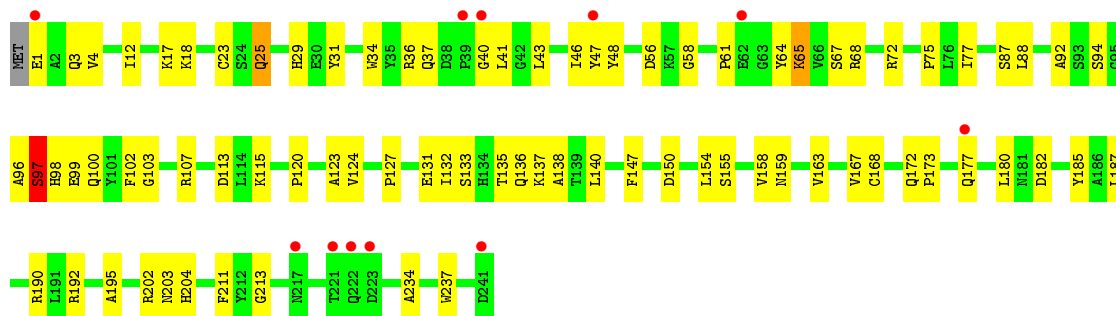


#### • Molecule 2: T36-5 TCR beta chain





• Molecule 2: T36-5 TCR beta chain



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	45.77Å 65.04Å 92.28Å 91.30° 94.19° 105.93°	Depositor
Resolution (Å)	50.00 – 2.50 29.61 – 2.50	Depositor EDS
% Data completeness (in resolution range)	91.5 (50.00-2.50) 90.7 (29.61-2.50)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.04	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.08 (at 2.51Å)	Xtriage
Refinement program	CNS 1.3	Depositor
R, $R_{free}$	0.230 , 0.271 0.231 , 0.271	Depositor DCC
$R_{free}$ test set	3220 reflections (9.99%)	DCC
Wilson B-factor (Å <sup>2</sup> )	57.3	Xtriage
Anisotropy	0.298	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 63.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 32246 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	7080	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	69.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.03% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>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	0.65	4/1587 (0.3%)	0.68	0/2149
1	C	0.42	0/1587	0.72	1/2149 (0.0%)
2	B	0.41	0/1986	0.69	1/2705 (0.0%)
2	D	0.43	0/1986	0.70	1/2705 (0.0%)
All	All	0.48	4/7146 (0.1%)	0.70	3/9708 (0.0%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	39	TYR	CD2-CE2	-7.77	1.27	1.39
1	A	39	TYR	CE2-CZ	-7.60	1.28	1.38
1	A	39	TYR	CD1-CE1	-7.07	1.28	1.39
1	A	39	TYR	CE1-CZ	-5.59	1.31	1.38

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	6	GLN	N-CA-C	-8.19	88.88	111.00
2	B	97	SER	N-CA-C	-6.08	94.59	111.00
2	D	174	LEU	CA-CB-CG	5.15	127.14	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	1553	0	1461	75	0
1	C	1553	0	1461	57	0
2	B	1933	0	1845	72	0
2	D	1933	0	1845	54	0
3	A	12	0	0	6	0
3	B	31	0	0	1	0
3	C	15	0	0	2	0
3	D	50	0	0	8	0
All	All	7080	0	6612	238	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 18.

All (238) 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:177:GLN:HB3	2:D:180:LEU:HD13	1.21	1.18
2:D:177:GLN:CB	2:D:180:LEU:HD13	1.85	1.07
1:C:110:PRO:HG3	1:C:159:VAL:HG21	1.43	1.00
2:B:3:GLN:HG2	2:B:4:VAL:H	1.31	0.95
2:D:177:GLN:HB3	2:D:180:LEU:CD1	1.97	0.94
2:B:92:ALA:HB1	2:B:100:GLN:HG3	1.52	0.91
1:A:6:GLN:CD	1:A:7:ASN:H	1.85	0.80
1:C:110:PRO:HG3	1:C:159:VAL:CG2	2.14	0.78
1:A:6:GLN:NE2	1:A:7:ASN:H	1.80	0.78
1:C:178:LYS:HG2	1:C:180:ASP:HB2	1.66	0.77
2:D:98:HIS:HD2	3:D:345:HOH:O	1.66	0.76
1:A:97:LYS:HE2	2:B:48:TYR:OH	1.88	0.74
1:C:166:ASP:HB2	3:C:314:HOH:O	1.89	0.73
2:D:73:ASN:HB3	3:D:311:HOH:O	1.88	0.73
2:B:31:TYR:HB3	2:B:94:SER:HB3	1.70	0.72
2:B:3:GLN:HG2	2:B:4:VAL:N	2.05	0.72
1:A:42:LYS:HD2	1:A:42:LYS:N	2.04	0.72
1:C:77:ARG:HG3	1:C:77:ARG:HH11	1.55	0.71
2:B:3:GLN:CG	2:B:4:VAL:H	2.03	0.71
1:A:178:LYS:HG2	1:A:180:ASP:HB2	1.71	0.71
1:C:197:PHE:CZ	1:C:199:PRO:HG3	2.26	0.70
1:A:4:VAL:HG22	1:A:5:GLU:HG2	1.72	0.69
2:D:202:ARG:HG3	2:D:202:ARG:HH11	1.55	0.69
1:A:38:GLN:NE2	2:B:37:GLN:HE22	1.92	0.68
2:B:135:THR:O	2:B:136:GLN:HB2	1.93	0.68
2:B:58:GLY:O	2:B:61:PRO:HD3	1.94	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:3:GLU:HG3	1:C:4:VAL:H	1.58	0.68
2:D:180:LEU:HD12	2:D:180:LEU:N	2.09	0.67
1:C:113:GLN:HA	1:C:113:GLN:HE21	1.59	0.67
1:A:162:MET:HE1	2:B:192:ARG:HD3	1.76	0.66
1:A:2:LYS:NZ	1:A:3:GLU:H	1.93	0.66
1:C:191:ILE:O	1:C:191:ILE:HG13	1.94	0.66
2:D:98:HIS:CD2	3:D:345:HOH:O	2.46	0.66
1:C:5:GLU:OE2	1:C:5:GLU:HA	1.95	0.66
1:A:62:THR:OG1	1:A:77:ARG:NH2	2.30	0.65
1:C:42:LYS:HD2	1:C:42:LYS:N	2.11	0.65
2:B:88:LEU:HD13	2:B:107:ARG:HG3	1.79	0.65
1:C:113:GLN:HG2	3:C:311:HOH:O	1.96	0.64
2:D:167:VAL:HA	2:D:190:ARG:O	1.98	0.64
2:D:60:VAL:HG23	2:D:60:VAL:O	1.98	0.63
1:C:33:PHE:O	1:C:50:ILE:HG22	1.98	0.63
2:B:202:ARG:HH11	2:B:202:ARG:HG3	1.64	0.63
2:D:159:ASN:ND2	2:D:204:HIS:H	1.97	0.62
2:D:177:GLN:HB2	2:D:180:LEU:HD13	1.79	0.62
1:A:176:SER:HA	3:A:305:HOH:O	1.99	0.62
1:A:6:GLN:HB2	1:A:102:GLN:HB2	1.82	0.62
1:C:77:ARG:NH1	1:C:77:ARG:HG3	2.14	0.62
2:B:127:PRO:HG2	2:B:138:ALA:HB1	1.82	0.62
2:B:100:GLN:HB3	2:B:102:PHE:CE2	2.34	0.61
2:B:167:VAL:HA	2:B:190:ARG:O	2.00	0.61
2:D:92:ALA:HB1	2:D:100:GLN:HG3	1.83	0.61
1:A:81:PRO:HA	1:A:108:VAL:HB	1.83	0.61
2:D:58:GLY:O	2:D:61:PRO:HD3	2.01	0.60
1:A:159:VAL:O	2:B:168:CYS:SG	2.60	0.59
1:C:64:GLN:HB2	1:C:73:SER:OG	2.02	0.59
1:A:110:PRO:HG3	1:A:159:VAL:HG21	1.84	0.59
1:C:89:TRP:CZ2	1:C:101:GLY:HA3	2.37	0.59
1:A:2:LYS:HZ2	1:A:3:GLU:H	1.51	0.59
1:C:113:GLN:CA	1:C:113:GLN:HE21	2.16	0.58
1:A:152:VAL:HG13	1:A:152:VAL:O	2.03	0.58
1:A:89:TRP:CZ2	1:A:101:GLY:HA3	2.39	0.58
1:C:149:ASP:HB3	1:C:152:VAL:HG12	1.84	0.57
1:C:197:PHE:CD1	1:C:198:PHE:N	2.73	0.57
1:C:197:PHE:CE2	1:C:199:PRO:HG3	2.39	0.57
1:A:64:GLN:HB2	1:A:73:SER:OG	2.05	0.57
1:C:113:GLN:HA	1:C:113:GLN:NE2	2.19	0.57
2:D:159:ASN:HD21	2:D:204:HIS:H	1.51	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:34:PHE:HD1	1:C:49:SER:HG	1.52	0.57
1:C:60:ARG:NH2	1:C:83:ASP:OD2	2.37	0.57
1:A:97:LYS:HE2	2:B:48:TYR:CZ	2.39	0.57
2:B:97:SER:O	2:B:99:GLU:HG2	2.05	0.56
2:B:154:LEU:HD23	2:B:155:SER:N	2.19	0.56
1:A:60:ARG:HA	1:A:77:ARG:NH1	2.21	0.56
1:A:160:LEU:HB3	2:B:168:CYS:HB2	1.87	0.56
1:C:185:ASN:HA	1:C:188:ASN:ND2	2.20	0.56
1:A:6:GLN:HE21	1:A:6:GLN:N	2.04	0.56
2:B:135:THR:HG21	2:B:192:ARG:NH2	2.20	0.56
1:C:173:VAL:HG12	1:C:174:ALA:N	2.21	0.56
1:C:153:TYR:O	1:C:174:ALA:HA	2.06	0.56
2:B:4:VAL:HB	2:B:103:GLY:HA2	1.88	0.55
1:C:98:LEU:HD13	2:D:100:GLN:OE1	2.07	0.55
2:B:120:PRO:HB3	2:B:147:PHE:HB3	1.88	0.55
2:D:135:THR:HG21	2:D:192:ARG:NH2	2.22	0.55
1:A:6:GLN:CD	1:A:7:ASN:N	2.57	0.55
1:C:185:ASN:HA	1:C:188:ASN:HD21	1.72	0.55
1:C:175:TRP:CD2	2:D:143:LEU:HD21	2.41	0.54
2:D:36:ARG:HB3	2:D:46:ILE:HD11	1.90	0.54
1:A:34:PHE:HD1	1:A:49:SER:HG	1.55	0.54
1:A:180:ASP:HB3	3:A:312:HOH:O	2.06	0.54
2:D:220:TRP:CZ2	2:D:227:PRO:HD3	2.44	0.53
1:A:6:GLN:CG	1:A:7:ASN:N	2.71	0.53
1:A:33:PHE:CD1	1:A:72:VAL:HG21	2.42	0.53
2:D:202:ARG:NH1	2:D:202:ARG:HG3	2.20	0.53
1:A:38:GLN:HE22	2:B:37:GLN:HE22	1.55	0.53
1:A:2:LYS:HB2	1:A:2:LYS:NZ	2.24	0.53
2:B:177:GLN:HB3	2:B:180:LEU:HD13	1.91	0.53
2:B:150:ASP:HB3	2:B:185:TYR:CD2	2.42	0.53
2:B:12:ILE:CD1	2:B:213:GLY:HA2	2.38	0.53
2:D:88:LEU:HD13	2:D:107:ARG:HG3	1.91	0.53
1:C:117:PRO:HG3	1:C:138:PHE:HA	1.91	0.52
2:B:127:PRO:CG	2:B:138:ALA:HB1	2.38	0.52
2:B:135:THR:HG21	2:B:192:ARG:HH21	1.74	0.52
1:A:149:ASP:HB3	1:A:152:VAL:HG12	1.91	0.51
1:A:69:SER:O	1:A:70:GLN:HB2	2.09	0.51
2:B:36:ARG:HB3	2:B:46:ILE:HD11	1.93	0.51
1:C:149:ASP:HB3	1:C:152:VAL:CG1	2.41	0.51
1:A:157:LYS:HA	1:A:171:SER:O	2.11	0.50
1:A:98:LEU:HD13	2:B:100:GLN:OE1	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:23:CYS:HB2	2:B:34:TRP:CZ2	2.46	0.50
1:A:60:ARG:NH2	1:A:83:ASP:OD2	2.45	0.50
1:A:110:PRO:HG3	1:A:159:VAL:CG2	2.42	0.50
1:A:153:TYR:O	1:A:174:ALA:HA	2.11	0.50
1:A:149:ASP:HB3	1:A:152:VAL:CG1	2.42	0.50
2:B:25:GLN:OE1	2:B:29:HIS:N	2.28	0.50
1:A:159:VAL:HG22	1:A:170:ASN:OD1	2.12	0.49
1:A:185:ASN:HA	1:A:188:ASN:ND2	2.28	0.49
2:D:162:GLU:HB2	3:D:339:HOH:O	2.12	0.49
1:C:4:VAL:HG22	1:C:5:GLU:N	2.27	0.49
1:C:69:SER:O	1:C:70:GLN:HB2	2.11	0.49
2:B:173:PRO:HG3	2:B:187:LEU:HD13	1.95	0.49
2:B:127:PRO:HD3	2:B:140:LEU:HG	1.93	0.49
1:A:181:PHE:HA	3:A:306:HOH:O	2.12	0.49
2:D:214:LEU:CD1	2:D:227:PRO:HG2	2.43	0.48
1:A:124:ASP:OD1	1:A:124:ASP:N	2.44	0.48
1:C:5:GLU:CA	1:C:5:GLU:OE2	2.60	0.48
1:A:21:LEU:HD21	1:A:104:THR:HG21	1.94	0.48
2:B:202:ARG:NH1	2:B:202:ARG:HG3	2.28	0.48
2:B:64:TYR:O	2:B:65:LYS:HD2	2.14	0.48
1:A:92:TYR:OH	1:A:96:GLY:HA2	2.13	0.48
2:B:72:ARG:HH11	2:B:72:ARG:HG3	1.79	0.48
2:B:113:ASP:C	2:B:115:LYS:H	2.18	0.48
1:C:38:GLN:HE22	2:D:37:GLN:HE22	1.60	0.48
1:C:38:GLN:NE2	2:D:37:GLN:HE22	2.12	0.48
1:A:97:LYS:O	1:A:99:ILE:HD12	2.14	0.48
1:C:157:LYS:HA	1:C:171:SER:O	2.14	0.47
1:A:103:GLY:N	3:A:304:HOH:O	2.48	0.47
1:C:33:PHE:CD1	1:C:72:VAL:HG21	2.50	0.47
2:B:131:GLU:HG2	2:B:131:GLU:O	2.15	0.47
2:B:25:GLN:HB3	2:B:25:GLN:HE21	1.56	0.47
2:B:123:ALA:HB2	3:B:322:HOH:O	2.15	0.47
1:A:97:LYS:CE	2:B:48:TYR:OH	2.61	0.47
2:D:172:GLN:NE2	2:D:173:PRO:CD	2.78	0.47
2:B:135:THR:CG2	2:B:137:LYS:HB2	2.44	0.47
2:D:135:THR:HG22	2:D:137:LYS:HG3	1.97	0.47
1:A:37:ARG:HB2	1:A:47:ILE:HD13	1.97	0.47
1:A:2:LYS:HZ2	1:A:3:GLU:N	2.11	0.46
1:A:25:TYR:CZ	1:A:70:GLN:HG2	2.50	0.46
2:D:64:TYR:O	2:D:65:LYS:HD2	2.15	0.46
1:C:159:VAL:O	2:D:168:CYS:SG	2.74	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:2:LYS:HZ3	1:A:2:LYS:HB2	1.79	0.46
1:A:61:PHE:CD1	1:A:76:ILE:HG12	2.50	0.46
1:A:97:LYS:HE3	1:A:97:LYS:HB2	1.63	0.46
2:B:87:SER:OG	2:B:88:LEU:N	2.49	0.46
2:B:31:TYR:HB3	2:B:94:SER:CB	2.45	0.46
1:A:159:VAL:N	2:B:168:CYS:SG	2.89	0.46
2:B:68:ARG:HG2	2:B:68:ARG:HH11	1.81	0.46
2:D:131:GLU:O	2:D:135:THR:HB	2.16	0.46
2:D:177:GLN:HB2	2:D:183:SER:HB2	1.98	0.45
2:D:96:ALA:O	2:D:97:SER:C	2.55	0.45
2:D:177:GLN:HA	2:D:177:GLN:OE1	2.16	0.45
2:B:97:SER:HB3	2:B:99:GLU:HG2	1.98	0.45
1:A:92:TYR:CZ	1:A:96:GLY:HA2	2.51	0.45
2:B:204:HIS:HB2	2:B:237:TRP:CZ3	2.51	0.45
2:D:127:PRO:HD2	2:D:198:TRP:CZ2	2.51	0.45
2:D:150:ASP:HB3	2:D:185:TYR:CD2	2.51	0.45
2:D:154:LEU:HD23	2:D:155:SER:N	2.31	0.45
2:D:135:THR:O	2:D:136:GLN:HB2	2.15	0.45
2:B:12:ILE:HD11	2:B:213:GLY:HA2	1.98	0.45
2:B:67:SER:O	2:B:75:PRO:HD2	2.17	0.45
2:D:224:ARG:NH1	3:D:340:HOH:O	2.50	0.45
1:A:97:LYS:HE2	2:B:48:TYR:CE2	2.52	0.44
1:C:122:LEU:N	1:C:122:LEU:HD12	2.31	0.44
1:C:26:SER:OG	1:C:27:ASP:N	2.50	0.44
2:B:40:GLY:C	2:B:41:LEU:HD22	2.37	0.44
1:A:191:ILE:O	1:A:191:ILE:HG13	2.17	0.44
1:C:9:GLY:HA2	1:C:10:PRO:HD3	1.70	0.44
2:B:120:PRO:HD3	2:B:211:PHE:CD1	2.52	0.44
1:A:4:VAL:HG22	1:A:5:GLU:N	2.33	0.44
2:B:147:PHE:CE1	2:B:185:TYR:HB2	2.53	0.44
1:C:162:MET:CE	2:D:192:ARG:HD3	2.48	0.44
2:B:3:GLN:CG	2:B:4:VAL:N	2.71	0.44
2:D:175:LYS:HE2	2:D:178:PRO:O	2.18	0.44
2:D:17:LYS:O	2:D:81:PRO:HG2	2.18	0.44
1:A:60:ARG:HA	1:A:77:ARG:HH12	1.82	0.43
1:A:102:GLN:NE2	3:A:308:HOH:O	2.39	0.43
2:B:154:LEU:HD23	2:B:154:LEU:C	2.38	0.43
2:D:172:GLN:NE2	2:D:173:PRO:HD2	2.33	0.43
2:D:131:GLU:HG2	2:D:131:GLU:O	2.17	0.43
1:C:71:TYR:CD1	1:C:71:TYR:C	2.92	0.43
1:A:38:GLN:HE22	2:B:37:GLN:NE2	2.16	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:2:LYS:N	1:A:3:GLU:OE2	2.52	0.43
1:A:71:TYR:CD1	1:A:71:TYR:O	2.71	0.43
2:D:46:ILE:HG22	2:D:47:TYR:HD2	1.83	0.43
1:A:48:MET:HG2	1:A:63:ALA:HB2	2.00	0.43
1:C:37:ARG:HE	1:C:37:ARG:HB3	1.54	0.43
1:C:81:PRO:HA	1:C:108:VAL:HB	2.00	0.43
2:D:84:ASN:H	2:D:84:ASN:ND2	2.16	0.43
1:C:92:TYR:OH	1:C:96:GLY:HA2	2.19	0.42
1:C:112:ILE:HD11	1:C:170:ASN:ND2	2.34	0.42
1:C:4:VAL:HG22	1:C:5:GLU:H	1.83	0.42
2:B:37:GLN:HB2	2:B:43:LEU:HD12	2.01	0.42
1:A:21:LEU:O	1:A:73:SER:HB2	2.19	0.42
2:D:39:PRO:HD2	3:D:332:HOH:O	2.18	0.42
1:A:171:SER:OG	2:B:190:ARG:HD2	2.19	0.42
2:B:96:ALA:O	2:B:98:HIS:CE1	2.72	0.42
2:B:158:VAL:CG2	2:B:163:VAL:HG21	2.50	0.42
1:A:37:ARG:HG3	1:A:87:TYR:CE1	2.55	0.42
2:B:124:VAL:HG23	2:B:234:ALA:HB3	2.01	0.42
1:C:4:VAL:HG22	1:C:5:GLU:CD	2.40	0.42
1:C:163:ARG:HG2	1:C:163:ARG:H	1.53	0.42
2:D:6:GLN:HE22	2:D:105:GLY:HA2	1.84	0.42
1:A:169:SER:HA	3:A:303:HOH:O	2.18	0.42
1:A:177:ASN:N	1:A:177:ASN:OD1	2.53	0.42
1:A:197:PHE:CZ	1:A:199:PRO:HG3	2.55	0.42
1:C:175:TRP:CG	2:D:143:LEU:HD21	2.55	0.42
2:D:172:GLN:NE2	2:D:173:PRO:HD3	2.35	0.42
2:B:77:ILE:HD12	2:B:77:ILE:H	1.84	0.41
1:A:24:THR:HG22	1:A:71:TYR:HB3	2.03	0.41
2:D:125:PHE:HE2	2:D:143:LEU:HB2	1.84	0.41
1:A:71:TYR:CD1	1:A:71:TYR:C	2.94	0.41
1:C:116:ASP:N	1:C:117:PRO:HD3	2.35	0.41
2:B:41:LEU:HD13	2:B:41:LEU:HA	1.93	0.41
1:C:177:ASN:OD1	1:C:177:ASN:N	2.52	0.41
1:C:127:SER:HB3	3:D:341:HOH:O	2.20	0.41
1:C:159:VAL:HG22	1:C:170:ASN:OD1	2.21	0.41
1:A:101:GLY:O	1:A:102:GLN:C	2.59	0.41
2:B:46:ILE:HG22	2:B:47:TYR:HD2	1.84	0.41
2:B:172:GLN:HA	2:B:173:PRO:HD3	1.85	0.41
2:D:38:ASP:HB3	3:D:332:HOH:O	2.20	0.41
2:B:159:ASN:HD21	2:B:203:ASN:HA	1.86	0.41
1:A:24:THR:HA	1:A:71:TYR:HA	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:99:ILE:N	1:C:99:ILE:HD12	2.36	0.41
2:D:41:LEU:HB3	2:D:44:ARG:HD3	2.03	0.41
2:D:1:GLU:O	2:D:1:GLU:OE2	2.39	0.40
2:B:132:ILE:HG23	2:B:195:ALA:HB1	2.03	0.40
2:B:124:VAL:HG23	2:B:234:ALA:CB	2.52	0.40
1:A:195:ASP:OD2	1:A:195:ASP:O	2.39	0.40
2:B:135:THR:HG22	2:B:137:LYS:HB2	2.03	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	197/205 (96%)	178 (90%)	17 (9%)	2 (1%)	19	34
1	C	197/205 (96%)	178 (90%)	18 (9%)	1 (0%)	34	55
2	B	239/242 (99%)	227 (95%)	12 (5%)	0	100	100
2	D	239/242 (99%)	230 (96%)	9 (4%)	0	100	100
All	All	872/894 (98%)	813 (93%)	56 (6%)	3 (0%)	46	68

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	40	SER
1	A	70	GLN
1	A	4	VAL

### 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	177/183 (97%)	161 (91%)	16 (9%)	12	22
1	C	177/183 (97%)	161 (91%)	16 (9%)	12	22
2	B	214/215 (100%)	205 (96%)	9 (4%)	36	62
2	D	214/215 (100%)	204 (95%)	10 (5%)	32	56
All	All	782/796 (98%)	731 (94%)	51 (6%)	21	39

All (51) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	5	GLU
1	C	32	SER
1	C	40	SER
1	C	42	LYS
1	C	45	GLU
1	C	48	MET
1	C	117	PRO
1	C	124	ASP
1	C	125	SER
1	C	148	LYS
1	C	150	SER
1	C	151	ASP
1	C	163	ARG
1	C	164	SER
1	C	178	LYS
1	C	180	ASP
2	D	17	LYS
2	D	18	LYS
2	D	25	GLN
2	D	65	LYS
2	D	93	SER
2	D	99	GLU
2	D	133	SER
2	D	167	VAL
2	D	168	CYS
2	D	190	ARG
1	A	2	LYS
1	A	6	GLN

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Mol	Chain	Res	Type
1	A	26	SER
1	A	32	SER
1	A	42	LYS
1	A	45	GLU
1	A	48	MET
1	A	75	LEU
1	A	114	ASN
1	A	124	ASP
1	A	125	SER
1	A	150	SER
1	A	151	ASP
1	A	163	ARG
1	A	164	SER
1	A	180	ASP
2	B	1	GLU
2	B	17	LYS
2	B	18	LYS
2	B	25	GLN
2	B	56	ASP
2	B	65	LYS
2	B	97	SER
2	B	133	SER
2	B	182	ASP

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

Mol	Chain	Res	Type
1	C	6	GLN
1	C	22	ASN
1	C	64	GLN
1	C	113	GLN
1	C	188	ASN
2	D	6	GLN
2	D	37	GLN
2	D	84	ASN
2	D	151	HIS
2	D	159	ASN
2	D	172	GLN
2	D	203	ASN
2	D	217	ASN
2	D	230	GLN
1	A	6	GLN

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Mol	Chain	Res	Type
1	A	22	ASN
1	A	38	GLN
1	A	111	ASN
1	A	113	GLN
1	A	188	ASN
1	A	189	ASN
2	B	45	GLN
2	B	159	ASN
2	B	172	GLN
2	B	203	ASN
2	B	217	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 [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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	199/205 (97%)	0.80	21 (10%) 8 8	50, 78, 107, 143	0
1	C	199/205 (97%)	0.90	26 (13%) 5 4	41, 71, 95, 131	0
2	B	241/242 (99%)	0.44	11 (4%) 36 41	38, 67, 90, 97	0
2	D	241/242 (99%)	0.45	22 (9%) 11 12	33, 60, 83, 97	0
All	All	880/894 (98%)	0.63	80 (9%) 11 12	33, 69, 96, 143	0

All (80) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	4	VAL	10.6
1	A	4	VAL	9.3
1	C	8	SER	6.2
2	D	1	GLU	6.0
1	C	6	GLN	6.0
1	C	3	GLU	5.6
2	B	241	ASP	5.6
1	A	8	SER	5.5
1	A	3	GLU	5.5
2	B	40	GLY	5.2
1	A	7	ASN	5.0
1	A	166	ASP	4.8
1	C	200	SER	4.6
1	A	6	GLN	4.3
2	D	2	ALA	4.1
1	C	7	ASN	3.9
1	A	5	GLU	3.9
1	C	5	GLU	3.9
2	B	221	THR	3.7
1	A	2	LYS	3.6
1	C	2	LYS	3.5

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Mol	Chain	Res	Type	RSRZ
1	A	179	SER	3.5
2	B	39	PRO	3.4
2	D	241	ASP	3.4
1	C	52	SER	3.4
2	D	41	LEU	3.3
2	B	1	GLU	3.3
2	D	139	THR	3.3
2	B	177	GLN	3.2
1	C	180	ASP	3.2
2	B	223	ASP	3.2
2	D	221	THR	3.1
2	D	223	ASP	3.0
1	C	163	ARG	3.0
2	B	217	ASN	3.0
2	D	129	GLU	2.9
1	C	20	SER	2.9
1	A	127	SER	2.9
1	A	163	ARG	2.8
2	D	181	ASN	2.8
1	A	51	TYR	2.8
1	C	178	LYS	2.7
2	B	62	GLU	2.6
1	A	180	ASP	2.6
1	A	92	TYR	2.6
2	D	127	PRO	2.5
1	C	21	LEU	2.4
1	A	29	GLY	2.4
2	D	217	ASN	2.4
1	C	92	TYR	2.4
1	C	129	ASP	2.4
1	C	19	ALA	2.4
2	D	47	TYR	2.4
1	C	134	LEU	2.4
2	D	140	LEU	2.3
1	C	9	GLY	2.3
2	B	47	TYR	2.3
2	D	57	LYS	2.3
1	A	191	ILE	2.3
2	B	222	GLN	2.2
1	A	20	SER	2.2
1	A	132	VAL	2.2
1	C	166	ASP	2.2

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Mol	Chain	Res	Type	RSRZ
2	D	33	SER	2.2
2	D	40	GLY	2.2
1	C	179	SER	2.2
1	A	194	GLU	2.2
2	D	115	LYS	2.2
1	C	51	TYR	2.1
1	C	161	ASP	2.1
1	C	194	GLU	2.1
2	D	222	GLN	2.1
1	C	167	PHE	2.1
2	D	97	SER	2.1
1	A	167	PHE	2.1
1	A	181	PHE	2.1
2	D	142	CYS	2.0
1	C	173	VAL	2.0
2	D	141	VAL	2.0
2	D	35	TYR	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.