



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 10:20 PM GMT

PDB ID : 1T5W  
Title : HLA-DR1 in complex with a synthetic peptide (AAYSDQATPLLLSPR)  
Authors : Zavala-Ruiz, Z.; Strug, I.; Anderson, M.W.; Gorski, J.; Stern, L.J.  
Deposited on : 2004-05-05  
Resolution : 2.40 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

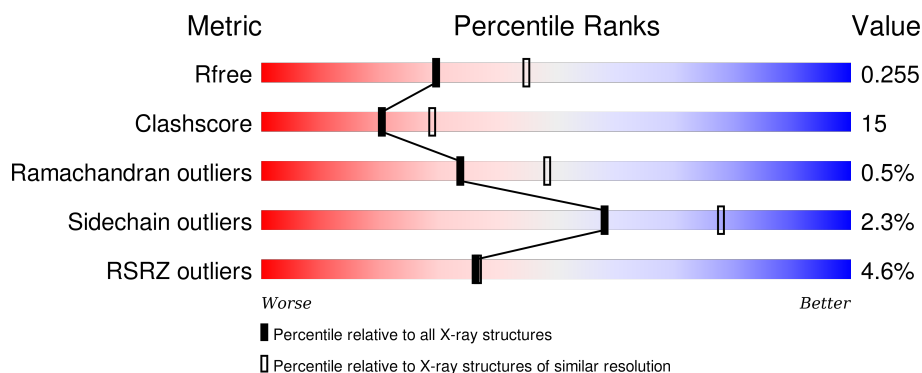
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.40 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	2919 (2.40-2.40)
Clashscore	102246	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)
RSRZ outliers	91569	2928 (2.40-2.40)

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	180	<div> <div>0%</div> <div> <div></div> <div>79%</div> <div>21%</div> <div>•</div> </div> </div>
1	D	180	<div> <div>6%</div> <div> <div></div> <div>76%</div> <div>23%</div> <div>••</div> </div> </div>
2	B	190	<div> <div>3%</div> <div> <div></div> <div>72%</div> <div>25%</div> <div>•</div> </div> </div>
2	E	190	<div> <div>8%</div> <div> <div></div> <div>69%</div> <div>27%</div> <div>•</div> </div> </div>
3	C	15	<div> <div>7%</div> <div> <div></div> <div>47%</div> <div>33%</div> <div>7%</div> <div>13%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
3	F	15	<div><div></div><div></div><div></div><div></div><div></div></div> <div><div>7%</div><div>60%</div><div>20%</div><div>7%</div><div>13%</div></div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6407 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 HLA class II histocompatibility antigen, DR alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	180	Total	C	N	O	S	9	0	0
			1483	960	241	277	5			
1	D	179	Total	C	N	O	S	0	0	0
			1474	954	239	276	5			

- Molecule 2 is a protein called HLA class II histocompatibility antigen, DRB1-1 beta chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	190	Total	C	N	O	S	22	0	0
			1557	979	279	293	6			
2	E	190	Total	C	N	O	S	0	0	0
			1557	979	279	293	6			

- Molecule 3 is a protein called 15-mer peptide fragment of Regulatory protein MIG1.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	13	Total	C	N	O	0	0	0
			94	60	14	20			
3	F	13	Total	C	N	O	0	0	0
			94	60	14	20			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	0	ALA	-	INSERTION	UNP P27705
C	1	ALA	-	INSERTION	UNP P27705
C	2	TYR	-	INSERTION	UNP P27705
C	3	SER	-	INSERTION	UNP P27705
C	4	ASP	-	INSERTION	UNP P27705
C	5	GLN	-	INSERTION	UNP P27705
C	6	ALA	-	INSERTION	UNP P27705
F	0	ALA	-	INSERTION	UNP P27705

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Chain	Residue	Modelled	Actual	Comment	Reference
F	1	ALA	-	INSERTION	UNP P27705
F	2	TYR	-	INSERTION	UNP P27705
F	3	SER	-	INSERTION	UNP P27705
F	4	ASP	-	INSERTION	UNP P27705
F	5	GLN	-	INSERTION	UNP P27705
F	6	ALA	-	INSERTION	UNP P27705

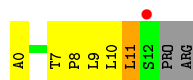
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	64	Total O 64 64	0	0
4	B	42	Total O 42 42	0	0
4	C	4	Total O 4 4	0	0
4	D	12	Total O 12 12	0	0
4	E	26	Total O 26 26	0	0

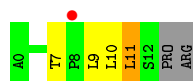




- Molecule 3: 15-mer peptide fragment of Regulatory protein MIG1



- Molecule 3: 15-mer peptide fragment of Regulatory protein MIG1



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	96.55Å 112.65Å 206.98Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	23.90 – 2.40 24.79 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.8 (23.90-2.40) 99.4 (24.79-2.30)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	0.07	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.07 (at 2.28Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, $R_{free}$	0.231 , 0.255 0.231 , 0.255	Depositor DCC
$R_{free}$ test set	3142 reflections (7.09%)	DCC
Wilson B-factor (Å <sup>2</sup> )	36.0	Xtriage
Anisotropy	0.689	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 44.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 50347 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	6407	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.56% 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	0.43	0/1528	0.65	0/2081
1	D	0.36	0/1519	0.62	0/2070
2	B	0.40	0/1597	0.66	2/2168 (0.1%)
2	E	0.35	0/1597	0.63	2/2168 (0.1%)
3	C	0.39	0/95	0.70	0/130
3	F	0.37	0/95	0.71	0/130
All	All	0.38	0/6431	0.64	4/8747 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	166	ARG	NE-CZ-NH2	-7.34	116.63	120.30
2	B	166	ARG	NE-CZ-NH1	-6.97	116.81	120.30
2	E	166	ARG	NE-CZ-NH1	6.58	123.59	120.30
2	B	166	ARG	NE-CZ-NH2	6.32	123.46	120.30

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	1483	0	1420	34	1
1	D	1474	0	1407	35	0
2	B	1557	0	1488	61	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	E	1557	0	1488	65	0
3	C	94	0	95	14	0
3	F	94	0	95	10	0
4	A	64	0	0	2	0
4	B	42	0	0	2	0
4	C	4	0	0	0	0
4	D	12	0	0	0	0
4	E	26	0	0	1	0
All	All	6407	0	5993	178	1

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 15.

All (178) 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:E:150:ASN:HD22	2:E:154:THR:HG22	1.08	1.13
2:B:150:ASN:HD22	2:B:154:THR:HG22	1.09	1.08
1:D:96:PRO:HG3	2:E:100:THR:HG21	1.37	1.01
2:E:65:LYS:H	2:E:65:LYS:HE3	1.35	0.91
2:B:65:LYS:HE3	2:B:65:LYS:H	1.32	0.91
1:D:147:LYS:HE3	1:D:149:HIS:HE1	1.38	0.86
2:B:100:THR:HG23	2:B:118:SER:HB3	1.58	0.84
2:E:150:ASN:ND2	2:E:154:THR:HG22	1.91	0.84
2:E:56:PRO:HB3	3:F:11:LEU:HD12	1.59	0.83
2:B:150:ASN:HD22	2:B:154:THR:CG2	1.90	0.82
1:A:147:LYS:HE3	1:A:149:HIS:HE1	1.45	0.81
2:B:150:ASN:ND2	2:B:154:THR:HG22	1.93	0.80
2:E:100:THR:HG23	2:E:118:SER:HB3	1.62	0.79
2:B:56:PRO:HB3	3:C:11:LEU:HD12	1.64	0.77
1:A:16:PRO:HD2	2:B:6:ARG:HD3	1.68	0.76
2:E:150:ASN:HD22	2:E:154:THR:CG2	1.96	0.75
1:D:16:PRO:HD2	2:E:6:ARG:HD3	1.70	0.73
1:A:96:PRO:HG3	2:B:100:THR:HG21	1.69	0.73
1:D:67:LYS:O	1:D:71:GLU:HG2	1.89	0.72
2:B:105:LYS:HZ3	2:B:107:GLN:HE21	1.36	0.71
2:E:94:ARG:HG3	2:E:94:ARG:HH11	1.56	0.71
2:B:8:LEU:O	2:B:32:TYR:O	2.08	0.70
2:E:93:ARG:HG2	2:E:123:TYR:CD1	2.26	0.70
2:E:97:PRO:HB3	2:E:122:PHE:HB3	1.74	0.70
2:B:93:ARG:HG2	2:B:123:TYR:CD1	2.27	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:105:LYS:HZ3	2:E:107:GLN:HE21	1.39	0.69
2:E:56:PRO:CB	3:F:11:LEU:HD12	2.23	0.69
2:B:97:PRO:HB3	2:B:122:PHE:HB3	1.76	0.68
2:E:8:LEU:O	2:E:32:TYR:O	2.11	0.68
2:B:94:ARG:HH11	2:B:94:ARG:HG3	1.57	0.68
1:D:81:PRO:HB3	2:E:5:PRO:HB2	1.76	0.67
1:A:67:LYS:O	1:A:71:GLU:HG2	1.95	0.66
2:B:65:LYS:CE	2:B:65:LYS:H	2.07	0.65
2:B:56:PRO:CB	3:C:11:LEU:HD12	2.26	0.64
2:E:152:ASP:OD1	2:E:154:THR:HB	1.99	0.63
2:B:2:ASP:OD1	2:B:4:ARG:HD3	1.98	0.63
2:E:170:VAL:HG22	2:E:189:ARG:HG2	1.81	0.62
2:E:113:ASN:HD22	2:E:114:LEU:H	1.47	0.62
1:D:147:LYS:HE3	1:D:149:HIS:CE1	2.27	0.62
1:D:100:ARG:HH11	1:D:100:ARG:CG	2.13	0.62
2:B:105:LYS:NZ	2:B:107:GLN:HE21	1.99	0.60
2:E:94:ARG:HG3	2:E:94:ARG:NH1	2.17	0.60
1:A:100:ARG:HH11	1:A:100:ARG:CG	2.15	0.60
2:E:105:LYS:NZ	2:E:107:GLN:HE21	1.99	0.59
1:D:99:LEU:HD21	1:D:180:PHE:CE2	2.37	0.58
2:B:94:ARG:HG3	2:B:94:ARG:NH1	2.18	0.58
2:B:55:ARG:HB2	2:B:56:PRO:HD3	1.86	0.58
1:A:147:LYS:HE3	1:A:149:HIS:CE1	2.34	0.57
2:B:113:ASN:HD22	2:B:114:LEU:H	1.50	0.57
2:E:55:ARG:HB2	2:E:56:PRO:HD3	1.86	0.57
2:E:145:THR:CG2	2:E:158:LEU:H	2.17	0.57
2:E:105:LYS:NZ	2:E:105:LYS:HB2	2.20	0.56
1:D:118:ASN:HB2	1:D:166:GLU:HB2	1.87	0.56
1:A:77:SER:O	1:A:78:ASN:HB3	2.05	0.56
2:E:25:ARG:HD2	2:E:43:ASP:OD2	2.05	0.56
2:B:145:THR:CG2	2:B:158:LEU:H	2.18	0.56
2:B:61:TRP:CZ2	3:C:10:LEU:HD12	2.41	0.56
2:B:105:LYS:NZ	2:B:105:LYS:HB2	2.22	0.55
2:E:65:LYS:H	2:E:65:LYS:CE	2.12	0.55
2:B:25:ARG:HD2	2:B:43:ASP:OD2	2.07	0.55
1:A:110:ASP:OD1	1:A:146:ARG:HG2	2.06	0.55
2:B:152:ASP:OD1	2:B:154:THR:HB	2.06	0.55
1:A:89:VAL:HG12	1:A:176:LYS:HG3	1.89	0.54
2:B:61:TRP:HZ2	3:C:10:LEU:HD12	1.73	0.54
1:D:110:ASP:OD1	1:D:146:ARG:HG2	2.07	0.54
2:E:2:ASP:OD1	2:E:6:ARG:NH2	2.41	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:60:TYR:CD2	3:F:11:LEU:HD23	2.43	0.54
2:E:23:ARG:HG2	2:E:23:ARG:HH11	1.72	0.54
1:D:77:SER:O	1:D:78:ASN:HB3	2.08	0.53
2:E:35:GLU:CG	2:E:51:THR:HG21	2.38	0.53
2:E:100:THR:HG22	4:E:193:HOH:O	2.09	0.52
1:A:97:VAL:HG12	4:A:218:HOH:O	2.09	0.52
2:B:35:GLU:CG	2:B:51:THR:HG21	2.40	0.52
1:A:100:ARG:HH11	1:A:100:ARG:HG3	1.75	0.52
2:E:61:TRP:CZ2	3:F:10:LEU:HD12	2.45	0.52
2:B:23:ARG:HH11	2:B:23:ARG:HG2	1.75	0.52
1:A:69:ASN:HB3	3:C:10:LEU:HD22	1.93	0.51
2:E:81:HIS:O	2:E:85:VAL:HG23	2.10	0.51
1:A:76:ARG:NH2	2:B:57:ASP:OD2	2.39	0.51
1:D:99:LEU:HD21	1:D:180:PHE:CD2	2.46	0.51
2:B:35:GLU:CD	2:B:51:THR:HG21	2.31	0.51
1:D:76:ARG:NH2	2:E:57:ASP:OD2	2.38	0.50
1:A:65:VAL:HG11	3:C:7:THR:OG1	2.12	0.50
1:A:100:ARG:NH1	1:A:100:ARG:HG3	2.26	0.50
2:B:51:THR:HA	2:E:51:THR:HA	1.92	0.50
1:D:100:ARG:NH1	1:D:100:ARG:CG	2.74	0.50
2:E:9:TRP:CH2	2:E:30:CYS:HB3	2.47	0.50
2:E:105:LYS:NZ	2:E:107:GLN:NE2	2.60	0.50
2:E:35:GLU:CD	2:E:51:THR:HG21	2.31	0.49
2:E:134:ASN:ND2	2:E:170:VAL:H	2.11	0.49
2:B:85:VAL:HG22	3:C:0:ALA:HB1	1.93	0.49
2:B:105:LYS:NZ	2:B:107:GLN:NE2	2.59	0.49
1:D:89:VAL:HG12	1:D:176:LYS:HG3	1.94	0.48
2:E:60:TYR:CE1	2:E:64:GLN:NE2	2.81	0.48
2:E:55:ARG:CB	2:E:56:PRO:HD3	2.43	0.48
1:D:12:PHE:C	1:D:12:PHE:CD1	2.87	0.48
2:E:113:ASN:ND2	2:E:114:LEU:H	2.10	0.47
2:E:145:THR:HG22	2:E:158:LEU:H	1.78	0.47
1:D:100:ARG:HG3	1:D:100:ARG:NH1	2.29	0.47
2:B:104:SER:O	2:B:105:LYS:HB2	2.15	0.47
4:A:187:HOH:O	2:B:120:SER:HB2	2.15	0.47
2:E:31:ILE:HD12	2:E:31:ILE:N	2.28	0.47
2:B:162:GLU:OE2	1:D:162:ASP:OD2	2.33	0.47
2:B:31:ILE:N	2:B:31:ILE:HD12	2.30	0.47
2:B:55:ARG:CB	2:B:56:PRO:HD3	2.45	0.46
2:B:113:ASN:ND2	2:B:114:LEU:H	2.14	0.46
2:B:145:THR:HG21	2:B:158:LEU:H	1.81	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:3:GLU:HB2	2:E:18:PHE:CE2	2.50	0.46
1:A:118:ASN:HB2	1:A:166:GLU:HB2	1.96	0.46
2:E:145:THR:HG23	2:E:146:GLY:O	2.16	0.46
2:B:134:ASN:ND2	2:B:170:VAL:H	2.14	0.46
2:B:2:ASP:OD1	2:B:6:ARG:NH2	2.49	0.46
2:B:100:THR:HG22	4:B:192:HOH:O	2.16	0.45
2:E:61:TRP:HZ2	3:F:10:LEU:HD12	1.81	0.45
1:D:69:ASN:HB3	3:F:10:LEU:HD22	1.98	0.45
1:D:96:PRO:HG3	2:E:100:THR:CG2	2.27	0.45
2:E:65:LYS:HG2	2:E:66:ASP:N	2.32	0.45
2:E:60:TYR:HB2	3:F:11:LEU:HD21	1.99	0.45
1:A:73:MET:HE2	3:C:10:LEU:HD21	1.99	0.45
2:B:117:CYS:HB2	2:B:131:TRP:CZ2	2.52	0.45
2:B:60:TYR:CD2	3:C:11:LEU:HD23	2.53	0.44
1:A:82:ILE:HG13	2:B:33:ASN:HB3	2.00	0.44
1:D:121:TRP:O	1:D:127:PRO:HA	2.17	0.44
1:D:141:GLU:HA	1:D:141:GLU:OE1	2.17	0.44
1:A:3:GLU:OE2	1:A:6:VAL:CG2	2.65	0.44
2:B:145:THR:HG23	2:B:146:GLY:O	2.17	0.44
2:E:18:PHE:HB2	2:E:23:ARG:HB3	2.00	0.44
1:A:65:VAL:CG1	3:C:7:THR:OG1	2.67	0.43
2:B:11:LEU:CD2	3:C:7:THR:HG22	2.48	0.43
2:B:65:LYS:HG2	2:B:66:ASP:N	2.34	0.43
3:C:11:LEU:HA	3:C:11:LEU:HD23	1.62	0.43
1:D:73:MET:HE2	3:F:10:LEU:HD21	2.01	0.43
1:D:94:ASN:HB3	1:D:106:ILE:HD11	2.01	0.43
2:B:9:TRP:CH2	2:B:30:CYS:HB3	2.54	0.43
4:B:228:HOH:O	3:C:8:PRO:HD2	2.18	0.43
1:D:82:ILE:HG13	2:E:33:ASN:HB3	2.00	0.43
1:D:50:ARG:HB2	1:D:50:ARG:NH2	2.34	0.43
1:A:50:ARG:CB	1:A:50:ARG:HH21	2.31	0.43
3:F:11:LEU:HD23	3:F:11:LEU:HA	1.51	0.43
1:A:50:ARG:NH2	1:A:50:ARG:HB2	2.34	0.43
1:A:94:ASN:HB3	1:A:106:ILE:HD11	2.00	0.43
2:B:90:THR:OG1	2:B:91:VAL:N	2.51	0.43
2:E:30:CYS:C	2:E:31:ILE:HD12	2.39	0.42
1:D:26:PHE:HB2	1:D:31:ILE:HD11	2.01	0.42
1:D:11:GLU:HG3	2:E:11:LEU:HB3	2.00	0.42
1:D:129:THR:O	1:D:129:THR:HG22	2.19	0.42
2:E:104:SER:O	2:E:105:LYS:HB2	2.19	0.42
1:D:50:ARG:HH21	1:D:50:ARG:CB	2.31	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:13:PHE:CE2	2:E:28:GLU:HG3	2.54	0.42
1:A:12:PHE:C	1:A:12:PHE:CD1	2.92	0.42
1:D:65:VAL:HG11	3:F:7:THR:OG1	2.19	0.42
1:A:121:TRP:O	1:A:127:PRO:HA	2.19	0.42
2:B:55:ARG:NE	2:E:55:ARG:HH21	2.17	0.42
2:B:2:ASP:CG	2:B:6:ARG:HH22	2.22	0.42
2:B:97:PRO:CB	2:B:122:PHE:HB3	2.48	0.42
1:A:69:ASN:CB	3:C:10:LEU:HD22	2.50	0.42
1:A:129:THR:O	1:A:129:THR:HG22	2.20	0.42
2:E:117:CYS:HB2	2:E:131:TRP:CZ2	2.54	0.42
2:E:113:ASN:HD22	2:E:114:LEU:N	2.16	0.41
1:D:87:PRO:HB3	1:D:112:PHE:HB3	2.02	0.41
2:B:18:PHE:HB2	2:B:23:ARG:HB3	2.02	0.41
2:E:68:LEU:O	2:E:72:ARG:HG3	2.19	0.41
1:A:11:GLU:HG3	2:B:11:LEU:HB3	2.03	0.41
1:A:76:ARG:HH22	2:B:57:ASP:CG	2.23	0.41
1:A:47:GLU:O	1:A:51:PHE:HD1	2.04	0.41
1:A:81:PRO:HB3	2:B:5:PRO:HB2	2.03	0.41
2:E:9:TRP:CZ2	2:E:30:CYS:HB3	2.56	0.41
2:B:145:THR:HG22	2:B:158:LEU:H	1.83	0.41
2:E:74:ALA:O	2:E:78:TYR:HB3	2.21	0.41
1:A:141:GLU:OE1	1:A:141:GLU:HA	2.21	0.41
1:A:162:ASP:OD2	2:E:162:GLU:OE2	2.39	0.41
1:A:143:HIS:HD2	2:B:12:LYS:NZ	2.19	0.40
1:D:17:ASP:O	1:D:18:GLN:HB2	2.21	0.40
2:E:2:ASP:CG	2:E:6:ARG:HH22	2.24	0.40
2:E:105:LYS:HZ3	2:E:105:LYS:HB2	1.87	0.40
2:B:113:ASN:HD22	2:B:114:LEU:N	2.18	0.40
2:E:145:THR:HG21	2:E:158:LEU:H	1.84	0.40
1:D:47:GLU:O	1:D:51:PHE:HD1	2.05	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:57:GLN:NE2	1:A:57:GLN:NE2[4_555]	1.83	0.37

## 5.3 Torsion angles

### 5.3.1 Protein backbone

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	178/180 (99%)	173 (97%)	5 (3%)	0	100	100
1	D	177/180 (98%)	172 (97%)	5 (3%)	0	100	100
2	B	188/190 (99%)	182 (97%)	4 (2%)	2 (1%)	17	25
2	E	188/190 (99%)	180 (96%)	6 (3%)	2 (1%)	17	25
3	C	11/15 (73%)	10 (91%)	1 (9%)	0	100	100
3	F	11/15 (73%)	9 (82%)	2 (18%)	0	100	100
All	All	753/770 (98%)	726 (96%)	23 (3%)	4 (0%)	34	48

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	32	TYR
2	E	32	TYR
2	B	105	LYS
2	E	105	LYS

### 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	165/165 (100%)	164 (99%)	1 (1%)	90	96
1	D	164/165 (99%)	163 (99%)	1 (1%)	90	96
2	B	171/171 (100%)	166 (97%)	5 (3%)	50	71

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	E	171/171 (100%)	166 (97%)	5 (3%)	50	71
3	C	10/12 (83%)	8 (80%)	2 (20%)	1	1
3	F	10/12 (83%)	8 (80%)	2 (20%)	1	1
All	All	691/696 (99%)	675 (98%)	16 (2%)	58	78

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	100	ARG
2	B	29	ARG
2	B	65	LYS
2	B	100	THR
2	B	113	ASN
2	B	154	THR
3	C	9	LEU
3	C	11	LEU
1	D	100	ARG
2	E	29	ARG
2	E	65	LYS
2	E	100	THR
2	E	113	ASN
2	E	154	THR
3	F	9	LEU
3	F	11	LEU

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

Mol	Chain	Res	Type
1	A	18	GLN
1	A	143	HIS
1	A	149	HIS
2	B	64	GLN
2	B	107	GLN
2	B	113	ASN
2	B	134	ASN
2	B	150	ASN
2	B	156	GLN
3	C	5	GLN
1	D	18	GLN
1	D	143	HIS

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Mol	Chain	Res	Type
1	D	149	HIS
2	E	64	GLN
2	E	107	GLN
2	E	113	ASN
2	E	134	ASN
2	E	150	ASN
2	E	156	GLN
3	F	5	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 [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	179/180 (99%)	-0.19	2 (1%) 82 82	18, 30, 69, 93	0
1	D	179/180 (99%)	0.24	10 (5%) 28 28	27, 46, 96, 138	0
2	B	190/190 (100%)	0.12	5 (2%) 59 58	17, 34, 75, 169	5 (2%)
2	E	190/190 (100%)	0.45	16 (8%) 14 13	24, 47, 96, 149	0
3	C	13/15 (86%)	0.54	1 (7%) 16 16	26, 42, 67, 104	0
3	F	13/15 (86%)	0.74	1 (7%) 16 16	47, 61, 70, 71	0
All	All	764/770 (99%)	0.18	35 (4%) 36 37	17, 41, 87, 169	5 (0%)

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	1	GLY	19.0
2	E	1	GLY	9.7
1	D	100	ARG	4.7
2	E	106	THR	4.6
2	E	107	GLN	4.3
2	E	18	PHE	3.6
1	D	99	LEU	3.5
1	D	37	ALA	3.5
1	D	181	ASP	3.5
2	E	3	THR	3.4
1	D	57	GLN	3.2
2	E	105	LYS	3.2
1	D	101	GLU	3.1
2	B	181	THR	3.0
2	E	168	GLY	3.0
2	E	167	SER	2.9
1	D	51	PHE	2.8
2	E	189	ARG	2.8
2	E	23	ARG	2.7

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Mol	Chain	Res	Type	RSRZ
2	E	19	ASN	2.7
2	E	65	LYS	2.6
3	C	12	SER	2.6
2	B	136	GLN	2.6
1	D	38	LYS	2.5
2	E	5	PRO	2.5
2	E	66	ASP	2.5
2	E	2	ASP	2.4
1	D	50	ARG	2.3
1	D	39	LYS	2.2
1	A	129	THR	2.2
2	B	182	SER	2.1
1	A	130	THR	2.1
3	F	8	PRO	2.1
2	B	2	ASP	2.1
2	E	11	LEU	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.