

# Full wwPDB X-ray Structure Validation Report (i)

#### Jul 26, 2023 – 04:21 AM EDT

PDB ID	:	1BD2
Title	:	COMPLEX BETWEEN HUMAN T-CELL RECEPTOR B7, VIRAL PEP-
		TIDE (TAX) AND MHC CLASS I MOLECULE HLA-A 0201
Authors	:	Ding, YH.; Smith, K.J.; Garboczi, D.N.; Utz, U.; Biddison, W.E.; Wiley,
		D.C.
Deposited on	:	1998-05-12
Resolution	:	2.50  Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at *validation@mail.wwpdb.org* A user guide is available at https://www.wwpdb.org/validation/2017/XrayValidationReportHelp with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references (1)) were used in the production of this report:

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	NOT EXECUTED
$\mathrm{EDS}$	:	NOT EXECUTED
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.34

### 1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:  $X\text{-}RAY \, DIFFRACTION$ 

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.



Matria	Whole archive	Similar resolution
wietric	$(\# { m Entries})$	$(\# { m Entries},  { m resolution}  { m range}({ m \AA}))$
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=5%

Note EDS was not executed.

Mol	Chain	Length	Quality of chai			
1	А	275	71%		28%	•
2	В	100	75%		22%	•••
3	С	9	33%	67%		_
4	D	204	48%	38%	5% •	7%
5	Е	244	62%	3	34%	•••



### 2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 6378 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-A 0201.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	А	275	Total	С	N	0	S	0	0	0
			2186	1369	392	416	9	_	-	-

• Molecule 2 is a protein called BETA-2 MICROGLOBULIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	В	99	Total 805	C 512	N 138	O 152	${ m S} { m 3}$	0	0	0

• Molecule 3 is a protein called TAX PEPTIDE.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf	Trace
3	С	9	Total         C         N         O           77         56         9         12	0	0	0

• Molecule 4 is a protein called T CELL RECEPTOR ALPHA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	190	Total 1421	C 899	N 230	O 285	${ m S} 7$	0	0	0

There are 13 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	56	ALA	GLU	conflict	GB 338766
D	?	-	LYS	deletion	GB 338766
D	93	MET	GLY	conflict	GB 338766
D	94	GLU	ALA	conflict	GB 338766
D	?	-	THR	deletion	GB 338766
D	102	GLN	SER	conflict	GB 338766
D	105	VAL	THR	conflict	GB 338766
D	108	GLN	THR	conflict	GB 338766



Chain	Residue	Modelled	Actual	Comment	Reference
D	113	THR	GLN	conflict	GB 338766
D	114	ILE	VAL	conflict	GB 338766
D	115	ASN	THR	conflict	GB 338766
D	116	PRO	LEU	conflict	GB 338766
D	117	ASN	ASP	conflict	GB 338766

• Molecule 5 is a protein called T CELL RECEPTOR BETA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	Е	242	Total 1850	C 1172	N 318	O 353	${ m S} 7$	0	0	0

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
Е	96	TYR	PHE	conflict	GB 3002925
Е	98	GLY	ARG	conflict	GB 3002925
E	99	GLY	GLN	conflict	GB 3002925
E	100	GLY	PRO	conflict	GB 3002925
Е	101	PHE	SER	conflict	GB 3002925
Е	?	-	ASN	deletion	GB 3002925
Е	107	TYR	PHE	conflict	GB 3002925
Е	117	THR	LEU	conflict	GB 3002925
Ē	192	ALA	CYS	conflict	GB 3002925
Е	206	ASP	ASN	conflict	GB 3002925

• Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	А	9	Total O 9 9	0	0
6	В	5	Total O 5 5	0	0
6	D	13	Total O 13 13	0	0
6	Е	12	$\begin{array}{ccc} \text{Total} & \text{O} \\ 12 & 12 \end{array}$	0	0



#### Residue-property plots (i) 3

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. 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. 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.

Chain A: 71% 28% • Molecule 2: BETA-2 MICROGLOBULIN Chain B: 75% 22% • Molecule 3: TAX PEPTIDE Chain C: 33% 67% • Molecule 4: T CELL RECEPTOR ALPHA Chain D: 48% 38% 5% • 7%

Note EDS was not executed.

- Molecule 1: HLA-A 0201



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 $\bullet$  Molecule 5: T CELL RECEPTOR BETA





### 4 Data and refinement statistics (i)

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

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants	63.80Å 73.30Å 217.70Å	Depositor
a, b, c, $\alpha$ , $\beta$ , $\gamma$	$90.00^{\circ}$ $90.00^{\circ}$ $90.00^{\circ}$	Depositor
Resolution (Å)	8.00 - 2.50	Depositor
% Data completeness	90.6 (8.00-2.50)	Depositor
(in resolution range)	50.0 (0.00 2.00)	Depositor
$R_{merge}$	0.11	Depositor
R <sub>sym</sub>	(Not available)	Depositor
Refinement program	X-PLOR 3.851	Depositor
$R, R_{free}$	0.238 , $0.312$	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	6378	wwPDB-VP
Average B, all atoms $(Å^2)$	41.0	wwPDB-VP



# 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).

Mal	Mol Chain		Bond lengths		nd angles
	Unam	RMSZ	# Z  > 5	RMSZ	# Z  > 5
1	А	0.47	0/2250	0.69	2/3058~(0.1%)
2	В	0.45	0/828	0.70	0/1121
3	С	0.56	0/80	0.88	0/108
4	D	0.55	0/1452	0.86	4/1980~(0.2%)
5	Е	0.50	0/1905	0.75	0/2604
All	All	0.50	0/6515	0.75	6/8871~(0.1%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
4	D	196	SER	N-CA-C	5.61	126.15	111.00
4	D	199	PRO	N-CA-C	5.59	126.64	112.10
4	D	187	PHE	N-CA-C	-5.29	96.70	111.00
4	D	199	PRO	N-CA-CB	5.16	109.49	103.30
1	А	28	VAL	N-CA-C	-5.14	97.13	111.00
1	А	65	ARG	NE-CZ-NH1	-5.01	117.79	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	А	2186	0	2005	60	0
2	В	805	0	756	17	0



Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	С	77	0	79	12	0
4	D	1421	0	1274	80	0
5	Е	1850	0	1699	58	0
6	А	9	0	0	1	0
6	В	5	0	0	0	0
6	D	13	0	0	4	0
6	Е	12	0	0	0	0
All	All	6378	0	5813	208	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 17.

All (208) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom 1	Atom 2	Interatomic	Clash
Atom-1	Atom-2	distance (Å)	overlap (Å)
4:D:17:ILE:HG13	4:D:78:VAL:HG22	1.40	1.01
1:A:82:ARG:NH1	1:A:89:GLU:HG3	1.78	0.97
1:A:131:ARG:HE	1:A:157:ARG:HH12	1.18	0.88
4:D:151:SER:OG	4:D:195:ASN:HB2	1.73	0.87
1:A:131:ARG:NE	1:A:157:ARG:HH12	1.78	0.81
4:D:151:SER:CB	4:D:195:ASN:HB2	2.13	0.79
1:A:82:ARG:HH11	1:A:89:GLU:HG3	1.47	0.79
2:B:73:THR:CG2	2:B:75:LYS:HG2	2.14	0.77
1:A:131:ARG:HE	1:A:157:ARG:NH1	1.84	0.76
5:E:21:LEU:HD12	5:E:77:LEU:HD23	1.68	0.74
5:E:117:THR:HG23	5:E:154:TYR:OH	1.88	0.74
4:D:151:SER:HB2	4:D:195:ASN:HD22	1.54	0.72
2:B:22:PHE:CE2	2:B:69:GLU:HG2	2.26	0.70
4:D:15:GLY:HA2	6:D:215:HOH:O	1.91	0.70
4:D:67:ASN:HD22	4:D:70:ALA:HB3	1.56	0.70
2:B:73:THR:HG22	2:B:75:LYS:H	1.58	0.69
4:D:149:ASN:O	4:D:150:VAL:HG12	1.93	0.68
1:A:63:GLU:OE1	3:C:2:LEU:HD12	1.94	0.68
4:D:166:LEU:HB3	5:E:174:SER:HB2	1.76	0.67
1:A:45:MET:CE	3:C:2:LEU:HD11	2.26	0.66
4:D:58:GLY:O	4:D:61:ARG:HB2	1.95	0.66
4:D:194:ASN:O	4:D:195:ASN:CG	2.34	0.65
5:E:125:PRO:HD3	5:E:233:PRO:HB3	1.79	0.64
4:D:5:GLN:NE2	4:D:109:GLY:H	1.96	0.64
1:A:82:ARG:HH12	1:A:89:GLU:HG3	1.62	0.63
4:D:158:VAL:HG13	4:D:182:SER:HB2	1.80	0.63



		Interatomic	Clash
Atom-1	Atom-2	distance (Å)	overlap (Å)
2:B:73:THR:HG22	2:B:75:LYS:HG2	1.79	0.62
2:B:25:CYS:HB2	2:B:39:LEU:HD21	1.80	0.62
1:A:263:HIS:HD2	1:A:265:GLY:H	1.49	0.60
2:B:25:CYS:O	2:B:65:LEU:HD12	2.01	0.60
1:A:263:HIS:CD2	1:A:265:GLY:H	2.19	0.60
5:E:37:GLN:HB2	5:E:43:LEU:HD23	1.81	0.60
4:D:5:GLN:HE21	4:D:107:GLY:HA3	1.67	0.60
5:E:89:VAL:HG22	5:E:113:ARG:HG2	1.83	0.59
5:E:10:PHE:HB3	5:E:157:HIS:CD2	2.38	0.59
4:D:51:SER:O	4:D:68:LYS:HD2	2.03	0.59
5:E:182:GLU:HG2	5:E:190:ARG:O	2.03	0.59
1:A:6:ARG:HD2	1:A:100:GLY:HA3	1.84	0.58
4:D:152:GLN:HG2	4:D:153:SER:H	1.67	0.58
5:E:12:VAL:HG22	5:E:155:PRO:HG3	1.85	0.58
4:D:79:PRO:HB2	6:D:215:HOH:O	2.04	0.57
5:E:84:PRO:HA	5:E:116:VAL:HG13	1.84	0.57
4:D:173:PHE:HE2	4:D:175:SER:HB3	1.69	0.57
1:A:65:ARG:NH1	6:A:276:HOH:O	2.38	0.57
4:D:108:GLN:OE1	4:D:108:GLN:N	2.37	0.57
4:D:194:ASN:O	4:D:195:ASN:ND2	2.38	0.57
5:E:13:LEU:HD11	5:E:19:MET:HB2	1.86	0.57
4:D:153:SER:HB3	4:D:158:VAL:O	2.06	0.56
5:E:133:PRO:HG3	5:E:146:LEU:HD12	1.86	0.56
4:D:144:PHE:CE1	4:D:176:ASN:HB3	2.41	0.56
5:E:12:VAL:CG2	5:E:155:PRO:HG3	2.36	0.56
1:A:204:TRP:HE3	1:A:206:LEU:HD21	1.71	0.56
1:A:187:THR:HG22	1:A:204:TRP:O	2.05	0.56
5:E:8:PRO:O	5:E:112:THR:HB	2.06	0.55
5:E:205:GLN:HA	5:E:245:ARG:O	2.06	0.55
4:D:142:THR:HA	4:D:176:ASN:O	2.07	0.55
5:E:131:PHE:HB2	5:E:147:VAL:HG13	1.87	0.55
4:D:91:ALA:HB1	4:D:104:LEU:CD1	2.37	0.55
5:E:65:TYR:HB3	5:E:77:LEU:HD11	1.89	0.55
1:A:208:PHE:CE1	1:A:241:PHE:HB2	2.42	0.55
4:D:115:ASN:HB3	4:D:146:SER:HB3	1.89	0.54
4:D:145:ASP:OD2	4:D:147:GLN:HB2	2.07	0.54
1:A:80:THR:HG21	3:C:9:VAL:OXT	2.07	0.54
1:A:187:THR:HB	1:A:272:LEU:HD11	1.89	0.54
1:A:234:ARG:HG2	1:A:242:GLN:HB2	1.90	0.54
1:A:51:TRP:CZ3	1:A:171:TYR:HB3	2.42	0.54
5:E:207:PRO:HA	5:E:244:GLY:O	2.08	0.54



		Interatomic	Clash
Atom-1	Atom-2	distance (Å)	overlap (Å)
5:E:148:CYS:HB2	5:E:162:TRP:CZ2	2.43	0.54
1:A:81:LEU:HD23	1:A:84:TYR:HD2	1.73	0.53
4:D:187:PHE:CD1	4:D:187:PHE:C	2.82	0.53
5:E:83:ALA:O	5:E:116:VAL:HG11	2.08	0.53
4:D:1:GLN:HG2	4:D:26:ASN:ND2	2.23	0.53
4:D:5:GLN:HE21	4:D:107:GLY:CA	2.21	0.53
1:A:187:THR:HG21	1:A:261:VAL:HG21	1.90	0.52
4:D:128:LEU:HD22	5:E:147:VAL:HG12	1.91	0.52
1:A:117:ALA:HB2	2:B:60:TRP:CE2	2.43	0.52
4:D:151:SER:CB	4:D:195:ASN:HD22	2.22	0.52
4:D:67:ASN:ND2	4:D:70:ALA:HB3	2.24	0.52
4:D:56:ALA:HB2	4:D:63:THR:HG23	1.90	0.52
4:D:61:ARG:CG	4:D:79:PRO:HD2	2.41	0.51
4:D:151:SER:HB2	4:D:195:ASN:ND2	2.22	0.51
4:D:173:PHE:CE2	4:D:175:SER:HB3	2.45	0.51
4:D:118:ILE:HB	4:D:121:PRO:HG3	1.91	0.51
4:D:150:VAL:O	4:D:150:VAL:HG13	2.10	0.51
1:A:194:VAL:HG12	1:A:194:VAL:O	2.10	0.51
5:E:6:GLN:HE22	5:E:92:CYS:H	1.58	0.51
5:E:21:LEU:HB2	5:E:77:LEU:HB3	1.93	0.51
1:A:131:ARG:NE	1:A:157:ARG:NH1	2.51	0.51
4:D:158:VAL:HA	4:D:182:SER:HB2	1.93	0.51
4:D:197:ILE:HD13	4:D:197:ILE:N	2.27	0.50
4:D:145:ASP:C	4:D:147:GLN:H	2.15	0.50
2:B:73:THR:HG22	2:B:75:LYS:N	2.26	0.50
1:A:211:ALA:HB2	1:A:241:PHE:CE2	2.46	0.50
5:E:4:VAL:HG13	5:E:109:GLY:HA2	1.92	0.50
5:E:123:VAL:O	5:E:230:ARG:NH2	2.45	0.50
1:A:72:GLN:NE2	1:A:75:ARG:HD3	2.27	0.50
5:E:234:VAL:O	5:E:236:GLN:HG2	2.12	0.50
5:E:8:PRO:HG2	5:E:11:GLN:HE21	1.77	0.49
3:C:5:TYR:N	3:C:5:TYR:CD1	2.80	0.49
5:E:124:PHE:CE1	5:E:230:ARG:NH1	2.81	0.49
5:E:141:THR:O	5:E:142:GLN:HB2	2.13	0.49
1:A:66:LYS:NZ	3:C:1:LEU:HG	2.28	0.49
4:D:123:PRO:O	4:D:200:GLU:HB3	2.13	0.49
4:D:128:LEU:HD22	5:E:147:VAL:CG1	2.42	0.49
1:A:159:TYR:HA	1:A:163:THR:HG23	1.94	0.48
2:B:29:GLY:HA2	2:B:61:SER:HB2	1.95	0.48
5:E:31:TYR:CE1	5:E:106:GLN:NE2	2.80	0.48
1:A:4:SER:HB3	1:A:102:ASP:OD1	2.14	0.48



	i agem	Interatomic	Clash
Atom-1	Atom-2	distance (Å)	overlap (Å)
1:A:73:THR:HA	3:C:8:TYR:CE1	2.48	0.48
1:A:49:ALA:O	1:A:52:ILE:HG22	2.13	0.48
1:A:73:THR:HA	3:C:8:TYR:HE1	1.79	0.48
1:A:204:TRP:CE3	1:A:206:LEU:HD21	2.48	0.48
4:D:34:TRP:CH2	4:D:90:CYS:HB2	2.48	0.48
4:D:47:ILE:HD13	4:D:64:VAL:HG23	1.95	0.48
4:D:155:ASP:HB3	4:D:158:VAL:HG23	1.96	0.48
5:E:37:GLN:HB2	5:E:43:LEU:CD2	2.43	0.47
1:A:56:GLY:O	1:A:59:TYR:HB3	2.14	0.47
1:A:13:SER:HA	1:A:20:PRO:HB3	1.96	0.47
1:A:213:ILE:CG2	1:A:243:LYS:HD2	2.44	0.47
3:C:5:TYR:CD2	4:D:96:ALA:HB2	2.50	0.47
5:E:131:PHE:N	5:E:131:PHE:CD1	2.83	0.47
5:E:31:TYR:HE1	5:E:106:GLN:NE2	2.13	0.47
2:B:6:LYS:HE2	2:B:29:GLY:HA3	1.98	0.46
4:D:87:VAL:HG22	4:D:111:ARG:HG2	1.97	0.46
5:E:6:GLN:HE21	5:E:109:GLY:HA3	1.80	0.46
5:E:31:TYR:HE1	5:E:106:GLN:HE22	1.63	0.46
5:E:241:GLU:OE2	5:E:243:TRP:CH2	2.69	0.46
1:A:211:ALA:HB2	1:A:241:PHE:CZ	2.51	0.46
1:A:5:MET:HB2	1:A:168:LEU:HD13	1.97	0.46
1:A:66:LYS:O	1:A:69:ALA:HB3	2.16	0.46
1:A:135:ALA:HB1	1:A:140:ALA:CB	2.46	0.46
4:D:32:PHE:HB3	4:D:73:LEU:HD22	1.97	0.46
4:D:61:ARG:HG2	4:D:78:VAL:O	2.16	0.45
4:D:138:VAL:HG11	5:E:149:LEU:HD22	1.98	0.45
1:A:44:ARG:HG2	1:A:44:ARG:HH11	1.81	0.45
2:B:11:SER:HB2	2:B:21:ASN:ND2	2.31	0.45
1:A:45:MET:HE1	3:C:2:LEU:HD11	1.98	0.45
5:E:29:HIS:HE1	5:E:107:TYR:CD2	2.35	0.45
1:A:158:ALA:HA	4:D:52:ILE:HD13	1.98	0.45
2:B:40:LEU:HD23	2:B:45:ARG:HA	1.99	0.45
4:D:85:SER:HA	4:D:112:LEU:O	2.17	0.45
4:D:143:ASP:OD1	5:E:198:ARG:NH2	2.49	0.45
1:A:51:TRP:HZ3	1:A:171:TYR:HB3	1.81	0.45
4:D:128:LEU:N	4:D:128:LEU:HD12	2.31	0.45
5:E:173:VAL:HA	5:E:196:ARG:O	2.17	0.45
1:A:259:CYS:O	1:A:271:THR:HA	2.16	0.44
2:B:51:HIS:HA	2:B:65:LEU:O	2.17	0.44
4:D:138:VAL:HG13	5:E:131:PHE:CE2	2.52	0.44
4:D:150:VAL:O	4:D:150:VAL:CG1	2.66	0.44



		Interatomic	Clash
Atom-1	Atom-2	distance $(\text{\AA})$	overlap (Å)
5:E:22:GLN:HA	5:E:22:GLN:OE1	2.17	0.44
5:E:32:MET:HA	5:E:93:ALA:O	2.17	0.44
5:E:170:HIS:O	5:E:173:VAL:HG13	2.17	0.44
4:D:81:GLN:O	4:D:84:ASP:HB2	2.17	0.44
1:A:168:LEU:O	1:A:172:LEU:HG	2.18	0.44
4:D:61:ARG:HG2	4:D:79:PRO:HD2	2.00	0.44
1:A:28:VAL:HG21	1:A:51:TRP:HH2	1.83	0.43
5:E:160:LEU:C	5:E:160:LEU:HD23	2.39	0.43
4:D:32:PHE:HB2	4:D:49:ILE:HG22	2.00	0.43
4:D:157:ASP:HB2	4:D:184:LYS:HE3	1.99	0.43
4:D:116:PRO:HG2	4:D:165:VAL:HG11	2.01	0.43
4:D:155:ASP:HB3	4:D:158:VAL:CG2	2.48	0.43
1:A:218:GLN:O	1:A:257:TYR:HA	2.18	0.43
4:D:10:LEU:HD23	4:D:112:LEU:HD13	2.00	0.43
1:A:133:TRP:HB2	1:A:144:LYS:HG3	2.00	0.43
2:B:2:GLN:HA	2:B:31:HIS:O	2.18	0.43
2:B:73:THR:HG21	2:B:75:LYS:HG2	1.99	0.43
5:E:14:LYS:HA	5:E:117:THR:O	2.17	0.43
4:D:17:ILE:CG1	4:D:78:VAL:HG22	2.30	0.43
5:E:65:TYR:CE2	5:E:90:TYR:HE2	2.37	0.43
1:A:249:VAL:HG13	1:A:257:TYR:HE1	1.84	0.42
1:A:249:VAL:HG13	1:A:257:TYR:CE1	2.54	0.42
4:D:49:ILE:HG21	4:D:73:LEU:HD11	2.01	0.42
4:D:93:MET:SD	4:D:104:LEU:CD2	3.06	0.42
5:E:80:LEU:HD23	5:E:80:LEU:HA	1.86	0.42
1:A:241:PHE:CD1	1:A:241:PHE:N	2.86	0.42
1:A:266:LEU:HD13	1:A:270:LEU:HG	2.02	0.42
4:D:46:LEU:O	4:D:47:ILE:HB	2.19	0.42
5:E:175:THR:HB	5:E:195:SER:OG	2.19	0.42
5:E:176:ASP:HA	5:E:177:PRO:HD3	1.83	0.42
4:D:127:GLN:HB2	6:D:219:HOH:O	2.19	0.42
4:D:165:VAL:HG21	6:D:211:HOH:O	2.19	0.42
3:C:4:GLY:C	3:C:5:TYR:CD1	2.93	0.42
1:A:165:VAL:HG11	1:A:169:ARG:NH2	2.34	0.42
4:D:198:ILE:HG22	4:D:199:PRO:N	2.34	0.42
4:D:141:PHE:O	4:D:177:SER:HA	2.21	0.41
1:A:47:PRO:HG3	1:A:60:TRP:CZ2	2.55	0.41
5:E:13:LEU:HD11	5:E:19:MET:CB	2.50	0.41
5:E:178:GLN:HA	5:E:179:PRO:HD3	1.93	0.41
4:D:34:TRP:CZ3	4:D:90:CYS:HB2	2.55	0.41
4:D:150:VAL:HG11	4:D:163:LYS:H	1.84	0.41



Atom 1	Atom 2	Interatomic	Clash
Atom-1	Atom-2	distance (Å)	overlap (Å)
1:A:52:ILE:HD12	1:A:52:ILE:HA	1.89	0.41
1:A:192:HIS:O	1:A:200:THR:N	2.52	0.41
4:D:161:THR:HG21	5:E:194:SER:OG	2.21	0.41
4:D:187:PHE:CD1	4:D:187:PHE:O	2.74	0.41
4:D:26:ASN:ND2	4:D:94:GLU:OE1	2.54	0.41
4:D:198:ILE:CG2	4:D:199:PRO:N	2.83	0.41
5:E:126:PRO:HD3	5:E:217:PHE:CD1	2.55	0.41
3:C:5:TYR:OH	4:D:31:TYR:HB2	2.21	0.41
2:B:37:VAL:HG22	2:B:82:VAL:HG22	2.03	0.40
5:E:49:SER:HB2	5:E:75:PHE:CE1	2.56	0.40
4:D:66:LEU:HD13	4:D:73:LEU:HD13	2.03	0.40
1:A:147:TRP:CZ2	3:C:9:VAL:HG12	2.56	0.40
4:D:85:SER:O	4:D:86:ALA:HB2	2.22	0.40
5:E:133:PRO:CG	5:E:146:LEU:HD12	2.49	0.40
2:B:51:HIS:HB3	2:B:66:TYR:CD2	2.56	0.40
5:E:232:LYS:HA	5:E:233:PRO:HD3	1.85	0.40
1:A:44:ARG:HA	1:A:64:THR:HG23	2.02	0.40
1:A:189:MET:SD	1:A:217:TRP:HH2	2.45	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	А	273/275~(99%)	250~(92%)	22 (8%)	1 (0%)	34 54
2	В	97/100~(97%)	91~(94%)	5 (5%)	1 (1%)	15 28
3	С	7/9~(78%)	7~(100%)	0	0	100 100
4	D	184/204~(90%)	161 (88%)	14 (8%)	9~(5%)	2 2
5	Ε	240/244~(98%)	224 (93%)	14 (6%)	2(1%)	19 35
All	All	801/832~(96%)	733~(92%)	55~(7%)	13~(2%)	9 17



Mol	Chain	Res	Type
4	D	192	ALA
4	D	195	ASN
4	D	199	PRO
4	D	58	GLY
5	Е	246	ALA
4	D	196	SER
4	D	202	THR
2	В	15	ALA
5	Е	156	ASP
1	А	136	ALA
4	D	150	VAL
4	D	47	ILE
4	D	197	ILE

All (13) Ramachandran outliers are listed below:

#### 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	Percer	ntiles
1	А	219/231~(95%)	208~(95%)	11 (5%)	24	46
2	В	89/95~(94%)	85~(96%)	4 (4%)	27	51
3	С	8/8~(100%)	8 (100%)	0	100	100
4	D	149/184~(81%)	131 (88%)	18 (12%)	5	9
5	Ε	191/209~(91%)	172~(90%)	19 (10%)	8	15
All	All	656/727~(90%)	604~(92%)	52 (8%)	12	24

All (52) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	А	6	ARG
1	А	23	ILE
1	А	35	ARG
1	А	86	ASN
1	А	132	SER



Mol	Chain	Res	Type
1	А	163	THR
1	А	196	ASP
1	А	212	GLU
1	А	216	THR
1	А	234	ARG
1	А	251	SER
2	В	34	ASP
2	В	61	SER
2	В	70	PHE
2	В	75	LYS
4	D	13	GLN
4	D	18	SER
4	D	27	SER
4	D	33	LEU
4	D	46	LEU
4	D	50	SER
4	D	69	SER
4	D	93	MET
4	D	104	LEU
4	D	108	GLN
4	D	147	GLN
4	D	150	VAL
4	D	153	SER
4	D	183	ASN
4	D	187	PHE
4	D	197	ILE
4	D	202	THR
4	D	203	PHE
5	Е	20	THR
5	Е	25	GLN
5	Ε	56	ASP
5	Е	67	VAL
5	Е	70	SER
5	Е	72	THR
5	Е	106	GLN
5	Е	112	THR
5	Е	117	THR
5	Е	118	GLU
5	Ε	128	VAL
5	Е	147	VAL
5	Е	173	VAL
5	Е	180	LEU



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Mol	Chain	Res	Type
5	Е	196	ARG
5	Е	198	ARG
5	Е	224	ASP
5	Е	239	SER
5	Е	247	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	А	72	GLN
1	А	87	GLN
1	А	224	GLN
1	А	255	GLN
1	А	263	HIS
4	D	5	GLN
4	D	26	ASN
4	D	55	ASN
4	D	67	ASN
4	D	76	HIS
4	D	183	ASN
4	D	195	ASN
5	Е	6	GLN
5	Е	11	GLN
5	Е	29	HIS
5	Е	86	GLN
5	Е	106	GLN
5	Е	157	HIS
5	Е	216	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 monosaccharides in this entry.



#### 5.6 Ligand geometry (i)

There are no ligands in this entry.

#### 5.7 Other polymers (i)

There are no such residues in this entry.

#### 5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.



### 6 Fit of model and data (i)

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

EDS was not executed - this section is therefore empty.

#### 6.2 Non-standard residues in protein, DNA, RNA chains (i)

EDS was not executed - this section is therefore empty.

#### 6.3 Carbohydrates (i)

EDS was not executed - this section is therefore empty.

#### 6.4 Ligands (i)

EDS was not executed - this section is therefore empty.

#### 6.5 Other polymers (i)

EDS was not executed - this section is therefore empty.

