



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:18 PM GMT

PDB ID : 3WS3  
Title : Crystal Structure of H-2D in complex with an insulin derived peptide  
Authors : Kumar, P.R.; Mukherjee, G.; Samanta, D.; DiLorenzo, T.P.; Almo, S.C.; Immune Function Network; New York Structural Genomics Research Consortium (NYSGRC)  
Deposited on : 2014-02-28  
Resolution : 2.33 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

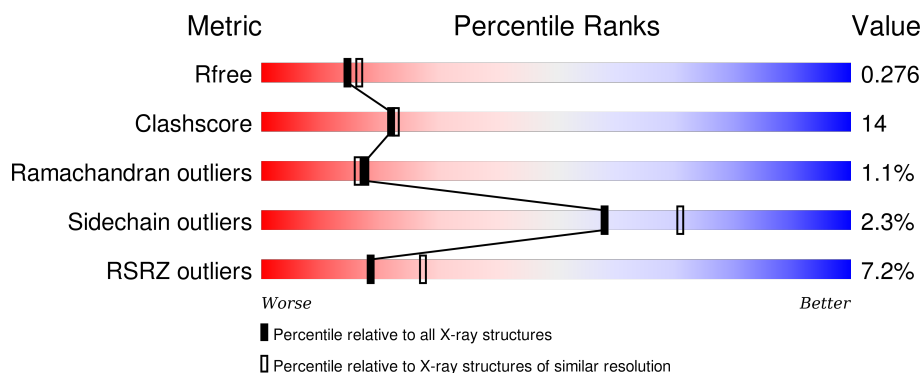
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.33 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	1406 (2.36-2.32)
Clashscore	102246	1509 (2.36-2.32)
Ramachandran outliers	100387	1490 (2.36-2.32)
Sidechain outliers	100360	1491 (2.36-2.32)
RSRZ outliers	91569	1412 (2.36-2.32)

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	273	<div> <div>8%</div> <div>63%</div> <div>31%</div> <div>...</div> </div>
1	C	273	<div> <div>10%</div> <div>70%</div> <div>27%</div> <div>..</div> </div>
2	B	100	<div> <div>3%</div> <div>73%</div> <div>27%</div> </div>
2	D	100	<div> <div>3%</div> <div>70%</div> <div>29%</div> <div>.</div> </div>
3	E	9	<div> <div>78%</div> <div>22%</div> </div>

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Mol	Chain	Length	Quality of chain
3	F	9	 56% 44%

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 6355 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 H-2 class I histocompatibility antigen, D-B alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	266	Total	C	N	O	S	0	1	0
			2205	1394	388	414	9			
1	C	271	Total	C	N	O	S	0	0	0
			2233	1410	395	419	9			

- Molecule 2 is a protein called Beta-2-microglobulin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	100	Total	C	N	O	S	0	0	0
			829	529	139	153	8			
2	D	100	Total	C	N	O	S	0	0	0
			829	529	139	153	8			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	20	MET	-	EXPRESSION TAG	UNP P01887
B	105	ASP	ALA	SEE REMARK 999	UNP P01887
D	20	MET	-	EXPRESSION TAG	UNP P01887
D	105	ASP	ALA	SEE REMARK 999	UNP P01887

- Molecule 3 is a protein called Insulin derived 9-mer peptide.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	9	Total	C	N	O	S	0	0	0
			77	49	11	16	1			
3	F	9	Total	C	N	O	S	0	0	0
			77	49	11	16	1			

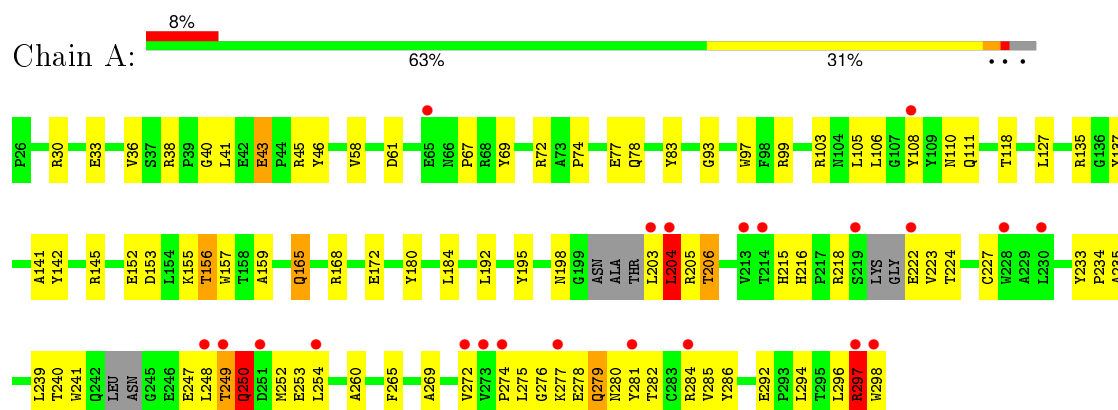
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	36	Total 36	O 36	0	0
4	B	11	Total 11	O 11	0	0
4	C	33	Total 33	O 33	0	0
4	D	23	Total 23	O 23	0	0
4	E	2	Total 2	O 2	0	0

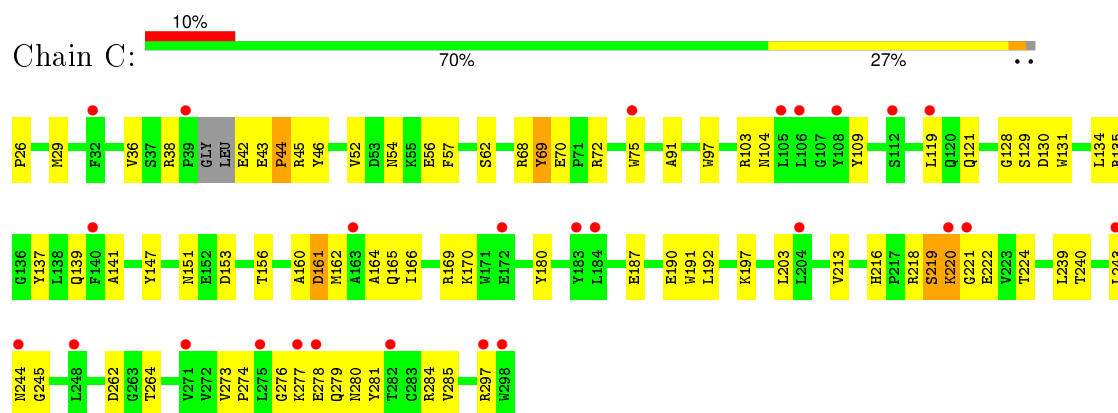
### 3 Residue-property plots [i](#)

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.

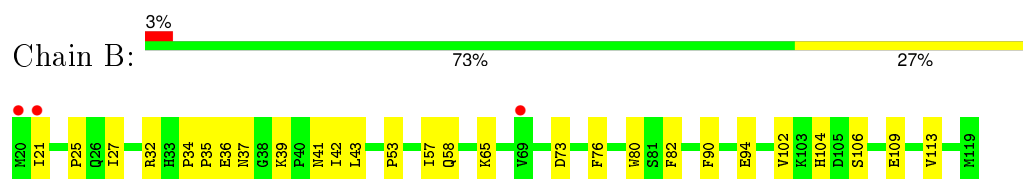
- Molecule 1: H-2 class I histocompatibility antigen, D-B alpha chain



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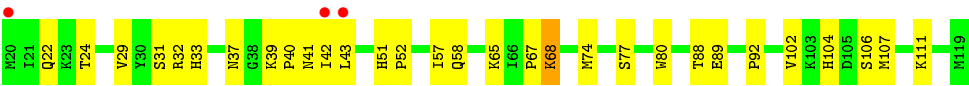


- Molecule 2: Beta-2-microglobulin



- Molecule 2: Beta-2-microglobulin





- Molecule 3: Insulin derived 9-mer peptide



- Molecule 3: Insulin derived 9-mer peptide



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	87.22Å 101.13Å 117.54Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.45 – 2.33 46.45 – 1.93	Depositor EDS
% Data completeness (in resolution range)	74.6 (46.45-2.33) 88.0 (46.45-1.93)	Depositor EDS
$R_{merge}$	0.15	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.19 (at 1.94Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.4_1496)	Depositor
R, $R_{free}$	0.217 , 0.294 0.269 , 0.276	Depositor DCC
$R_{free}$ test set	2196 reflections (5.17%)	DCC
Wilson B-factor (Å <sup>2</sup> )	25.1	Xtriage
Anisotropy	0.519	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 48.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	12 of 69183 reflections (0.017%)	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	6355	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 32.65 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 9.0861e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<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

### 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	0.55	1/2268 (0.0%)	0.77	4/3073 (0.1%)
1	C	0.49	0/2298	0.66	1/3117 (0.0%)
2	B	0.44	0/855	0.60	0/1158
2	D	0.49	0/855	0.66	0/1158
3	E	0.44	0/78	0.46	0/103
3	F	0.40	0/78	0.58	0/103
All	All	0.50	1/6432 (0.0%)	0.69	5/8712 (0.1%)

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	3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	297	ARG	CG-CD	-5.42	1.38	1.51

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	204	LEU	CB-CG-CD1	8.15	124.86	111.00
1	A	204	LEU	CA-CB-CG	7.60	132.78	115.30
1	A	297	ARG	NE-CZ-NH1	-7.28	116.66	120.30
1	C	130	ASP	N-CA-C	6.04	127.29	111.00
1	A	206	THR	N-CA-C	5.14	124.88	111.00

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	203	LEU	Peptide
1	A	204	LEU	Peptide
1	A	250	GLN	Peptide

## 5.2 Too-close contacts

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	2205	0	2069	85	0
1	C	2233	0	2103	54	0
2	B	829	0	802	16	1
2	D	829	0	802	19	0
3	E	77	0	70	2	0
3	F	77	0	70	5	0
4	A	36	0	0	5	0
4	B	11	0	0	1	0
4	C	33	0	0	4	1
4	D	23	0	0	2	0
4	E	2	0	0	0	0
All	All	6355	0	5916	169	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 14.

All (169) 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:145:ARG:HD2	2:B:21:ILE:HD13	1.55	0.87
1:C:129:SER:HA	1:C:131:TRP:H	1.42	0.84
1:C:103:ARG:NH2	4:C:317:HOH:O	2.11	0.80
1:C:68:ARG:NH1	1:C:70:GLU:OE2	2.15	0.79
1:A:279:GLN:H	1:A:297:ARG:NH2	1.83	0.76
1:A:249:THR:HG23	1:A:250:GLN:H	1.50	0.76
1:C:151:ASN:HD22	1:C:156:THR:HG23	1.51	0.76
1:A:285:VAL:HG22	1:A:294:LEU:HB2	1.69	0.75
1:C:244:ASN:N	1:C:245:GLY:HA3	2.02	0.74
1:C:279:GLN:O	1:C:297:ARG:NH1	2.21	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:56:GLU:OE2	1:C:72:ARG:NH1	2.21	0.74
2:B:104:HIS:ND1	2:B:106:SER:OG	2.20	0.73
1:A:135:ARG:HD3	1:A:137:TYR:CZ	2.24	0.73
2:D:31:SER:OG	2:D:33:HIS:O	2.08	0.72
1:A:281:TYR:O	1:A:297:ARG:NH1	2.26	0.69
1:A:33:GLU:OE2	3:E:2:GLN:NE2	2.26	0.68
1:A:106:LEU:HD23	1:A:111:GLN:HB2	1.78	0.65
1:C:43:GLU:HG3	1:C:44:PRO:HD2	1.78	0.64
1:C:161:ASP:OD1	4:C:326:HOH:O	2.15	0.64
2:D:67:PRO:O	4:D:211:HOH:O	2.14	0.64
1:C:219:SER:OG	1:C:220:LYS:N	2.19	0.64
1:C:160:ALA:O	1:C:164:ALA:HB3	1.98	0.64
1:A:204:LEU:HD22	1:A:205:ARG:H	1.62	0.63
1:C:129:SER:HA	1:C:131:TRP:N	2.14	0.62
1:A:38:ARG:HH21	1:A:45:ARG:HB2	1.64	0.62
1:C:29:MET:HB2	1:C:192:LEU:HD13	1.82	0.62
1:A:274:PRO:HG2	1:A:277:LYS:NZ	2.14	0.62
1:C:216:HIS:HB2	1:C:224:THR:HB	1.82	0.61
1:A:141:ALA:HB2	2:B:80:TRP:CE2	2.36	0.60
1:A:280:ASN:H	1:A:297:ARG:HH22	1.47	0.60
1:C:121:GLN:HE22	3:F:5:ASN:HD21	1.48	0.60
1:C:26:PRO:N	1:C:129:SER:HG	2.00	0.60
2:B:36:GLU:OE1	2:B:39:LYS:NZ	2.30	0.60
1:A:284:ARG:HG3	1:A:286:TYR:HE1	1.67	0.59
1:A:218:ARG:HD2	1:A:272:VAL:HG21	1.84	0.59
2:D:107:MET:SD	2:D:111:LYS:HG2	2.42	0.59
1:A:153:ASP:O	1:A:155:LYS:HG3	2.03	0.59
1:A:296:LEU:O	4:A:330:HOH:O	2.17	0.58
1:A:78:GLN:HE21	1:A:198:ASN:HB3	1.68	0.58
1:A:280:ASN:N	1:A:297:ARG:HH12	2.01	0.58
1:A:38:ARG:NH2	1:A:45:ARG:HB2	2.19	0.58
1:A:38:ARG:HH11	1:A:41:LEU:HD22	1.69	0.57
1:A:279:GLN:H	1:A:297:ARG:CZ	2.17	0.57
1:A:218:ARG:N	1:A:222:GLU:O	2.36	0.57
1:C:103:ARG:NH1	1:C:104:ASN:OD1	2.37	0.57
1:A:157:TRP:HB2	1:A:168:ARG:HG3	1.86	0.57
1:A:127:LEU:HG	1:A:192:LEU:HD23	1.87	0.56
1:C:219:SER:O	1:C:221:GLY:N	2.38	0.56
1:A:97:TRP:CZ2	3:E:8:GLY:HA2	2.41	0.56
1:C:141:ALA:HB2	2:D:80:TRP:CE2	2.40	0.56
1:A:156:THR:OG1	1:A:157:TRP:N	2.40	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:161:ASP:O	1:C:165:GLN:NE2	2.39	0.54
1:A:43:GLU:HG3	1:A:99:ARG:NH2	2.23	0.53
1:A:43:GLU:OE1	1:A:99:ARG:NE	2.41	0.53
2:D:39:LYS:HG3	2:D:40:PRO:HD2	1.89	0.53
1:C:187:GLU:N	1:C:187:GLU:OE1	2.30	0.53
1:A:135:ARG:HG3	1:A:152:GLU:HG3	1.91	0.52
1:A:45:ARG:NH2	1:A:61:ASP:OD2	2.40	0.52
1:C:162:MET:O	1:C:166:ILE:HG12	2.09	0.52
1:A:110:ASN:O	1:A:110:ASN:ND2	2.42	0.52
1:C:239:LEU:HD22	1:C:285:VAL:HG22	1.91	0.52
2:B:53:PRO:HD3	2:B:82:PHE:CE2	2.45	0.52
1:A:204:LEU:HD22	1:A:205:ARG:HG2	1.92	0.52
1:A:58:VAL:HB	1:A:69:TYR:CD1	2.45	0.52
1:C:276:GLY:N	1:C:278:GLU:OE2	2.41	0.52
1:C:109:TYR:OH	1:C:161:ASP:OD2	2.11	0.51
2:D:22:GLN:HB2	2:D:51:HIS:O	2.10	0.51
1:A:233:TYR:CD1	1:A:234:PRO:HA	2.45	0.51
1:A:43:GLU:HG3	1:A:99:ARG:HH21	1.75	0.51
1:A:165:GLN:HG3	1:A:168:ARG:HE	1.76	0.51
1:C:97:TRP:CZ2	3:F:8:GLY:HA2	2.46	0.51
1:A:284:ARG:HG3	1:A:286:TYR:CE1	2.46	0.51
2:D:41:ASN:OD1	2:D:42:ILE:N	2.35	0.51
1:A:250:GLN:HG2	1:A:252:MET:HB2	1.93	0.51
1:C:75:TRP:CH2	1:C:203:LEU:HD21	2.45	0.51
1:A:279:GLN:N	1:A:297:ARG:NH2	2.55	0.51
1:C:243:LEU:HD22	1:C:281:TYR:CZ	2.46	0.51
1:A:241:TRP:HA	1:A:282:THR:O	2.11	0.50
1:C:277:LYS:O	1:C:280:ASN:HB2	2.11	0.50
2:B:41:ASN:OD1	2:B:42:ILE:N	2.40	0.50
1:A:215:HIS:CD2	1:A:223:VAL:HG21	2.47	0.50
1:A:159:ALA:HB3	1:A:165:GLN:NE2	2.27	0.50
1:A:33:GLU:OE2	1:A:46:TYR:OH	2.31	0.49
1:C:213:VAL:HG12	4:C:325:HOH:O	2.12	0.49
1:A:279:GLN:NE2	1:A:298:TRP:OXT	2.45	0.49
1:A:281:TYR:N	1:A:297:ARG:NH1	2.61	0.49
1:A:78:GLN:NE2	1:A:198:ASN:HB3	2.28	0.49
2:B:73:ASP:O	4:B:201:HOH:O	2.20	0.49
1:C:191:TRP:CE2	3:F:1:TYR:CD1	3.01	0.49
2:B:25:PRO:HD3	2:B:106:SER:OG	2.13	0.49
1:A:103:ARG:O	1:A:106:LEU:HB2	2.12	0.49
1:A:204:LEU:HA	1:A:205:ARG:HD3	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:32:ARG:CZ	2:D:42:ILE:HD12	2.42	0.49
1:A:83:TYR:OH	1:A:195:TYR:OH	2.29	0.49
1:A:260:ALA:HB1	2:B:32:ARG:HG3	1.94	0.48
1:C:222:GLU:HA	1:C:274:PRO:HA	1.93	0.48
1:A:281:TYR:H	1:A:297:ARG:NH1	2.12	0.48
2:D:58:GLN:HG2	2:D:65:LYS:HD2	1.95	0.48
1:C:169:ARG:HG2	1:C:170:LYS:N	2.27	0.48
1:A:227:CYS:HB2	1:A:241:TRP:CZ2	2.49	0.47
1:A:40:GLY:O	4:A:325:HOH:O	2.20	0.47
1:A:74:PRO:O	1:A:77:GLU:HG3	2.13	0.47
1:A:274:PRO:HG2	1:A:277:LYS:HZ1	1.79	0.47
2:D:57:ILE:HG12	2:D:102:VAL:HG22	1.96	0.47
1:C:160:ALA:C	1:C:165:GLN:HE22	2.17	0.47
1:A:275:LEU:HD23	1:A:276:GLY:N	2.29	0.47
1:A:233:TYR:CG	1:A:234:PRO:HA	2.50	0.47
1:C:69:TYR:CE2	1:C:91:ALA:HB2	2.50	0.47
1:C:128:GLY:N	1:C:134:LEU:HG	2.30	0.47
1:A:105:LEU:HD13	1:A:142:TYR:CG	2.50	0.46
2:B:57:ILE:HG12	2:B:102:VAL:HG22	1.96	0.46
1:A:281:TYR:H	1:A:297:ARG:CZ	2.28	0.46
1:A:93:GLY:HA3	4:A:315:HOH:O	2.15	0.46
2:B:58:GLN:HB3	2:B:65:LYS:NZ	2.30	0.46
1:A:292:GLU:CD	1:C:197:LYS:HE2	2.36	0.46
1:C:180:TYR:HE1	3:F:5:ASN:HD22	1.64	0.46
1:A:281:TYR:O	1:A:297:ARG:HD2	2.16	0.46
1:A:218:ARG:HH11	1:A:218:ARG:HG2	1.81	0.46
2:D:22:GLN:HE21	2:D:106:SER:HA	1.81	0.46
1:A:105:LEU:HD23	1:A:108:TYR:CD2	2.50	0.46
1:C:38:ARG:NH2	1:C:45:ARG:HB2	2.30	0.46
1:C:240:THR:HG22	1:C:284:ARG:HB2	1.97	0.46
1:A:67:PRO:HG2	4:A:327:HOH:O	2.16	0.45
2:B:37:ASN:ND2	2:B:94:GLU:HG3	2.31	0.45
2:D:37:ASN:HA	2:D:92:PRO:O	2.16	0.45
1:A:30:ARG:NH2	1:A:137:TYR:CD2	2.85	0.45
2:D:29:VAL:HG12	2:D:43:LEU:HD11	1.99	0.45
1:A:281:TYR:H	1:A:297:ARG:NH2	2.15	0.45
2:B:34:PRO:HA	2:B:35:PRO:HD3	1.79	0.45
1:C:52:VAL:HG23	1:C:57:PHE:CD2	2.52	0.44
1:C:147:TYR:CE2	3:F:9:LEU:HD22	2.52	0.44
1:A:248:LEU:HA	1:A:248:LEU:HD12	1.67	0.44
2:B:76:PHE:HB3	2:B:82:PHE:CD1	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:180:TYR:O	1:A:184:LEU:HG	2.18	0.44
2:D:32:ARG:NH2	2:D:42:ILE:HD12	2.33	0.44
1:C:119:LEU:HD12	1:C:141:ALA:O	2.18	0.44
1:C:218:ARG:HG2	1:C:219:SER:HB3	2.00	0.43
1:A:240:THR:OG1	1:A:284:ARG:NH1	2.51	0.43
1:A:218:ARG:H	1:A:223:VAL:HA	1.84	0.43
1:C:273:VAL:HG23	1:C:274:PRO:O	2.19	0.43
1:A:234:PRO:HB2	4:A:323:HOH:O	2.19	0.43
2:D:88:THR:HG23	2:D:89:GLU:O	2.19	0.43
1:A:278:GLU:O	1:A:279:GLN:HB2	2.18	0.43
1:A:239:LEU:HD22	1:A:285:VAL:HG12	2.01	0.43
1:C:153:ASP:OD2	1:C:156:THR:HG22	2.19	0.42
1:A:216:HIS:HB2	1:A:224:THR:HB	2.00	0.42
1:C:243:LEU:HA	1:C:244:ASN:HA	1.68	0.42
2:D:24:THR:HA	2:D:106:SER:OG	2.19	0.42
1:A:36:VAL:HG13	1:A:118:THR:HG22	2.02	0.42
1:C:128:GLY:H	1:C:134:LEU:HG	1.84	0.42
1:C:54:ASN:ND2	4:C:303:HOH:O	2.20	0.42
2:D:68:LYS:HA	2:D:68:LYS:HD2	1.62	0.42
2:D:74:MET:HB2	4:D:208:HOH:O	2.19	0.42
2:D:52:PRO:HD2	2:D:104:HIS:CE1	2.55	0.42
1:C:262:ASP:OD1	1:C:264:THR:OG1	2.29	0.41
1:C:36:VAL:O	1:C:44:PRO:HA	2.20	0.41
1:A:83:TYR:HH	1:A:195:TYR:HH	1.60	0.41
1:A:135:ARG:NH2	1:A:152:GLU:HB2	2.36	0.41
1:C:46:TYR:H	1:C:62:SER:HB3	1.86	0.41
1:A:205:ARG:HD3	1:A:205:ARG:N	2.36	0.41
1:A:241:TRP:H	1:A:252:MET:HE1	1.86	0.41
1:A:253:GLU:O	1:A:269:ALA:HA	2.20	0.41
2:B:27:ILE:HB	2:B:113:VAL:HG21	2.02	0.41
1:A:72:ARG:HD2	1:A:72:ARG:HA	1.94	0.41
1:A:235:ALA:HB2	1:A:265:PHE:CE2	2.56	0.41
2:B:43:LEU:HB2	2:B:90:PHE:CD2	2.55	0.41
1:A:135:ARG:HD3	1:A:137:TYR:CE2	2.57	0.40
1:C:135:ARG:HD3	1:C:137:TYR:OH	2.21	0.40
1:C:46:TYR:H	1:C:62:SER:CB	2.35	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)
2:B:94:GLU:OE1	4:C:317:HOH:O[2_555]	2.06	0.14

## 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	259/273 (95%)	239 (92%)	16 (6%)	4 (2%)	13	10
1	C	267/273 (98%)	250 (94%)	13 (5%)	4 (2%)	13	10
2	B	98/100 (98%)	95 (97%)	3 (3%)	0	100	100
2	D	98/100 (98%)	95 (97%)	3 (3%)	0	100	100
3	E	7/9 (78%)	6 (86%)	1 (14%)	0	100	100
3	F	7/9 (78%)	6 (86%)	1 (14%)	0	100	100
All	All	736/764 (96%)	691 (94%)	37 (5%)	8 (1%)	17	16

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	206	THR
1	A	249	THR
1	C	220	LYS
1	A	279	GLN
1	C	44	PRO
1	C	219	SER
1	C	161	ASP
1	A	247	GLU

### 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	228/232 (98%)	220 (96%)	8 (4%)	43	55
1	C	231/232 (100%)	227 (98%)	4 (2%)	68	80
2	B	95/95 (100%)	94 (99%)	1 (1%)	80	89
2	D	95/95 (100%)	93 (98%)	2 (2%)	61	75
3	E	8/8 (100%)	8 (100%)	0	100	100
3	F	8/8 (100%)	8 (100%)	0	100	100
All	All	665/670 (99%)	650 (98%)	15 (2%)	58	71

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	43	GLU
1	A	156	THR
1	A	165	GLN
1	A	172	GLU
1	A	204	LEU
1	A	250	GLN
1	A	254	LEU
1	A	297	ARG
2	B	109	GLU
1	C	42	GLU
1	C	69	TYR
1	C	139	GLN
1	C	190	GLU
2	D	68	LYS
2	D	77	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	165	GLN
1	A	215	HIS
1	A	280	ASN
1	C	117	HIS
1	C	121	GLN
1	C	165	GLN
1	C	212	HIS

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Mol	Chain	Res	Type
2	D	22	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	266/273 (97%)	0.73	22 (8%)	14 22	32, 54, 110, 135	0
1	C	271/273 (99%)	0.68	26 (9%)	10 16	35, 65, 104, 131	0
2	B	100/100 (100%)	0.43	3 (3%)	54 64	36, 53, 76, 117	0
2	D	100/100 (100%)	0.62	3 (3%)	54 64	31, 48, 78, 137	0
3	E	9/9 (100%)	0.79	0	100 100	41, 51, 55, 61	0
3	F	9/9 (100%)	0.58	0	100 100	44, 51, 65, 69	0
All	All	755/764 (98%)	0.66	54 (7%)	18 27	31, 56, 103, 137	0

All (54) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	203	LEU	8.9
2	D	20	MET	7.8
1	A	204	LEU	6.5
1	C	204	LEU	5.7
1	C	277	LYS	5.6
1	C	297	ARG	5.4
1	A	281	TYR	4.9
1	A	273	VAL	4.2
1	A	214	THR	4.2
1	C	221	GLY	3.9
1	A	251	ASP	3.7
1	A	228	TRP	3.5
1	C	106	LEU	3.5
1	C	108	TYR	3.5
2	D	43	LEU	3.4
1	C	105	LEU	3.3
1	A	254	LEU	3.2
1	C	39	PRO	3.1
1	A	230	LEU	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	213	VAL	3.0
1	A	249	THR	2.9
1	C	282	THR	2.9
1	C	172	GLU	2.9
1	C	248	LEU	2.8
2	B	21	ILE	2.8
1	C	278	GLU	2.7
1	A	274	PRO	2.7
1	A	65	GLU	2.7
1	C	163	ALA	2.7
1	A	297	ARG	2.7
1	C	183	TYR	2.6
1	A	284	ARG	2.6
1	C	140	PHE	2.6
1	A	248	LEU	2.6
1	C	275	LEU	2.5
1	A	108	TYR	2.5
1	C	112	SER	2.5
1	C	271	VAL	2.5
2	B	69	VAL	2.5
1	C	32	PHE	2.4
1	C	244	ASN	2.4
1	C	298	TRP	2.4
1	C	243	LEU	2.4
1	A	219	SER	2.4
1	A	298	TRP	2.3
1	A	277	LYS	2.2
1	C	220	LYS	2.2
1	C	119	LEU	2.2
1	C	75	TRP	2.2
1	C	184	LEU	2.2
2	D	42	ILE	2.1
1	A	272	VAL	2.1
2	B	20	MET	2.0
1	A	222	GLU	2.0

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

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.