



# Full wwPDB X-ray Structure Validation Report ⓘ

Sep 9, 2023 – 02:50 PM EDT

PDB ID : 4HKJ  
Title : Structure of Cowpox CPXV203 in complex with MHCI (H-2Kb)  
Authors : McCoy IV, W.H.; Fremont, D.H.; Center for Structural Genomics of Infectious Diseases (CSGID)  
Deposited on : 2012-10-15  
Resolution : 3.00 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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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 : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.35.1  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.35.1

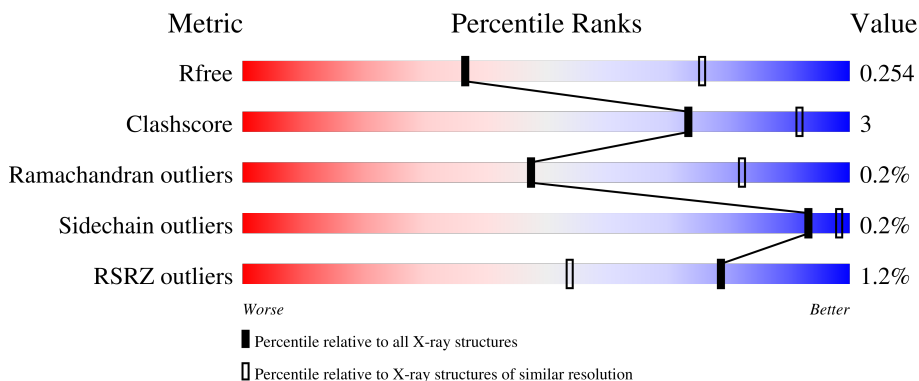
# 1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.00 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.





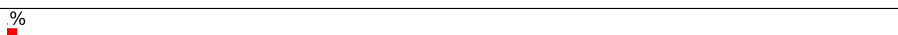
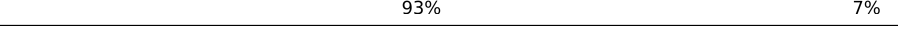


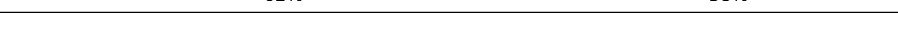



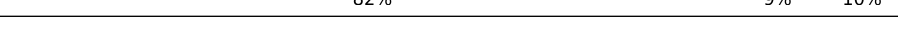
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	2092 (3.00-3.00)
Clashscore	141614	2416 (3.00-3.00)
Ramachandran outliers	138981	2333 (3.00-3.00)
Sidechain outliers	138945	2336 (3.00-3.00)
RSRZ outliers	127900	1990 (3.00-3.00)

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  $\geq 3$ , 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  $\leq 5\%$ . 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	280	 88% 11%
1	E	280	 88% 11%
1	I	280	 88% 11%
1	M	280	 88% 11%
2	B	100	 91% 9%

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Mol	Chain	Length	Quality of chain
2	F	100	 90% 10%
2	J	100	 4% 90% 10%
2	N	100	 % 93% 7%
3	C	8	 88% 12%
3	G	8	 62% 38%
3	K	8	 62% 38%
3	O	8	 88% 12%
4	D	206	 82% 9% 10%
4	H	206	 83% 7% 10%
4	L	206	 83% 8% 10%
4	P	206	 83% 7% 10%

## 2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 18660 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, K-B alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	277	2254	1423	396	426	9	0	0	0
1	E	277	2254	1423	396	426	9	0	0	0
1	I	277	2254	1423	396	426	9	0	0	0
1	M	277	2254	1423	396	426	9	0	0	0

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	100	837	533	141	159	4	0	0	0
2	F	100	837	533	141	159	4	0	0	0
2	J	100	837	533	141	159	4	0	0	0
2	N	100	837	533	141	159	4	0	0	0

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	0	MET	-	initiating methionine	UNP P61769
F	0	MET	-	initiating methionine	UNP P61769
J	0	MET	-	initiating methionine	UNP P61769
N	0	MET	-	initiating methionine	UNP P61769

- Molecule 3 is a protein called Ovalbumin.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	8	Total	C	N	O	0	0	0
			68	45	10	13			
3	G	8	Total	C	N	O	0	0	0
			68	45	10	13			
3	K	8	Total	C	N	O	0	0	0
			68	45	10	13			
3	O	8	Total	C	N	O	0	0	0
			68	45	10	13			

- Molecule 4 is a protein called CPXV203 protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
4	D	186	Total	C	N	O	S	Se	0	0	0
			1488	928	253	294	10	3			
4	H	186	Total	C	N	O	S	Se	0	0	0
			1488	928	253	294	10	3			
4	L	186	Total	C	N	O	S	Se	0	0	0
			1488	928	253	294	10	3			
4	P	186	Total	C	N	O	S	Se	0	0	0
			1488	928	253	294	10	3			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	0	ALA	-	expression tag	UNP Q8QMP2
H	0	ALA	-	expression tag	UNP Q8QMP2
L	0	ALA	-	expression tag	UNP Q8QMP2
P	0	ALA	-	expression tag	UNP Q8QMP2

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	14	Total	O	0	0
			14	14		
5	B	3	Total	O	0	0
			3	3		
5	C	1	Total	O	0	0
			1	1		
5	D	6	Total	O	0	0
			6	6		
5	E	5	Total	O	0	0
			5	5		

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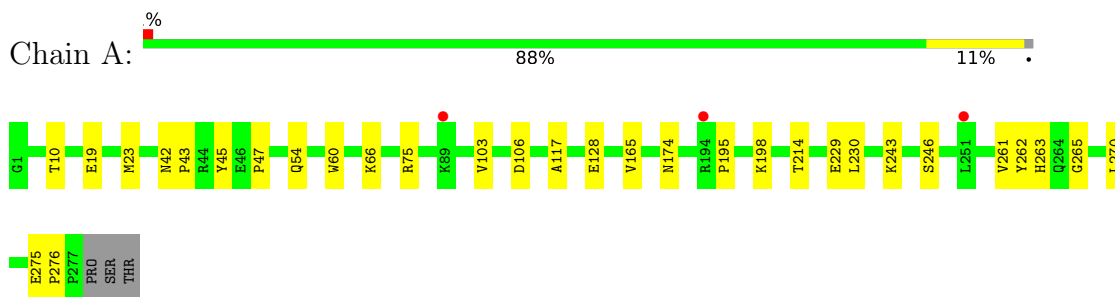
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<b>Mol</b>	<b>Chain</b>	<b>Residues</b>	<b>Atoms</b>	<b>ZeroOcc</b>	<b>AltConf</b>
5	F	3	Total O 3 3	0	0
5	H	4	Total O 4 4	0	0
5	I	5	Total O 5 5	0	0
5	J	4	Total O 4 4	0	0
5	K	1	Total O 1 1	0	0
5	L	4	Total O 4 4	0	0
5	M	11	Total O 11 11	0	0
5	N	4	Total O 4 4	0	0
5	P	7	Total O 7 7	0	0

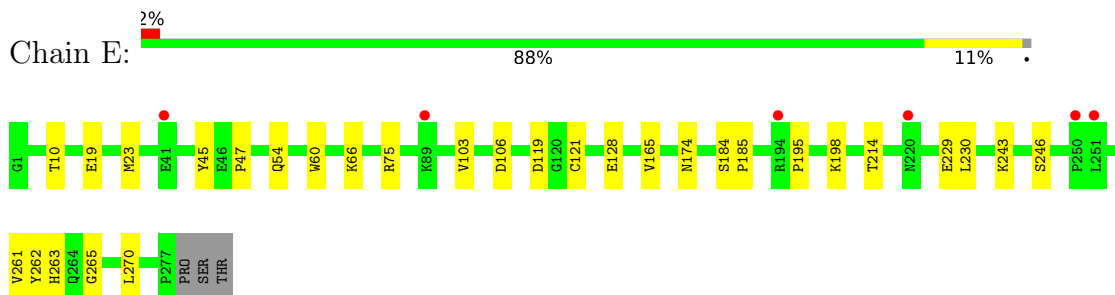
### 3 Residue-property plots [i](#)

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 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.

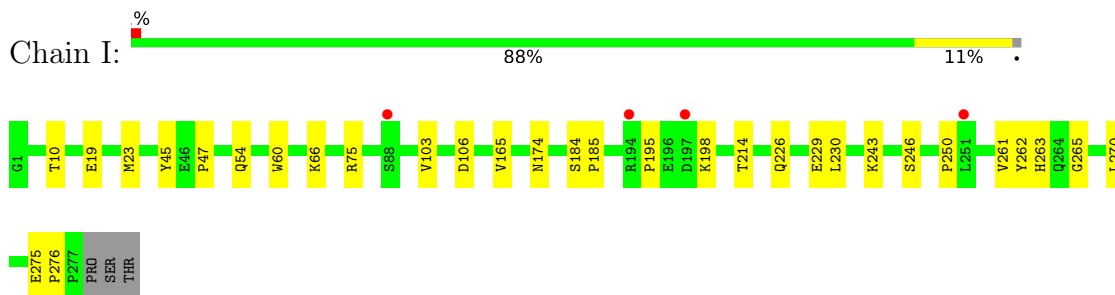
- Molecule 1: H-2 class I histocompatibility antigen, K-B alpha chain



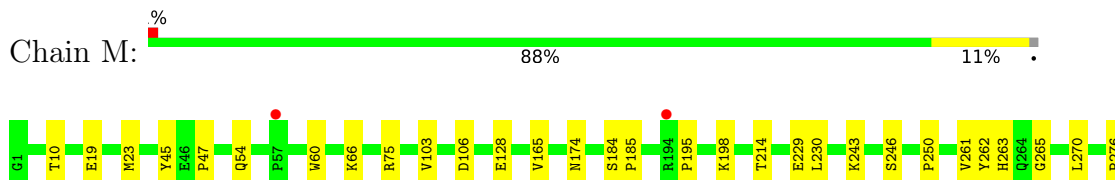
- Molecule 1: H-2 class I histocompatibility antigen, K-B alpha chain



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

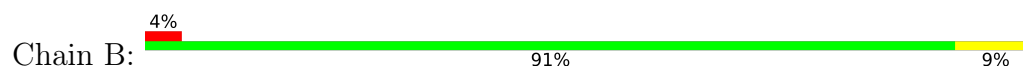


- Molecule 1: H-2 class I histocompatibility antigen, K-B alpha chain

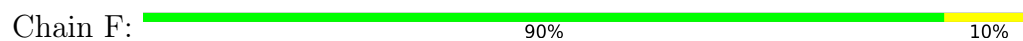




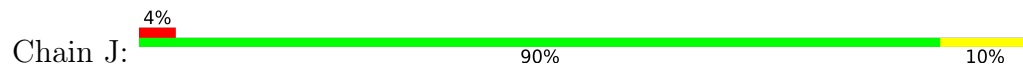
- Molecule 2: Beta-2-microglobulin



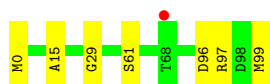
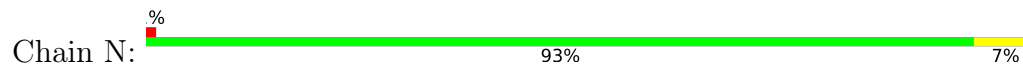
- Molecule 2: Beta-2-microglobulin



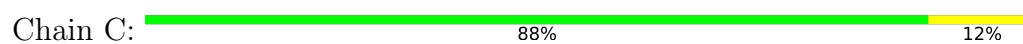
- Molecule 2: Beta-2-microglobulin



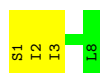
- Molecule 2: Beta-2-microglobulin



- Molecule 3: Ovalbumin



- Molecule 3: Ovalbumin



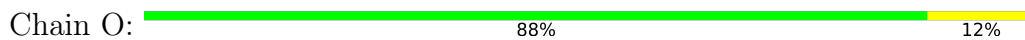
- Molecule 3: Ovalbumin



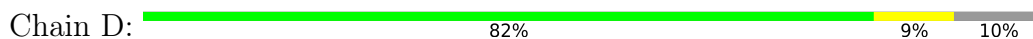




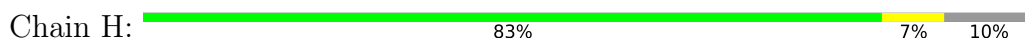
• Molecule 3: Ovalbumin



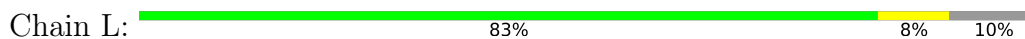
• Molecule 4: CPXV203 protein



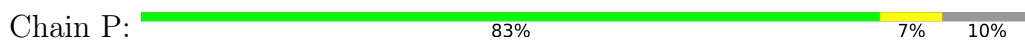
• Molecule 4: CPXV203 protein



• Molecule 4: CPXV203 protein



• Molecule 4: CPXV203 protein



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	88.31Å 88.25Å 106.42Å 76.18° 69.29° 66.69°	Depositor
Resolution (Å)	49.36 – 3.00 49.44 – 3.00	Depositor EDS
% Data completeness (in resolution range)	87.2 (49.36-3.00) 82.9 (49.44-3.00)	Depositor EDS
$R_{merge}$	0.16	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.09 (at 3.01Å)	Xtrriage
Refinement program	PHENIX 1.7.3_928	Depositor
R, $R_{free}$	0.230 , 0.253 0.233 , 0.254	Depositor DCC
$R_{free}$ test set	2398 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	65.4	Xtrriage
Anisotropy	0.247	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 76.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.45$ , $\langle L^2 \rangle = 0.28$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.90	EDS
Total number of atoms	18660	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	83.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.32% 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, 0.333 respectively for untwinned datasets, and 0.375, 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	0.20	0/2317	0.37	0/3149
1	E	0.20	0/2317	0.37	0/3149
1	I	0.20	0/2317	0.37	0/3149
1	M	0.20	0/2317	0.37	0/3149
2	B	0.21	0/860	0.37	0/1162
2	F	0.21	0/860	0.37	0/1162
2	J	0.21	0/860	0.37	0/1162
2	N	0.21	0/860	0.37	0/1162
3	C	0.22	0/68	0.33	0/88
3	G	0.21	0/68	0.32	0/88
3	K	0.22	0/68	0.33	0/88
3	O	0.22	0/68	0.33	0/88
4	D	0.21	0/1516	0.39	0/2044
4	H	0.21	0/1516	0.39	0/2044
4	L	0.21	0/1516	0.39	0/2044
4	P	0.21	0/1516	0.39	0/2044
All	All	0.21	0/19044	0.38	0/25772

There are no bond length outliers.

There are no bond angle outliers.

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	2254	0	2143	18	0
1	E	2254	0	2143	18	0
1	I	2254	0	2143	18	0
1	M	2254	0	2143	18	0
2	B	837	0	803	6	0
2	F	837	0	803	7	0
2	J	837	0	803	6	0
2	N	837	0	803	4	0
3	C	68	0	74	1	0
3	G	68	0	74	2	0
3	K	68	0	74	2	0
3	O	68	0	74	1	0
4	D	1488	0	1419	10	0
4	H	1488	0	1419	8	0
4	L	1488	0	1419	9	0
4	P	1488	0	1419	9	0
5	A	14	0	0	0	0
5	B	3	0	0	0	0
5	C	1	0	0	0	0
5	D	6	0	0	0	0
5	E	5	0	0	0	0
5	F	3	0	0	0	0
5	H	4	0	0	0	0
5	I	5	0	0	0	0
5	J	4	0	0	0	0
5	K	1	0	0	0	0
5	L	4	0	0	0	0
5	M	11	0	0	0	0
5	N	4	0	0	0	0
5	P	7	0	0	0	0
All	All	18660	0	17756	118	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 3.

All (118) 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:N:96:ASP:HB3	2:N:99:MET:HB2	1.79	0.65
1:I:262:TYR:HB3	1:M:106:ASP:HB2	1.78	0.65
2:F:96:ASP:HB3	2:F:99:MET:HB2	1.80	0.64
2:J:96:ASP:HB3	2:J:99:MET:HB2	1.79	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:96:ASP:HB3	2:B:99:MET:HB2	1.80	0.63
2:N:29:GLY:HA2	2:N:61:SER:HB3	1.82	0.62
2:B:29:GLY:HA2	2:B:61:SER:HB3	1.82	0.61
2:N:0:MET:N	4:P:125:GLU:O	2.33	0.61
2:J:29:GLY:HA2	2:J:61:SER:HB3	1.82	0.61
1:E:66:LYS:NZ	3:G:1:SER:OG	2.33	0.61
2:F:29:GLY:HA2	2:F:61:SER:HB3	1.83	0.60
1:I:66:LYS:NZ	3:K:1:SER:OG	2.36	0.58
1:M:195:PRO:HB2	1:M:198:LYS:HB2	1.88	0.56
2:J:0:MET:N	4:L:125:GLU:O	2.37	0.56
1:A:195:PRO:HB2	1:A:198:LYS:HB2	1.88	0.56
2:B:0:MET:N	4:D:125:GLU:O	2.39	0.55
1:E:195:PRO:HB2	1:E:198:LYS:HB2	1.88	0.55
1:A:262:TYR:HB3	1:E:106:ASP:HB2	1.89	0.55
1:I:195:PRO:HB2	1:I:198:LYS:HB2	1.88	0.55
1:E:214:THR:HB	1:E:262:TYR:HB2	1.91	0.53
1:A:106:ASP:HB2	1:E:262:TYR:HB3	1.90	0.53
1:M:229:GLU:HB3	1:M:246:SER:HB3	1.91	0.53
1:I:214:THR:HB	1:I:262:TYR:HB2	1.90	0.53
1:A:229:GLU:HB3	1:A:246:SER:HB3	1.91	0.53
1:I:229:GLU:HB3	1:I:246:SER:HB3	1.91	0.52
1:A:214:THR:HB	1:A:262:TYR:HB2	1.90	0.52
1:M:214:THR:HB	1:M:262:TYR:HB2	1.91	0.52
1:M:66:LYS:NZ	3:O:1:SER:OG	2.43	0.51
1:A:261:VAL:HB	1:A:270:LEU:HB2	1.93	0.51
1:I:106:ASP:HB2	1:M:262:TYR:HB3	1.92	0.51
1:E:261:VAL:HB	1:E:270:LEU:HB2	1.92	0.51
1:M:261:VAL:HB	1:M:270:LEU:HB2	1.93	0.51
1:E:229:GLU:HB3	1:E:246:SER:HB3	1.91	0.51
1:M:10:THR:HB	1:M:23:MET:HB2	1.93	0.51
1:A:66:LYS:NZ	3:C:1:SER:OG	2.44	0.50
1:A:10:THR:HB	1:A:23:MET:HB2	1.93	0.50
1:E:10:THR:HB	1:E:23:MET:HB2	1.93	0.50
1:I:261:VAL:HB	1:I:270:LEU:HB2	1.93	0.50
2:F:0:MET:N	4:H:125:GLU:O	2.44	0.49
1:I:10:THR:HB	1:I:23:MET:HB2	1.93	0.49
4:P:189:ASP:N	4:P:189:ASP:OD1	2.47	0.48
1:I:263:HIS:CD2	1:I:265:GLY:H	2.32	0.48
1:M:263:HIS:CD2	1:M:265:GLY:H	2.32	0.47
1:I:54:GLN:NE2	1:I:174:ASN:O	2.47	0.47
4:D:56:VAL:HG22	4:D:105:ILE:HG12	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:P:56:VAL:HG22	4:P:105:ILE:HG12	1.96	0.47
1:A:263:HIS:CD2	1:A:265:GLY:H	2.32	0.47
4:L:56:VAL:HG22	4:L:105:ILE:HG12	1.96	0.47
1:E:54:GLN:NE2	1:E:174:ASN:O	2.47	0.46
1:E:263:HIS:CD2	1:E:265:GLY:H	2.32	0.46
1:E:103:VAL:HG21	1:E:165:VAL:HG13	1.98	0.46
4:H:189:ASP:N	4:H:189:ASP:OD1	2.47	0.46
1:I:103:VAL:HG21	1:I:165:VAL:HG13	1.98	0.46
4:H:56:VAL:HG22	4:H:105:ILE:HG12	1.97	0.46
1:M:54:GLN:NE2	1:M:174:ASN:O	2.47	0.46
1:M:128:GLU:OE2	4:P:159:SER:OG	2.28	0.46
4:D:189:ASP:N	4:D:189:ASP:OD1	2.48	0.46
1:M:47:PRO:HG3	1:M:60:TRP:CZ2	2.51	0.46
1:A:47:PRO:HG3	1:A:60:TRP:CZ2	2.51	0.46
1:E:47:PRO:HG3	1:E:60:TRP:CZ2	2.51	0.46
1:I:47:PRO:HG3	1:I:60:TRP:CZ2	2.51	0.45
1:M:103:VAL:HG21	1:M:165:VAL:HG13	1.97	0.45
4:P:150:MSE:HE3	4:P:155:ARG:HG3	1.99	0.45
4:H:150:MSE:HE3	4:H:155:ARG:HG3	1.99	0.45
1:A:103:VAL:HG21	1:A:165:VAL:HG13	1.98	0.45
4:D:150:MSE:HE3	4:D:155:ARG:HG3	1.99	0.45
4:L:189:ASP:N	4:L:189:ASP:OD1	2.47	0.45
1:A:54:GLN:NE2	1:A:174:ASN:O	2.47	0.44
4:L:150:MSE:HE3	4:L:155:ARG:HG3	1.99	0.44
1:I:226:GLN:HA	4:L:76:PHE:HD1	1.82	0.44
1:M:230:LEU:HD11	1:M:243:LYS:HE3	2.00	0.43
1:A:230:LEU:HD11	1:A:243:LYS:HE3	2.00	0.43
1:A:19:GLU:OE1	1:A:75:ARG:NH2	2.52	0.43
1:A:128:GLU:OE2	4:D:159:SER:OG	2.33	0.43
1:E:19:GLU:OE1	1:E:75:ARG:NH2	2.51	0.43
4:P:152:GLU:OE1	4:P:155:ARG:NH2	2.45	0.43
1:E:230:LEU:HD11	1:E:243:LYS:HE3	2.00	0.43
1:I:19:GLU:OE1	1:I:75:ARG:NH2	2.52	0.43
1:E:184:SER:HA	1:E:185:PRO:HD3	1.89	0.42
2:B:15:ALA:HB3	2:B:97:ARG:HB2	2.01	0.42
1:I:230:LEU:HD11	1:I:243:LYS:HE3	2.01	0.42
2:J:15:ALA:HB3	2:J:97:ARG:HB2	2.01	0.42
1:M:19:GLU:OE1	1:M:75:ARG:NH2	2.52	0.42
2:N:15:ALA:HB3	2:N:97:ARG:HB2	2.01	0.42
2:F:15:ALA:HB3	2:F:97:ARG:HB2	2.01	0.41
1:A:42:ASN:HA	1:A:43:PRO:HD3	1.92	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:28:TYR:O	4:D:111:GLN:NE2	2.54	0.41
4:H:121:LEU:HD23	4:H:123:TYR:CZ	2.56	0.41
1:M:276:PRO:HA	1:M:277:PRO:HD3	1.89	0.41
4:D:128:TYR:HB3	4:D:149:PHE:HD2	1.85	0.41
1:E:121:CYS:SG	2:F:1:ILE:HG23	2.61	0.41
4:P:128:TYR:HB3	4:P:149:PHE:HD2	1.85	0.41
4:H:128:TYR:HB3	4:H:149:PHE:HD2	1.85	0.41
4:D:73:GLN:HG2	4:D:83:ILE:HG12	2.02	0.41
1:A:117:ALA:HB2	2:B:60:TRP:CE2	2.55	0.41
4:D:121:LEU:HD23	4:D:123:TYR:CZ	2.56	0.41
1:E:119:ASP:HB3	2:F:0:MET:HB2	2.03	0.41
4:L:128:TYR:HB3	4:L:149:PHE:HD2	1.85	0.41
4:P:121:LEU:HD23	4:P:123:TYR:CZ	2.56	0.41
3:G:2:ILE:HG12	3:G:3:ILE:H	1.86	0.41
1:I:198:LYS:HG2	1:I:250:PRO:HA	2.02	0.41
1:I:275:GLU:HA	1:I:276:PRO:HD2	1.98	0.41
2:F:24:ASN:HB3	2:F:65:LEU:HD11	2.03	0.41
1:M:184:SER:HA	1:M:185:PRO:HD3	1.89	0.40
4:P:28:TYR:O	4:P:111:GLN:NE2	2.54	0.40
2:J:24:ASN:HB3	2:J:65:LEU:HD11	2.04	0.40
4:L:25:LYS:HB2	4:L:141:ASN:HB2	2.03	0.40
1:E:128:GLU:OE2	4:H:159:SER:OG	2.36	0.40
4:H:28:TYR:O	4:H:111:GLN:NE2	2.54	0.40
1:I:184:SER:HA	1:I:185:PRO:HD3	1.89	0.40
4:L:121:LEU:HD23	4:L:123:TYR:CZ	2.56	0.40
1:A:275:GLU:HA	1:A:276:PRO:HD2	1.98	0.40
3:K:2:ILE:HG12	3:K:3:ILE:H	1.86	0.40
2:B:38:ASP:OD1	2:B:38:ASP:N	2.55	0.40
4:D:25:LYS:HB2	4:D:141:ASN:HB2	2.04	0.40
2:J:38:ASP:OD1	2:J:38:ASP:N	2.55	0.40
4:L:28:TYR:O	4:L:111:GLN:NE2	2.54	0.40
1:M:198:LYS:HG2	1:M:250:PRO:HA	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	275/280 (98%)	267 (97%)	8 (3%)	0	100	100
1	E	275/280 (98%)	267 (97%)	8 (3%)	0	100	100
1	I	275/280 (98%)	267 (97%)	8 (3%)	0	100	100
1	M	275/280 (98%)	267 (97%)	8 (3%)	0	100	100
2	B	98/100 (98%)	93 (95%)	5 (5%)	0	100	100
2	F	98/100 (98%)	94 (96%)	4 (4%)	0	100	100
2	J	98/100 (98%)	94 (96%)	4 (4%)	0	100	100
2	N	98/100 (98%)	94 (96%)	4 (4%)	0	100	100
3	C	6/8 (75%)	6 (100%)	0	0	100	100
3	G	6/8 (75%)	6 (100%)	0	0	100	100
3	K	6/8 (75%)	6 (100%)	0	0	100	100
3	O	6/8 (75%)	6 (100%)	0	0	100	100
4	D	184/206 (89%)	178 (97%)	5 (3%)	1 (0%)	29	68
4	H	184/206 (89%)	177 (96%)	6 (3%)	1 (0%)	29	68
4	L	184/206 (89%)	177 (96%)	6 (3%)	1 (0%)	29	68
4	P	184/206 (89%)	177 (96%)	6 (3%)	1 (0%)	29	68
All	All	2252/2376 (95%)	2176 (97%)	72 (3%)	4 (0%)	47	82

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
4	D	63	PRO
4	H	63	PRO
4	L	63	PRO
4	P	63	PRO

### 5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was



analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	235/238 (99%)	234 (100%)	1 (0%)	91	97
1	E	235/238 (99%)	234 (100%)	1 (0%)	91	97
1	I	235/238 (99%)	234 (100%)	1 (0%)	91	97
1	M	235/238 (99%)	234 (100%)	1 (0%)	91	97
2	B	95/95 (100%)	95 (100%)	0	100	100
2	F	95/95 (100%)	95 (100%)	0	100	100
2	J	95/95 (100%)	95 (100%)	0	100	100
2	N	95/95 (100%)	95 (100%)	0	100	100
3	C	8/8 (100%)	8 (100%)	0	100	100
3	G	8/8 (100%)	8 (100%)	0	100	100
3	K	8/8 (100%)	8 (100%)	0	100	100
3	O	8/8 (100%)	8 (100%)	0	100	100
4	D	170/184 (92%)	170 (100%)	0	100	100
4	H	170/184 (92%)	170 (100%)	0	100	100
4	L	170/184 (92%)	170 (100%)	0	100	100
4	P	170/184 (92%)	170 (100%)	0	100	100
All	All	2032/2100 (97%)	2028 (100%)	4 (0%)	93	98

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	45	TYR
1	E	45	TYR
1	I	45	TYR
1	M	45	TYR

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

Mol	Chain	Res	Type
4	D	73	GLN
4	D	75	HIS
4	H	73	GLN
4	H	75	HIS
4	L	73	GLN

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Mol	Chain	Res	Type
4	L	75	HIS
4	P	73	GLN
4	P	75	HIS

### 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

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	277/280 (98%)	-0.19	3 (1%) 80 56	44, 80, 125, 157	0
1	E	277/280 (98%)	-0.18	6 (2%) 62 33	43, 79, 125, 157	0
1	I	277/280 (98%)	-0.16	4 (1%) 75 49	45, 80, 125, 157	0
1	M	277/280 (98%)	-0.20	2 (0%) 87 69	44, 81, 125, 157	0
2	B	100/100 (100%)	0.06	4 (4%) 38 15	50, 89, 137, 145	0
2	F	100/100 (100%)	-0.19	0 100 100	49, 88, 138, 145	0
2	J	100/100 (100%)	0.04	4 (4%) 38 15	50, 88, 138, 145	0
2	N	100/100 (100%)	-0.02	1 (1%) 82 59	51, 98, 143, 155	0
3	C	8/8 (100%)	0.02	0 100 100	60, 74, 90, 90	0
3	G	8/8 (100%)	0.06	0 100 100	61, 74, 90, 92	0
3	K	8/8 (100%)	0.06	0 100 100	61, 74, 89, 90	0
3	O	8/8 (100%)	-0.16	0 100 100	55, 74, 89, 89	0
4	D	183/206 (88%)	-0.31	1 (0%) 91 75	48, 72, 117, 137	0
4	H	183/206 (88%)	-0.27	1 (0%) 91 75	48, 72, 117, 137	0
4	L	183/206 (88%)	-0.26	1 (0%) 91 75	48, 72, 117, 138	0
4	P	183/206 (88%)	-0.28	0 100 100	50, 74, 118, 137	0
All	All	2272/2376 (95%)	-0.18	27 (1%) 79 54	43, 79, 127, 157	0

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	I	194	ARG	3.9
2	B	21	ASN	3.6
1	M	57	PRO	3.6
2	J	75	LYS	3.2
2	N	68	THR	3.1

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Mol	Chain	Res	Type	RSRZ
1	E	194	ARG	3.1
2	B	68	THR	3.0
1	I	88	SER	3.0
1	M	194	ARG	2.9
4	L	41	VAL	2.9
1	A	194	ARG	2.7
1	I	197	ASP	2.7
1	E	251	LEU	2.7
1	I	251	LEU	2.5
1	E	220	ASN	2.5
2	B	16	GLU	2.5
1	E	250	PRO	2.4
2	J	21	ASN	2.4
4	H	112	GLU	2.4
1	A	251	LEU	2.3
2	J	16	GLU	2.3
1	E	41	GLU	2.2
1	E	89	LYS	2.2
2	B	72	PRO	2.2
1	A	89	LYS	2.2
2	J	70	PHE	2.2
4	D	43	GLY	2.1

## 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 monosaccharides 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.