



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 09:59 PM GMT

PDB ID : 1RJZ
Title : Mhc Class I Natural Mutant H-2Kbm8 Heavy Chain Complexed With beta-2 Microglobulin and Herpies Simplex Virus Mutant Glycoprotein B Peptide
Authors : Miley, M.J.; Messaoudi, I.; Nikolich-Zugich, J.; Fremont, D.H.
Deposited on : 2003-11-20
Resolution : 2.60 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

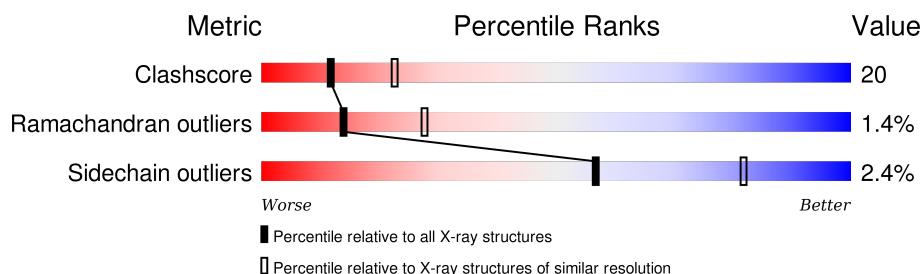
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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))
Clashscore	102246	2679 (2.60-2.60)
Ramachandran outliers	100387	2635 (2.60-2.60)
Sidechain outliers	100360	2635 (2.60-2.60)

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 ≥ 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	280	
1	D	280	
2	B	100	
2	E	100	
3	P	8	
3	Q	8	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6538 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
1	A	278	Total	C	N	O	S	0	0	0
			2257	1427	398	424	8			
1	D	278	Total	C	N	O	S	0	0	0
			2257	1427	398	424	8			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	22	PHE	TYR	ENGINEERED	UNP P01901
A	23	ILE	MET	ENGINEERED	UNP P01901
A	24	SER	GLU	ENGINEERED	UNP P01901
A	30	ASN	ASP	ENGINEERED	UNP P01901
D	22	PHE	TYR	ENGINEERED	UNP P01901
D	23	ILE	MET	ENGINEERED	UNP P01901
D	24	SER	GLU	ENGINEERED	UNP P01901
D	30	ASN	ASP	ENGINEERED	UNP P01901

- 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	E	100	Total	C	N	O	S	0	0	0
			829	529	139	153	8			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	0	MET	-	CLONING ARTIFACT	UNP P01887
E	0	MET	-	CLONING ARTIFACT	UNP P01887

- Molecule 3 is a protein called Glycoprotein B.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	P	8	Total	C	N	O	0	0	0
			68	43	11	14			
3	Q	8	Total	C	N	O	0	0	0
			68	43	11	14			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
P	2	GLU	SER	ENGINEERED	UNP P06436
Q	2	GLU	SER	ENGINEERED	UNP P06436

- Molecule 4 is water.

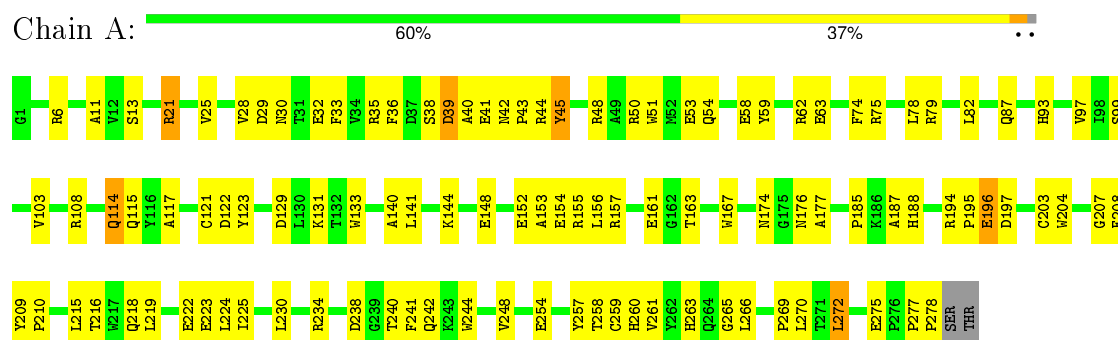
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	78	Total	O	0	0
			78	78		
4	B	40	Total	O	0	0
			40	40		
4	D	70	Total	O	0	0
			70	70		
4	E	33	Total	O	0	0
			33	33		
4	P	4	Total	O	0	0
			4	4		
4	Q	5	Total	O	0	0
			5	5		

3 Residue-property plots

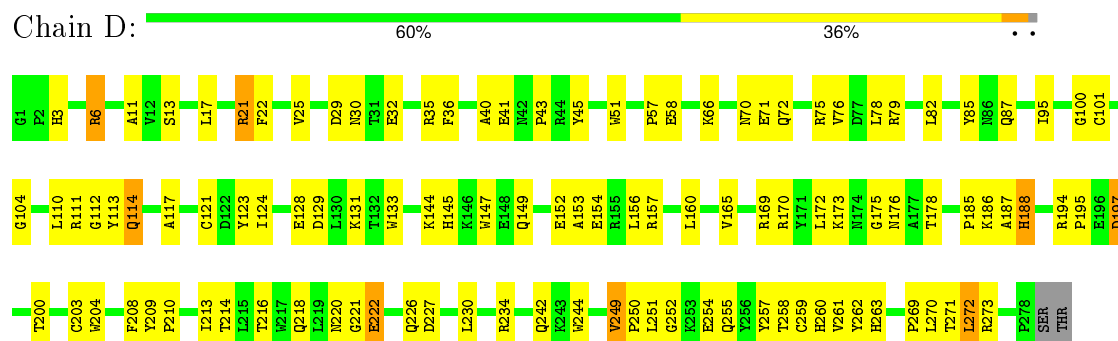
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

Note EDS was not executed.

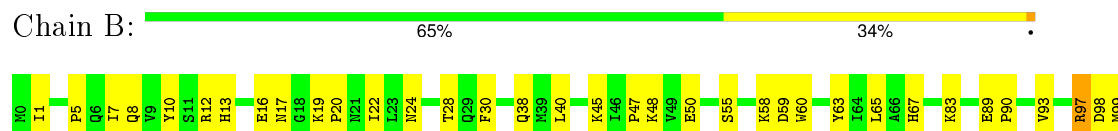
- Molecule 1: H-2 class I histocompatibility antigen, K-B alpha chain



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



- Molecule 2: Beta-2-microglobulin

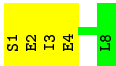


- Molecule 2: Beta-2-microglobulin





● Molecule 3: Glycoprotein B



● Molecule 3: Glycoprotein B



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	66.57Å 90.18Å 88.98Å 90.00° 111.32° 90.00°	Depositor
Resolution (Å)	19.94 – 2.60	Depositor
% Data completeness (in resolution range)	85.7 (19.94-2.60)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.230 , 0.288	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	6538	wwPDB-VP
Average B, all atoms (Å ²)	35.0	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).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.37	0/2321	0.61	0/3156
1	D	0.38	0/2321	0.62	0/3156
2	B	0.42	0/855	0.67	0/1158
2	E	0.40	0/855	0.66	0/1158
3	P	0.49	0/68	0.67	0/88
3	Q	0.52	0/68	0.66	0/88
All	All	0.39	0/6488	0.63	0/8804

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	2257	0	2153	112	0
1	D	2257	0	2153	97	0
2	B	829	0	805	26	0
2	E	829	0	805	24	0
3	P	68	0	68	7	0
3	Q	68	0	68	5	0
4	A	78	0	0	7	0
4	B	40	0	0	2	0
4	D	70	0	0	4	0
4	E	33	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	P	4	0	0	0	0
4	Q	5	0	0	0	0
All	All	6538	0	6052	249	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 20.

All (249) 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:121:CYS:SG	2:B:1:ILE:HG13	2.04	0.97
2:B:98:ASP:O	2:B:99:MET:HG3	1.70	0.91
1:A:141:LEU:HD11	1:D:57:PRO:HB2	1.57	0.86
1:D:6:ARG:HH22	2:E:58:LYS:NZ	1.76	0.83
1:A:234:ARG:HH11	2:B:8:GLN:NE2	1.77	0.83
1:A:21:ARG:HH11	1:A:21:ARG:HG3	1.44	0.83
1:D:21:ARG:HH11	1:D:21:ARG:HG3	1.44	0.81
1:A:234:ARG:HH11	2:B:8:GLN:HE22	1.29	0.81
1:D:121:CYS:SG	2:E:1:ILE:HG13	2.20	0.81
1:D:261:VAL:HB	1:D:270:LEU:HB2	1.61	0.80
1:D:194:ARG:HB3	1:D:195:PRO:HD2	1.64	0.80
1:A:194:ARG:HB3	1:A:195:PRO:HD2	1.66	0.78
1:D:170:ARG:HA	1:D:173:LYS:HE3	1.66	0.77
1:D:25:VAL:HG13	1:D:32:GLU:HG3	1.68	0.76
2:B:7:ILE:HB	2:B:93:VAL:HG21	1.68	0.75
1:D:234:ARG:HE	1:D:242:GLN:NE2	1.85	0.74
1:A:6:ARG:HD2	4:B:107:HOH:O	1.86	0.74
2:B:83:LYS:HG3	2:B:90:PRO:HG3	1.68	0.74
1:D:197:ASP:O	1:D:251:LEU:HB2	1.91	0.69
1:A:187:ALA:HA	1:A:204:TRP:O	1.92	0.69
1:A:261:VAL:HB	1:A:270:LEU:HB2	1.73	0.69
2:B:28:THR:HG22	2:B:63:TYR:HB2	1.74	0.67
2:E:28:THR:HG22	2:E:63:TYR:HB2	1.76	0.67
1:D:185:PRO:HB3	1:D:208:PHE:HB3	1.77	0.67
1:A:196:GLU:HG3	1:A:197:ASP:H	1.60	0.66
1:D:234:ARG:HE	1:D:242:GLN:HE21	1.43	0.66
1:A:185:PRO:HB3	1:A:208:PHE:HB3	1.77	0.66
1:D:259:CYS:HB3	1:D:272:LEU:CD1	2.26	0.66
1:A:21:ARG:NH1	1:A:21:ARG:HG3	2.10	0.65
1:A:144:LYS:O	1:A:148:GLU:HG3	1.96	0.65
1:A:272:LEU:N	1:A:272:LEU:HD12	2.11	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:6:ARG:HH22	2:E:58:LYS:HZ1	1.45	0.65
1:D:160:LEU:O	1:D:165:VAL:HG23	1.97	0.65
1:D:170:ARG:HH11	1:D:170:ARG:HG3	1.63	0.64
1:A:114:GLN:HA	1:A:114:GLN:HE21	1.62	0.64
1:D:104:GLY:N	1:D:110:LEU:HG	2.12	0.64
2:B:24:ASN:HB3	2:B:65:LEU:HD11	1.79	0.64
1:D:117:ALA:HB2	2:E:60:TRP:CE2	2.34	0.63
1:A:194:ARG:CZ	1:A:195:PRO:HD2	2.29	0.62
1:A:75:ARG:CZ	1:A:79:ARG:HH12	2.12	0.62
1:A:263:HIS:CD2	1:A:265:GLY:H	2.17	0.62
1:A:129:ASP:O	1:A:131:LYS:HG3	2.00	0.62
1:D:209:TYR:CD1	1:D:210:PRO:HA	2.35	0.62
1:A:114:GLN:HG2	1:A:156:LEU:CD1	2.31	0.60
1:A:42:ASN:O	1:A:44:ARG:NE	2.33	0.60
1:A:75:ARG:NH1	1:A:79:ARG:HH12	2.00	0.60
1:A:225:ILE:O	1:A:225:ILE:HG22	2.01	0.60
1:D:249:VAL:HB	1:D:257:TYR:CE1	2.37	0.60
1:A:114:GLN:HG2	1:A:156:LEU:HD13	1.83	0.59
2:B:12:ARG:NH1	2:B:22:ILE:HD12	2.17	0.59
1:A:82:LEU:HD21	1:A:93:HIS:CD2	2.37	0.59
1:D:129:ASP:O	1:D:131:LYS:HG3	2.02	0.59
1:D:25:VAL:CG1	1:D:32:GLU:HG3	2.34	0.58
1:D:220:ASN:HB2	4:D:305:HOH:O	2.03	0.58
1:D:21:ARG:NH1	1:D:21:ARG:HG3	2.13	0.58
1:D:145:HIS:O	1:D:149:GLN:HG2	2.04	0.57
1:D:6:ARG:HH22	2:E:58:LYS:HZ2	1.53	0.57
1:D:269:PRO:HD3	4:D:289:HOH:O	2.05	0.57
1:A:195:PRO:O	1:A:196:GLU:HB3	2.05	0.57
1:A:63:GLU:OE2	3:P:1:SER:HB2	2.05	0.57
2:B:38:GLN:NE2	2:B:45:LYS:HD2	2.20	0.57
1:A:154:GLU:OE1	1:A:154:GLU:N	2.39	0.56
1:D:262:TYR:CD1	1:D:269:PRO:HB3	2.40	0.56
1:A:260:HIS:HA	1:A:270:LEU:O	2.05	0.56
1:A:36:PHE:CZ	1:A:43:PRO:HB2	2.41	0.56
1:A:218:GLN:HG3	1:A:260:HIS:CD2	2.41	0.55
1:A:218:GLN:HG3	1:A:260:HIS:HD2	1.71	0.55
1:A:266:LEU:CD1	1:A:270:LEU:HG	2.36	0.55
1:D:6:ARG:NH2	2:E:58:LYS:NZ	2.50	0.55
1:A:50:ARG:O	1:A:53:GLU:HG3	2.07	0.55
1:D:255:GLN:HB2	4:D:290:HOH:O	2.04	0.55
1:A:82:LEU:HD22	1:A:87:GLN:HB2	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:194:ARG:HG2	1:A:194:ARG:HH11	1.71	0.54
2:B:17:ASN:OD1	2:B:97:ARG:NH2	2.41	0.54
1:A:103:VAL:HB	1:A:108:ARG:O	2.08	0.54
1:A:216:THR:HA	4:A:312:HOH:O	2.07	0.54
1:A:75:ARG:NH1	1:A:79:ARG:HH22	2.05	0.54
1:D:172:LEU:O	1:D:176:ASN:HB2	2.07	0.54
1:D:6:ARG:HB3	1:D:100:GLY:HA3	1.90	0.54
4:A:358:HOH:O	3:P:3:ILE:HG22	2.08	0.54
1:A:207:GLY:HA2	1:A:240:THR:CB	2.38	0.53
1:A:187:ALA:CB	1:A:272:LEU:HD21	2.39	0.53
1:A:117:ALA:HB2	2:B:60:TRP:CE2	2.44	0.53
1:D:133:TRP:HB2	1:D:144:LYS:HG3	1.90	0.53
1:D:194:ARG:HB3	1:D:195:PRO:CD	2.37	0.53
1:A:13:SER:HB3	1:A:78:LEU:HD13	1.90	0.53
1:D:176:ASN:C	1:D:178:THR:H	2.11	0.53
1:A:196:GLU:CG	1:A:197:ASP:H	2.21	0.53
3:P:3:ILE:HG12	3:P:4:GLU:N	2.24	0.53
1:D:70:ASN:CG	3:Q:5:PHE:HD2	2.12	0.52
1:D:194:ARG:HH11	1:D:194:ARG:HG2	1.73	0.52
1:A:82:LEU:CD2	1:A:87:GLN:HB2	2.39	0.52
1:A:6:ARG:NH1	4:A:335:HOH:O	2.42	0.52
1:D:25:VAL:HG21	1:D:35:ARG:NH2	2.25	0.52
1:A:266:LEU:HD13	1:A:270:LEU:HG	1.91	0.51
1:A:157:ARG:O	1:A:161:GLU:HG3	2.10	0.51
1:A:219:LEU:HD13	1:A:257:TYR:CZ	2.46	0.51
1:D:36:PHE:CZ	1:D:43:PRO:HB2	2.46	0.51
1:A:230:LEU:HD23	1:A:230:LEU:N	2.26	0.51
1:A:28:VAL:HG23	1:A:33:PHE:CD1	2.45	0.51
1:D:259:CYS:HB3	1:D:272:LEU:HD11	1.93	0.51
1:D:13:SER:HB3	1:D:78:LEU:HD13	1.92	0.51
1:A:35:ARG:HB3	1:A:48:ARG:HD3	1.93	0.51
2:E:15:PRO:HG3	2:E:97:ARG:HB2	1.93	0.50
1:A:242:GLN:HG2	4:A:325:HOH:O	2.11	0.50
1:D:213:ILE:HG12	1:D:214:THR:N	2.26	0.50
1:D:6:ARG:NH1	1:D:30:ASN:OD1	2.44	0.50
1:A:194:ARG:HG2	4:A:319:HOH:O	2.11	0.50
1:A:25:VAL:CG1	1:A:32:GLU:HG3	2.41	0.50
1:A:21:ARG:NE	1:A:39:ASP:HB3	2.26	0.50
1:D:114:GLN:HG2	1:D:156:LEU:CD1	2.42	0.50
1:A:123:TYR:CZ	1:A:140:ALA:HA	2.46	0.50
1:A:54:GLN:OE1	1:A:174:ASN:ND2	2.43	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:223:GLU:OE1	1:A:225:ILE:HD11	2.13	0.49
1:A:54:GLN:OE1	1:A:174:ASN:HB3	2.11	0.49
1:A:254:GLU:HB2	4:A:334:HOH:O	2.11	0.49
1:D:218:GLN:NE2	1:D:260:HIS:CD2	2.80	0.49
1:A:99:SER:CB	1:A:114:GLN:HE22	2.26	0.49
1:D:70:ASN:CG	3:Q:5:PHE:CD2	2.86	0.49
1:A:187:ALA:HB1	1:A:272:LEU:HD21	1.95	0.49
2:E:5:PRO:HG3	2:E:30:PHE:HB3	1.95	0.48
1:A:207:GLY:HA2	1:A:240:THR:HB	1.95	0.48
1:D:131:LYS:HZ1	1:D:157:ARG:HH12	1.59	0.48
1:A:224:LEU:HD11	1:A:257:TYR:HE2	1.78	0.48
1:A:74:PHE:HZ	1:A:97:VAL:HG21	1.78	0.48
1:D:75:ARG:NH1	1:D:79:ARG:HH22	2.11	0.48
2:E:28:THR:HG22	2:E:63:TYR:CB	2.44	0.48
1:A:133:TRP:HH2	1:A:156:LEU:HD12	1.78	0.48
1:D:131:LYS:NZ	1:D:157:ARG:HH12	2.11	0.48
2:B:10:TYR:CD2	2:B:10:TYR:N	2.82	0.48
1:A:196:GLU:HG3	1:A:197:ASP:N	2.28	0.48
1:D:11:ALA:HB3	1:D:95:ILE:CG2	2.44	0.47
2:B:40:LEU:HD23	2:B:45:LYS:HA	1.95	0.47
1:D:187:ALA:O	1:D:188:HIS:HB3	2.13	0.47
1:D:123:TYR:HD2	1:D:124:ILE:HG22	1.79	0.47
1:D:114:GLN:HG2	1:D:156:LEU:HD13	1.95	0.47
1:D:170:ARG:NH1	1:D:170:ARG:HG3	2.28	0.47
1:A:275:GLU:OE1	1:A:275:GLU:HA	2.15	0.47
2:E:40:LEU:HD22	2:E:43:GLY:O	2.13	0.47
1:D:230:LEU:N	1:D:230:LEU:HD23	2.29	0.47
1:D:194:ARG:HD2	1:D:200:THR:OG1	2.14	0.47
2:E:55:SER:HB3	4:E:125:HOH:O	2.15	0.47
1:D:114:GLN:HE21	1:D:114:GLN:HA	1.79	0.47
1:A:38:SER:C	1:A:40:ALA:H	2.18	0.47
1:A:156:LEU:HD23	3:P:3:ILE:HD13	1.97	0.46
1:A:194:ARG:HB3	1:A:195:PRO:CD	2.42	0.46
1:D:110:LEU:O	1:D:111:ARG:HB2	2.15	0.46
1:D:117:ALA:HB2	2:E:60:TRP:CZ2	2.51	0.46
1:D:258:THR:HG22	1:D:273:ARG:HG2	1.97	0.46
1:A:207:GLY:HA2	1:A:240:THR:OG1	2.16	0.46
1:D:66:LYS:HZ1	3:Q:1:SER:HB2	1.80	0.46
1:A:6:ARG:NH2	1:A:30:ASN:OD1	2.48	0.46
2:E:7:ILE:HB	2:E:93:VAL:HG21	1.98	0.46
1:A:194:ARG:HB3	1:A:194:ARG:CZ	2.46	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:260:HIS:CE1	1:D:271:THR:HG1	2.33	0.45
2:B:59:ASP:C	2:B:59:ASP:OD1	2.54	0.45
1:A:188:HIS:CE1	1:A:204:TRP:HB2	2.51	0.45
1:D:147:TRP:HB3	1:D:152:GLU:HB3	1.97	0.45
2:B:28:THR:HG22	2:B:63:TYR:CB	2.42	0.45
1:A:215:LEU:CD2	1:A:261:VAL:HG13	2.46	0.45
2:E:37:ILE:HD12	2:E:51:MET:CE	2.47	0.45
1:A:156:LEU:HD23	3:P:3:ILE:CD1	2.47	0.45
1:D:169:ARG:HG2	1:D:173:LYS:HE2	1.98	0.45
1:D:185:PRO:HD3	1:D:263:HIS:CD2	2.52	0.45
1:D:249:VAL:HG23	1:D:250:PRO:O	2.17	0.45
1:A:224:LEU:HD11	1:A:257:TYR:CE2	2.52	0.45
2:B:5:PRO:HB3	2:B:30:PHE:HB3	1.98	0.45
1:A:259:CYS:HB3	1:A:272:LEU:HD13	1.99	0.44
2:B:50:GLU:HB2	2:B:67:HIS:CE1	2.52	0.44
1:D:226:GLN:HG2	1:D:226:GLN:O	2.17	0.44
1:A:25:VAL:HG13	1:A:32:GLU:HG3	2.00	0.44
1:D:216:THR:OG1	1:D:260:HIS:HB2	2.18	0.44
2:E:38:GLN:NE2	2:E:45:LYS:HD2	2.32	0.44
1:A:62:ARG:HH22	1:A:167:TRP:HH2	1.65	0.44
1:D:144:LYS:O	1:D:145:HIS:C	2.56	0.44
1:A:230:LEU:HD23	1:A:230:LEU:H	1.83	0.44
1:A:195:PRO:O	1:A:196:GLU:CB	2.65	0.44
1:A:218:GLN:HB3	1:A:222:GLU:H	1.82	0.44
1:D:3:HIS:HA	1:D:29:ASP:OD1	2.18	0.44
1:D:252:GLY:N	1:D:254:GLU:OE1	2.51	0.44
2:B:12:ARG:HG2	2:B:13:HIS:CD2	2.53	0.44
1:A:121:CYS:O	1:A:122:ASP:C	2.55	0.44
1:D:22:PHE:CB	1:D:71:GLU:HG3	2.48	0.44
1:D:194:ARG:CZ	1:D:195:PRO:HD2	2.48	0.43
1:A:28:VAL:O	1:A:29:ASP:HB2	2.18	0.43
1:A:152:GLU:CD	1:A:155:ARG:HH21	2.21	0.43
1:D:221:GLY:O	1:D:222:GLU:HB2	2.18	0.43
2:B:19:LYS:HA	2:B:20:PRO:HD3	1.93	0.43
1:D:75:ARG:CZ	1:D:79:ARG:HH12	2.32	0.43
2:E:23:LEU:HB2	2:E:70:PHE:CD1	2.52	0.43
1:D:51:TRP:CD2	1:D:175:GLY:HA3	2.54	0.43
2:E:5:PRO:HD3	2:E:86:SER:OG	2.18	0.43
3:P:3:ILE:CG1	3:P:4:GLU:N	2.82	0.43
1:D:104:GLY:HA3	1:D:110:LEU:HD21	1.99	0.43
1:A:115:GLN:NE2	4:A:333:HOH:O	2.51	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:101:CYS:HA	1:D:112:GLY:HA2	2.01	0.43
1:D:194:ARG:NH1	1:D:194:ARG:HG2	2.33	0.43
1:A:272:LEU:N	1:A:272:LEU:CD1	2.79	0.43
2:E:38:GLN:HE22	2:E:45:LYS:HD2	1.82	0.43
1:A:163:THR:O	1:A:167:TRP:HD1	2.02	0.42
1:A:58:GLU:O	1:A:59:TYR:C	2.57	0.42
1:A:11:ALA:HB2	1:A:74:PHE:HD1	1.85	0.42
1:D:45:TYR:OH	3:Q:2:GLU:OE1	2.22	0.42
1:A:203:CYS:O	1:A:244:TRP:HA	2.19	0.42
1:D:186:LYS:HE3	4:D:346:HOH:O	2.18	0.42
1:A:216:THR:OG1	1:A:260:HIS:HB2	2.20	0.42
1:A:238:ASP:C	1:A:238:ASP:OD1	2.57	0.42
1:D:72:GLN:O	1:D:76:VAL:HG23	2.19	0.42
1:A:209:TYR:CD1	1:A:210:PRO:HA	2.55	0.42
1:A:6:ARG:NH1	2:B:58:LYS:HE2	2.35	0.42
2:E:55:SER:HB2	2:E:63:TYR:CZ	2.55	0.42
1:D:203:CYS:O	1:D:244:TRP:HA	2.20	0.42
2:B:89:GLU:HG2	2:B:90:PRO:HD2	2.01	0.42
1:D:11:ALA:HB3	1:D:95:ILE:HG23	2.02	0.42
1:D:185:PRO:CB	1:D:208:PHE:HB3	2.46	0.42
2:E:7:ILE:HG22	2:E:8:GLN:N	2.35	0.42
2:E:70:PHE:CE2	2:E:72:PRO:HB3	2.55	0.42
1:D:153:ALA:HB3	1:D:154:GLU:OE1	2.19	0.42
1:D:111:ARG:NH2	1:D:128:GLU:HG2	2.34	0.41
1:A:45:TYR:OH	3:P:2:GLU:OE1	2.24	0.41
1:D:82:LEU:CD2	1:D:87:GLN:HB2	2.50	0.41
1:D:111:ARG:CZ	1:D:128:GLU:HG2	2.50	0.41
1:A:11:ALA:CB	1:A:74:PHE:HD1	2.33	0.41
2:B:16:GLU:OE2	2:B:19:LYS:HD2	2.20	0.41
1:D:259:CYS:HB3	1:D:272:LEU:HD12	2.02	0.41
1:A:51:TRP:HB2	1:A:174:ASN:O	2.21	0.41
1:D:85:TYR:HB2	1:D:87:GLN:HG3	2.03	0.41
1:A:114:GLN:HA	1:A:114:GLN:NE2	2.33	0.41
1:D:40:ALA:O	1:D:41:GLU:C	2.58	0.41
2:B:58:LYS:NZ	4:B:107:HOH:O	2.54	0.41
1:A:114:GLN:HG2	1:A:156:LEU:HD11	2.03	0.41
1:D:187:ALA:HA	1:D:204:TRP:O	2.20	0.41
1:D:22:PHE:CG	1:D:71:GLU:HG3	2.56	0.41
2:E:10:TYR:CD2	2:E:10:TYR:N	2.89	0.41
2:B:89:GLU:CG	2:B:90:PRO:HD2	2.51	0.41
1:A:218:GLN:HA	1:A:222:GLU:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:45:TYR:HD1	1:A:63:GLU:HB3	1.85	0.41
1:D:66:LYS:HD2	3:Q:2:GLU:HG3	2.03	0.41
1:A:176:ASN:O	1:A:177:ALA:C	2.57	0.41
1:D:58:GLU:OE1	1:D:58:GLU:N	2.46	0.41
1:A:153:ALA:HB3	1:A:154:GLU:OE1	2.21	0.41
2:B:55:SER:HB2	2:B:63:TYR:CZ	2.56	0.40
1:A:99:SER:HB2	1:A:114:GLN:HE22	1.87	0.40
1:A:277:PRO:HA	1:A:278:PRO:HD2	1.96	0.40
1:A:208:PHE:CE1	1:A:241:PHE:HB2	2.57	0.40
1:A:194:ARG:HD3	1:A:248:VAL:HG13	2.03	0.40
1:A:266:LEU:HD11	1:A:270:LEU:HG	2.02	0.40
1:D:131:LYS:NZ	1:D:157:ARG:NH1	2.70	0.40
2:E:19:LYS:HA	2:E:20:PRO:HD3	1.95	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	276/280 (99%)	249 (90%)	23 (8%)	4 (1%)	14	28
1	D	276/280 (99%)	249 (90%)	24 (9%)	3 (1%)	17	36
2	B	98/100 (98%)	89 (91%)	7 (7%)	2 (2%)	9	18
2	E	98/100 (98%)	94 (96%)	2 (2%)	2 (2%)	9	18
3	P	6/8 (75%)	6 (100%)	0	0	100	100
3	Q	6/8 (75%)	6 (100%)	0	0	100	100
All	All	760/776 (98%)	693 (91%)	56 (7%)	11 (1%)	14	28

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	47	PRO
2	E	47	PRO
1	A	196	GLU
1	A	269	PRO
1	D	17	LEU
1	D	222	GLU
2	E	85	ASP
1	A	41	GLU
1	D	188	HIS
2	B	48	LYS
1	A	39	ASP

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	236/238 (99%)	231 (98%)	5 (2%)	61	85
1	D	236/238 (99%)	228 (97%)	8 (3%)	44	72
2	B	95/95 (100%)	94 (99%)	1 (1%)	80	93
2	E	95/95 (100%)	93 (98%)	2 (2%)	61	85
3	P	7/7 (100%)	7 (100%)	0	100	100
3	Q	7/7 (100%)	7 (100%)	0	100	100
All	All	676/680 (99%)	660 (98%)	16 (2%)	57	82

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

Mol	Chain	Res	Type
1	A	21	ARG
1	A	45	TYR
1	A	114	GLN
1	A	258	THR
1	A	272	LEU
2	B	97	ARG
1	D	6	ARG
1	D	21	ARG

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Mol	Chain	Res	Type
1	D	113	TYR
1	D	114	GLN
1	D	197	ASP
1	D	227	ASP
1	D	249	VAL
1	D	272	LEU
2	E	40	LEU
2	E	74	GLU

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	72	GLN
1	A	114	GLN
1	A	127	ASN
1	A	242	GLN
1	A	263	HIS
2	B	8	GLN
2	B	38	GLN
1	D	72	GLN
1	D	114	GLN
1	D	115	GLN
1	D	127	ASN
1	D	174	ASN
1	D	226	GLN
1	D	242	GLN
1	D	260	HIS
1	D	264	GLN
2	E	2	GLN
2	E	17	ASN
2	E	29	GLN
2	E	38	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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 ⓘ

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.