



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 11:04 PM GMT

PDB ID : 1WBZ
Title : CRYSTAL STRUCTURES OF MURINE MHC CLASS I H-2 DB AND KB MOLECULES IN COMPLEX WITH CTL EPITOPES FROM INFLUENZA A VIRUS: IMPLICATIONS FOR TCR REPERTOIRE SELECTION AND IMMUNODOMINANCE
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Deposited on : 2004-11-05
Resolution : 2.00 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
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) : 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

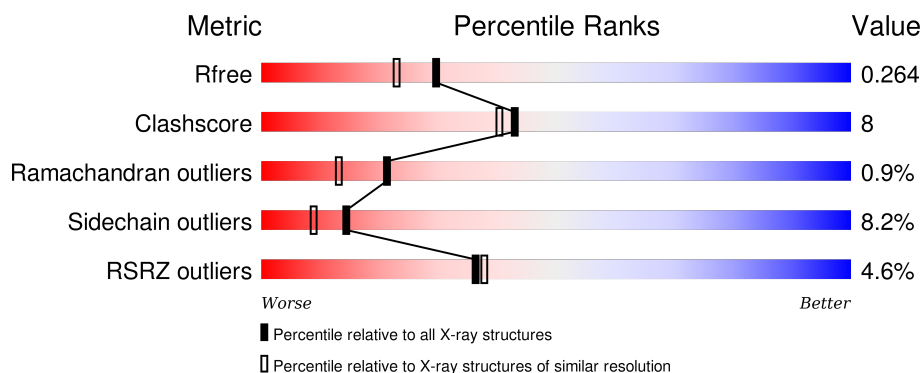
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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.



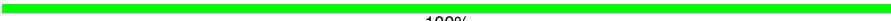
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	6249 (2.00-2.00)
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (2.00-2.00)

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 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.

Mol	Chain	Length	Quality of chain
1	A	275	<div> <div>4%</div> <div>84%</div> <div>14%</div> <div>.</div> </div>
1	C	275	<div> <div>8%</div> <div>76%</div> <div>19%</div> <div>.</div> <div>.</div> </div>
2	B	99	<div> <div>81%</div> <div>15%</div> <div>.</div> <div>.</div> </div>
2	D	99	<div> <div>4%</div> <div>79%</div> <div>15%</div> <div>6%</div> </div>
3	P	9	<div> <div>100%</div> </div>

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Mol	Chain	Length	Quality of chain
3	Q	9	 100%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 6952 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 PRECURSOR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	275	Total	C	N	O	S	0	0	0
			2240	1413	394	424	9			
1	C	275	Total	C	N	O	S	0	0	1
			2232	1408	394	421	9			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	99	Total	C	N	O	S	0	1	0
			826	527	138	154	7			
2	D	99	Total	C	N	O	S	0	0	0
			821	524	138	152	7			

- Molecule 3 is a protein called INFLUENZA A PEPTIDE.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	P	9	Total	C	N	O	0	0	0
			72	45	15	12			
3	Q	9	Total	C	N	O	0	0	0
			72	45	15	12			

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	234	Total	O	0	0
			234	234		
4	B	115	Total	O	0	0
			115	115		
4	C	226	Total	O	0	0
			226	226		
4	D	104	Total	O	0	0
			104	104		

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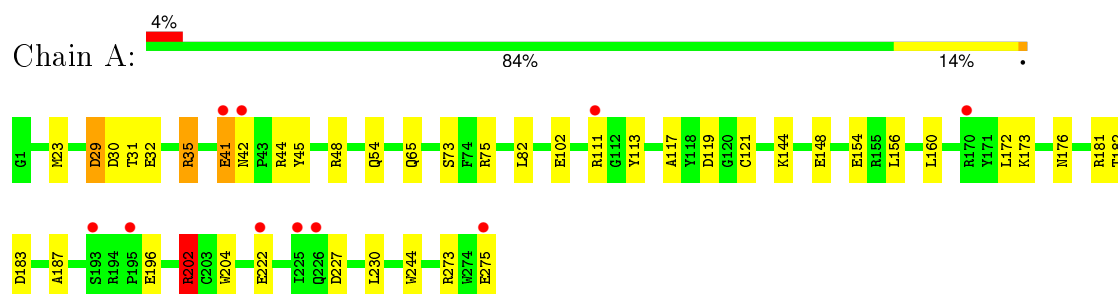
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	P	6	Total	O	0	0
			6	6		
4	Q	4	Total	O	0	0
			4	4		

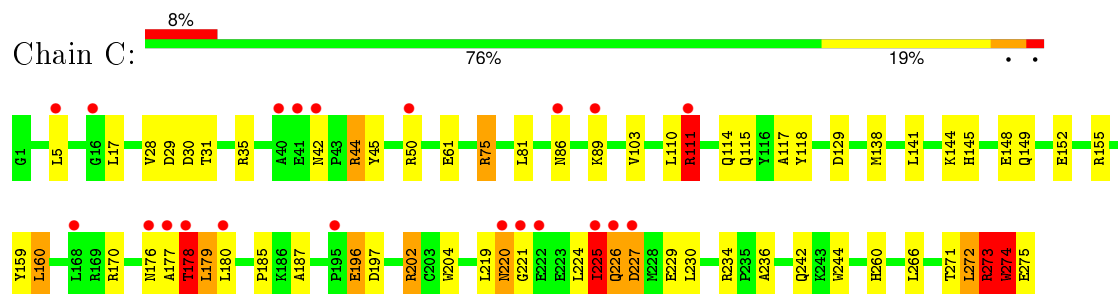
3 Residue-property plots

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.

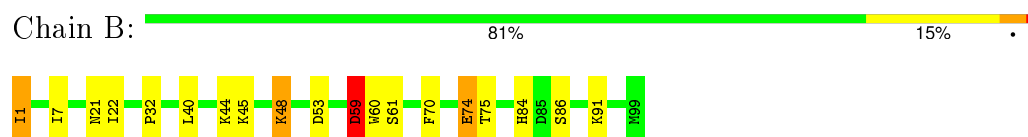
- Molecule 1: H-2 CLASS I HISTOCOMPATIBILITY ANTIGEN, K-B ALPHA CHAIN PRE-CURSOR



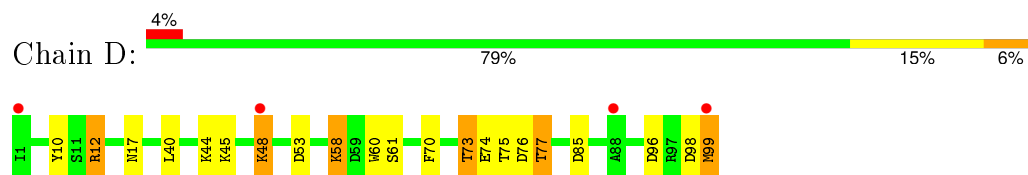
- Molecule 1: H-2 CLASS I HISTOCOMPATIBILITY ANTIGEN, K-B ALPHA CHAIN PRE-CURSOR



- Molecule 2: BETA-2MICROGLOBULIN



- Molecule 2: BETA-2MICROGLOBULIN



- Molecule 3: INFLUENZA A PEPTIDE

Chain P:  100%

There are no outlier residues recorded for this chain.

- Molecule 3: INFLUENZA A PEPTIDE

Chain Q:  100%

There are no outlier residues recorded for this chain.

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	89.17Å 91.26Å 67.18Å 90.00° 111.37° 90.00°	Depositor
Resolution (Å)	20.00 – 2.00 19.99 – 2.00	Depositor EDS
% Data completeness (in resolution range)	95.0 (20.00-2.00) 95.0 (19.99-2.00)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.22 (at 2.01Å)	Xtriage
Refinement program	REFMAC 5.2.0003	Depositor
R, R_{free}	0.208 , 0.254 0.213 , 0.264	Depositor DCC
R_{free} test set	3221 reflections (5.31%)	DCC
Wilson B-factor (Å ²)	29.5	Xtriage
Anisotropy	0.051	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 48.9	EDS
Estimated twinning fraction	0.019 for -h-l,-k,l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	1 of 63884 reflections (0.002%)	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6952	wwPDB-VP
Average B, all atoms (Å ²)	32.0	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 40.17 % 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 2.8474e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² 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	0.63	0/2301	0.94	9/3125 (0.3%)
1	C	0.65	0/2293	0.94	12/3115 (0.4%)
2	B	0.68	1/856 (0.1%)	0.88	3/1160 (0.3%)
2	D	0.63	0/847	0.91	3/1148 (0.3%)
3	P	0.60	0/73	0.66	0/97
3	Q	0.66	0/73	0.71	0/97
All	All	0.64	1/6443 (0.0%)	0.92	27/8742 (0.3%)

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	C	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	59	ASP	CB-CG	-5.28	1.40	1.51

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	202	ARG	NE-CZ-NH1	11.09	125.84	120.30
1	A	202	ARG	NE-CZ-NH2	-10.35	115.13	120.30
1	A	35	ARG	NE-CZ-NH2	-10.01	115.30	120.30
1	C	202	ARG	NE-CZ-NH1	9.66	125.13	120.30
1	A	29	ASP	CB-CG-OD2	8.86	126.27	118.30
1	C	202	ARG	NE-CZ-NH2	-8.39	116.11	120.30
1	A	35	ARG	NE-CZ-NH1	8.20	124.40	120.30
1	C	111	ARG	NE-CZ-NH1	7.73	124.17	120.30
2	D	53	ASP	CB-CG-OD2	7.48	125.03	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	274	TRP	CA-C-N	7.17	132.98	117.20
1	A	119	ASP	CB-CG-OD2	7.11	124.69	118.30
1	C	29	ASP	CB-CG-OD2	6.45	124.10	118.30
1	A	183	ASP	CB-CG-OD2	6.42	124.08	118.30
1	C	129	ASP	CB-CG-OD2	6.27	123.94	118.30
2	D	98	ASP	C-N-CA	5.85	136.32	121.70
2	B	74[A]	GLU	CB-CA-C	5.79	121.98	110.40
2	B	74[B]	GLU	CB-CA-C	5.79	121.98	110.40
1	C	75	ARG	NE-CZ-NH2	5.68	123.14	120.30
1	C	272	LEU	CA-CB-CG	5.59	128.15	115.30
1	C	35	ARG	NE-CZ-NH2	5.58	123.09	120.30
1	C	227	ASP	CB-CG-OD2	5.39	123.15	118.30
2	D	85	ASP	CB-CG-OD2	5.36	123.12	118.30
2	B	59	ASP	CB-CA-C	-5.32	99.77	110.40
1	C	274	TRP	CB-CA-C	5.25	120.90	110.40
1	A	30	ASP	CB-CG-OD2	5.09	122.88	118.30
1	A	29	ASP	CB-CA-C	5.07	120.54	110.40
1	C	274	TRP	CA-CB-CG	5.02	123.23	113.70

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	273	ARG	Peptide

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	2240	0	2129	27	0
1	C	2232	0	2123	51	0
2	B	826	0	798	14	0
2	D	821	0	796	17	0
3	P	72	0	77	0	0
3	Q	72	0	77	0	0
4	A	234	0	0	16	0
4	B	115	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	C	226	0	0	10	0
4	D	104	0	0	3	0
4	P	6	0	0	0	0
4	Q	4	0	0	0	0
All	All	6952	0	6000	100	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 8.

All (100) 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:23:MET:HG3	4:A:2032:HOH:O	1.45	1.12
1:A:275:GLU:O	4:A:2234:HOH:O	1.67	1.10
1:C:202:ARG:HD2	1:C:244:TRP:CD2	2.09	0.88
1:C:234:ARG:HE	1:C:242:GLN:HE21	1.25	0.84
1:C:204:TRP:HH2	2:D:99:MET:HB2	1.45	0.82
1:C:275:GLU:N	4:C:2224:HOH:O	2.16	0.78
1:C:275:GLU:N	4:C:2226:HOH:O	2.15	0.78
1:A:23:MET:SD	4:A:2032:HOH:O	2.41	0.75
1:A:275:GLU:O	4:A:2231:HOH:O	2.04	0.75
1:A:202:ARG:HD2	1:A:244:TRP:CD2	2.21	0.75
1:C:28:VAL:O	1:C:31:THR:HG22	1.86	0.74
1:C:234:ARG:HE	1:C:242:GLN:NE2	1.88	0.71
1:A:23:MET:CG	4:A:2032:HOH:O	2.13	0.70
1:C:31:THR:HG21	1:C:179:LEU:CD1	2.21	0.70
1:C:152:GLU:OE1	1:C:155:ARG:NH1	2.26	0.69
1:A:54:GLN:HG2	4:A:2057:HOH:O	1.95	0.67
1:C:260:HIS:ND1	1:C:271:THR:HG22	2.11	0.65
2:D:96:ASP:O	2:D:99:MET:HA	1.96	0.64
1:C:274:TRP:HD1	4:C:2154:HOH:O	1.80	0.63
1:C:111:ARG:HB2	1:C:111:ARG:HH11	1.63	0.63
1:C:176:ASN:O	1:C:178:THR:O	2.17	0.63
2:D:77:THR:HG23	4:D:2078:HOH:O	1.98	0.62
2:B:74[B]:GLU:HG2	2:B:75:THR:HG23	1.80	0.62
1:C:220:ASN:HB2	4:C:2184:HOH:O	2.00	0.61
1:C:229:GLU:HG2	4:C:2189:HOH:O	2.01	0.61
1:C:115:GLN:HE22	2:D:58:LYS:NZ	2.00	0.60
2:B:32:PRO:O	2:B:84:HIS:HE1	1.84	0.60
1:C:111:ARG:CB	1:C:111:ARG:HH11	2.15	0.60
1:C:31:THR:HG21	1:C:179:LEU:HD12	1.85	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:144:LYS:NZ	1:A:148:GLU:OE2	2.36	0.59
2:B:48:LYS:HD3	2:B:48:LYS:H	1.67	0.59
1:A:275:GLU:CB	4:A:2234:HOH:O	2.51	0.58
2:D:77:THR:CG2	4:D:2078:HOH:O	2.51	0.57
1:C:145:HIS:O	1:C:149:GLN:HG2	2.05	0.57
2:B:48:LYS:CD	2:B:48:LYS:H	2.17	0.57
1:A:154:GLU:HG3	4:A:2137:HOH:O	2.05	0.57
1:C:144:LYS:NZ	1:C:148:GLU:OE2	2.36	0.56
1:C:178:THR:O	1:C:179:LEU:HB2	2.05	0.56
2:D:40:LEU:HD23	2:D:45:LYS:HA	1.87	0.55
2:D:17:ASN:ND2	2:D:74:GLU:HG3	2.22	0.55
1:C:44:ARG:CG	1:C:44:ARG:HH21	2.20	0.54
1:C:86:ASN:HB2	4:C:2082:HOH:O	2.07	0.54
2:D:10:TYR:HB2	2:D:99:MET:HE3	1.89	0.53
1:C:226:GLN:HB3	4:C:2187:HOH:O	2.09	0.53
1:C:202:ARG:CD	1:C:244:TRP:CE3	2.91	0.53
1:C:202:ARG:HD3	1:C:244:TRP:CE3	2.44	0.52
1:C:275:GLU:N	4:C:2225:HOH:O	2.43	0.52
2:B:7:ILE:HD12	2:B:91:LYS:HE3	1.91	0.52
1:C:273:ARG:O	1:C:274:TRP:HB3	2.10	0.52
2:B:84:HIS:HD2	2:B:86:SER:OG	1.93	0.52
1:C:44:ARG:HH21	1:C:44:ARG:HG2	1.74	0.52
1:A:31:THR:HG22	4:A:2029:HOH:O	2.09	0.52
1:C:260:HIS:CE1	1:C:271:THR:HG22	2.45	0.51
2:D:48:LYS:O	2:D:48:LYS:HD2	2.11	0.51
2:D:48:LYS:HG2	4:D:2073:HOH:O	2.11	0.51
1:C:159:TYR:HD2	1:C:160:LEU:HD13	1.75	0.51
1:C:111:ARG:HH11	1:C:111:ARG:CG	2.24	0.51
1:C:138:MET:HG3	4:C:2126:HOH:O	2.11	0.50
2:B:59:ASP:HB3	2:B:61:SER:H	1.76	0.50
1:C:117:ALA:HB2	2:D:60:TRP:CE2	2.47	0.50
1:A:117:ALA:HB2	2:B:60:TRP:CE2	2.46	0.50
1:A:48:ARG:NH2	2:B:53:ASP:OD2	2.44	0.49
1:C:236:ALA:O	2:D:12:ARG:HD3	2.10	0.49
1:A:187:ALA:HA	1:A:204:TRP:O	2.11	0.49
1:C:202:ARG:CD	1:C:244:TRP:CD2	2.89	0.49
1:C:111:ARG:NH1	1:C:111:ARG:HB2	2.28	0.49
1:C:204:TRP:CH2	2:D:99:MET:HB2	2.35	0.48
1:C:185:PRO:HG2	1:C:266:LEU:HD11	1.96	0.48
2:D:48:LYS:O	2:D:48:LYS:CD	2.62	0.47
1:A:41:GLU:OE1	1:A:42:ASN:N	2.41	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:75:ARG:HD3	4:A:2081:HOH:O	2.14	0.46
1:A:23:MET:CE	4:A:2032:HOH:O	2.61	0.46
2:B:21:ASN:C	2:B:22:ILE:HD12	2.35	0.46
1:A:227:ASP:HB2	4:A:2197:HOH:O	2.15	0.46
1:C:187:ALA:HA	1:C:204:TRP:O	2.16	0.46
1:C:274:TRP:HZ2	4:C:2223:HOH:O	1.99	0.46
1:A:73:SER:HB3	4:A:2078:HOH:O	2.15	0.45
1:C:236:ALA:O	2:D:12:ARG:CD	2.65	0.45
2:B:40:LEU:HD23	2:B:45:LYS:HA	1.98	0.45
1:A:275:GLU:HB3	4:A:2234:HOH:O	2.14	0.44
1:C:81:LEU:HD13	1:C:118:TYR:CD1	2.53	0.44
2:B:48:LYS:HG2	4:B:2066:HOH:O	2.17	0.44
1:C:202:ARG:HD2	1:C:244:TRP:CE3	2.50	0.43
2:D:73:THR:HG22	2:D:75:THR:H	1.82	0.43
2:D:73:THR:HG22	2:D:76:ASP:H	1.82	0.43
1:C:31:THR:HG21	1:C:179:LEU:HD11	2.00	0.43
1:A:54:GLN:NE2	4:A:2056:HOH:O	2.47	0.43
1:C:225:ILE:C	1:C:227:ASP:H	2.21	0.43
1:A:121:CYS:SG	2:B:1:ILE:HG22	2.58	0.43
1:A:29:ASP:O	4:A:2024:HOH:O	2.21	0.43
1:A:202:ARG:CD	1:A:244:TRP:CE3	3.02	0.42
1:C:196:GLU:HB3	1:C:197:ASP:H	1.72	0.42
1:A:202:ARG:HD2	1:A:244:TRP:CE3	2.54	0.41
1:C:219:LEU:O	1:C:221:GLY:N	2.54	0.41
1:C:202:ARG:HD2	1:C:244:TRP:CG	2.53	0.41
2:B:22:ILE:N	2:B:22:ILE:HD12	2.36	0.41
1:A:172:LEU:O	1:A:176:ASN:HB2	2.20	0.41
1:A:32:GLU:OE2	1:A:35:ARG:HD2	2.21	0.41
1:C:44:ARG:NH1	1:C:61:GLU:OE1	2.53	0.41
1:C:170:ARG:HH11	1:C:170:ARG:HG3	1.86	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	273/275 (99%)	261 (96%)	11 (4%)	1 (0%)	39	33
1	C	273/275 (99%)	260 (95%)	7 (3%)	6 (2%)	8	3
2	B	98/99 (99%)	95 (97%)	3 (3%)	0	100	100
2	D	97/99 (98%)	95 (98%)	2 (2%)	0	100	100
3	P	7/9 (78%)	7 (100%)	0	0	100	100
3	Q	7/9 (78%)	7 (100%)	0	0	100	100
All	All	755/766 (99%)	725 (96%)	23 (3%)	7 (1%)	21	13

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	222	GLU
1	C	225	ILE
1	C	274	TRP
1	C	177	ALA
1	C	178	THR
1	C	220	ASN
1	C	224	LEU

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	233/233 (100%)	216 (93%)	17 (7%)	17	11
1	C	232/233 (100%)	208 (90%)	24 (10%)	9	5
2	B	95/94 (101%)	90 (95%)	5 (5%)	28	22
2	D	94/94 (100%)	85 (90%)	9 (10%)	10	6
3	P	8/8 (100%)	8 (100%)	0	100	100
3	Q	8/8 (100%)	8 (100%)	0	100	100
All	All	670/670 (100%)	615 (92%)	55 (8%)	14	9

All (55) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	41	GLU
1	A	44	ARG
1	A	45	TYR
1	A	65	GLN
1	A	82	LEU
1	A	102	GLU
1	A	111	ARG
1	A	113	TYR
1	A	156	LEU
1	A	160	LEU
1	A	173	LYS
1	A	181	ARG
1	A	182	THR
1	A	196	GLU
1	A	202	ARG
1	A	230	LEU
1	A	273	ARG
2	B	1	ILE
2	B	44	LYS
2	B	48	LYS
2	B	59	ASP
2	B	70	PHE
1	C	5	LEU
1	C	17	LEU
1	C	30	ASP
1	C	42	ASN
1	C	44	ARG
1	C	45	TYR
1	C	50	ARG
1	C	75	ARG
1	C	89	LYS
1	C	103	VAL
1	C	110	LEU
1	C	111	ARG
1	C	114	GLN
1	C	141	LEU
1	C	160	LEU
1	C	178	THR
1	C	179	LEU
1	C	180	LEU
1	C	196	GLU
1	C	225	ILE

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Mol	Chain	Res	Type
1	C	226	GLN
1	C	230	LEU
1	C	272	LEU
1	C	273	ARG
2	D	12	ARG
2	D	44	LYS
2	D	48	LYS
2	D	58	LYS
2	D	61	SER
2	D	70	PHE
2	D	73	THR
2	D	77	THR
2	D	99	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	65	GLN
1	A	87	GLN
1	A	114	GLN
1	A	145	HIS
1	A	218	GLN
1	A	264	GLN
2	B	29	GLN
2	B	84	HIS
1	C	42	ASN
1	C	87	GLN
1	C	114	GLN
1	C	115	GLN
1	C	218	GLN
1	C	226	GLN
1	C	242	GLN
2	D	2	GLN
2	D	8	GLN

5.3.3 RNA ⓘ

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, 95th 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 ²)	Q<0.9
1	A	275/275 (100%)	0.11	10 (3%) 46 48	18, 31, 44, 58	0
1	C	275/275 (100%)	0.27	21 (7%) 17 18	18, 32, 49, 57	0
2	B	99/99 (100%)	0.01	0 100 100	17, 27, 38, 46	0
2	D	99/99 (100%)	0.19	4 (4%) 42 44	21, 29, 40, 52	0
3	P	9/9 (100%)	-0.16	0 100 100	27, 31, 33, 33	0
3	Q	9/9 (100%)	-0.02	0 100 100	24, 31, 34, 37	0
All	All	766/766 (100%)	0.16	35 (4%) 36 38	17, 31, 44, 58	0

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	1	ILE	6.9
1	C	226	GLN	6.9
1	A	226	GLN	6.5
1	C	177	ALA	5.0
1	C	41	GLU	4.9
2	D	99	MET	4.6
1	A	225	ILE	4.5
1	C	42	ASN	4.4
1	C	176	ASN	4.2
1	C	16	GLY	3.9
1	C	225	ILE	3.9
1	C	227	ASP	3.5
1	A	275	GLU	3.5
1	C	180	LEU	3.3
1	A	41	GLU	3.0
1	A	222	GLU	2.9
1	C	221	GLY	2.8
1	C	86	ASN	2.8
1	C	178	THR	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	111	ARG	2.6
1	A	193	SER	2.6
1	C	40	ALA	2.5
1	A	195	PRO	2.5
1	C	50	ARG	2.4
1	A	42	ASN	2.4
1	C	89	LYS	2.4
1	C	220	ASN	2.3
1	C	195	PRO	2.3
1	C	111	ARG	2.1
1	C	5	LEU	2.1
1	A	170	ARG	2.1
1	C	222	GLU	2.1
1	C	168	LEU	2.1
2	D	48	LYS	2.0
2	D	88	ALA	2.0

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.