



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 17, 2017 – 03:02 PM EDT

PDB ID : 3WEX
Title : Crystal structure of HLA-DP5 in complex with Cry j 1-derived peptide (residues 214-222)
Authors : Kusano, S.; Kukimoto-Niino, M.; Shirouzu, M.; Yokoyama, S.
Deposited on : 2013-07-16
Resolution : 2.40 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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20029077
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
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) : rb-20029077

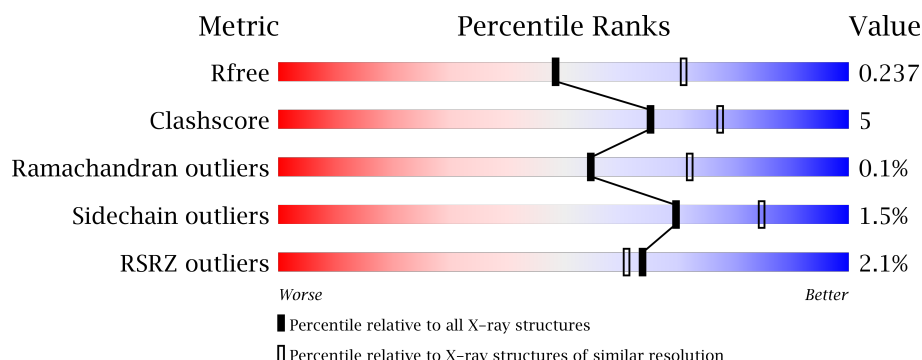
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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}	100719	3166 (2.40-2.40)
Clashscore	112137	3674 (2.40-2.40)
Ramachandran outliers	110173	3616 (2.40-2.40)
Sidechain outliers	110143	3617 (2.40-2.40)
RSRZ outliers	101464	3195 (2.40-2.40)

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	193	<div> <div>86%</div> <div>7% • 6%</div> </div>
1	C	193	<div> <div>82%</div> <div>10% •• 6%</div> </div>
1	E	193	<div> <div>80%</div> <div>13% • 6%</div> </div>
1	G	193	<div> <div>88%</div> <div>5% • 6%</div> </div>
2	B	221	<div> <div>76%</div> <div>12% 12%</div> </div>

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Mol	Chain	Length	Quality of chain
2	D	221	
2	F	221	
2	H	221	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	NAG	C	201	-	-	-	X
3	NAG	G	201	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 12695 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 MHC class II antigen.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	181	Total	C	N	O	S	0	0	0
			1483	955	245	278	5			
1	C	181	Total	C	N	O	S	0	0	0
			1483	955	245	278	5			
1	E	181	Total	C	N	O	S	0	0	0
			1483	955	245	278	5			
1	G	181	Total	C	N	O	S	0	0	0
			1483	955	245	278	5			

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	182	SER	-	EXPRESSION TAG	UNP I2G9G1
A	183	GLY	-	EXPRESSION TAG	UNP I2G9G1
A	184	PRO	-	EXPRESSION TAG	UNP I2G9G1
A	185	SER	-	EXPRESSION TAG	UNP I2G9G1
A	186	SER	-	EXPRESSION TAG	UNP I2G9G1
A	187	GLY	-	EXPRESSION TAG	UNP I2G9G1
A	188	GLU	-	EXPRESSION TAG	UNP I2G9G1
A	189	ASN	-	EXPRESSION TAG	UNP I2G9G1
A	190	LEU	-	EXPRESSION TAG	UNP I2G9G1
A	191	TYR	-	EXPRESSION TAG	UNP I2G9G1
A	192	PHE	-	EXPRESSION TAG	UNP I2G9G1
A	193	GLN	-	EXPRESSION TAG	UNP I2G9G1
C	182	SER	-	EXPRESSION TAG	UNP I2G9G1
C	183	GLY	-	EXPRESSION TAG	UNP I2G9G1
C	184	PRO	-	EXPRESSION TAG	UNP I2G9G1
C	185	SER	-	EXPRESSION TAG	UNP I2G9G1
C	186	SER	-	EXPRESSION TAG	UNP I2G9G1
C	187	GLY	-	EXPRESSION TAG	UNP I2G9G1
C	188	GLU	-	EXPRESSION TAG	UNP I2G9G1
C	189	ASN	-	EXPRESSION TAG	UNP I2G9G1
C	190	LEU	-	EXPRESSION TAG	UNP I2G9G1

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Chain	Residue	Modelled	Actual	Comment	Reference
C	191	TYR	-	EXPRESSION TAG	UNP I2G9G1
C	192	PHE	-	EXPRESSION TAG	UNP I2G9G1
C	193	GLN	-	EXPRESSION TAG	UNP I2G9G1
E	182	SER	-	EXPRESSION TAG	UNP I2G9G1
E	183	GLY	-	EXPRESSION TAG	UNP I2G9G1
E	184	PRO	-	EXPRESSION TAG	UNP I2G9G1
E	185	SER	-	EXPRESSION TAG	UNP I2G9G1
E	186	SER	-	EXPRESSION TAG	UNP I2G9G1
E	187	GLY	-	EXPRESSION TAG	UNP I2G9G1
E	188	GLU	-	EXPRESSION TAG	UNP I2G9G1
E	189	ASN	-	EXPRESSION TAG	UNP I2G9G1
E	190	LEU	-	EXPRESSION TAG	UNP I2G9G1
E	191	TYR	-	EXPRESSION TAG	UNP I2G9G1
E	192	PHE	-	EXPRESSION TAG	UNP I2G9G1
E	193	GLN	-	EXPRESSION TAG	UNP I2G9G1
G	182	SER	-	EXPRESSION TAG	UNP I2G9G1
G	183	GLY	-	EXPRESSION TAG	UNP I2G9G1
G	184	PRO	-	EXPRESSION TAG	UNP I2G9G1
G	185	SER	-	EXPRESSION TAG	UNP I2G9G1
G	186	SER	-	EXPRESSION TAG	UNP I2G9G1
G	187	GLY	-	EXPRESSION TAG	UNP I2G9G1
G	188	GLU	-	EXPRESSION TAG	UNP I2G9G1
G	189	ASN	-	EXPRESSION TAG	UNP I2G9G1
G	190	LEU	-	EXPRESSION TAG	UNP I2G9G1
G	191	TYR	-	EXPRESSION TAG	UNP I2G9G1
G	192	PHE	-	EXPRESSION TAG	UNP I2G9G1
G	193	GLN	-	EXPRESSION TAG	UNP I2G9G1

- Molecule 2 is a protein called MHC class II antigen.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	195	Total	C	N	O	S	0	0	0
			1610	1017	284	302	7			
2	D	195	Total	C	N	O	S	0	0	0
			1610	1017	284	302	7			
2	F	195	Total	C	N	O	S	0	0	0
			1610	1017	284	302	7			
2	H	195	Total	C	N	O	S	0	0	0
			1610	1017	284	302	7			

There are 148 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-19	LYS	-	EXPRESSION TAG	UNP I2FL84
B	-18	VAL	-	EXPRESSION TAG	UNP I2FL84
B	-17	THR	-	EXPRESSION TAG	UNP I2FL84
B	-16	VAL	-	EXPRESSION TAG	UNP I2FL84
B	-15	ALA	-	EXPRESSION TAG	UNP I2FL84
B	-14	PHE	-	EXPRESSION TAG	UNP I2FL84
B	-13	ASN	-	EXPRESSION TAG	UNP I2FL84
B	-12	GLN	-	EXPRESSION TAG	UNP I2FL84
B	-11	PHE	-	EXPRESSION TAG	UNP I2FL84
B	-10	GLY	-	EXPRESSION TAG	UNP I2FL84
B	-9	GLY	-	EXPRESSION TAG	UNP I2FL84
B	-8	SER	-	EXPRESSION TAG	UNP I2FL84
B	-7	LEU	-	EXPRESSION TAG	UNP I2FL84
B	-6	VAL	-	EXPRESSION TAG	UNP I2FL84
B	-5	PRO	-	EXPRESSION TAG	UNP I2FL84
B	-4	ARG	-	EXPRESSION TAG	UNP I2FL84
B	-3	GLY	-	EXPRESSION TAG	UNP I2FL84
B	-2	SER	-	EXPRESSION TAG	UNP I2FL84
B	-1	GLY	-	EXPRESSION TAG	UNP I2FL84
B	0	GLY	-	EXPRESSION TAG	UNP I2FL84
B	1	GLY	-	EXPRESSION TAG	UNP I2FL84
B	2	GLY	-	EXPRESSION TAG	UNP I2FL84
B	3	SER	-	EXPRESSION TAG	UNP I2FL84
B	4	PRO	-	EXPRESSION TAG	UNP I2FL84
B	5	GLU	-	EXPRESSION TAG	UNP I2FL84
B	190	SER	-	EXPRESSION TAG	UNP I2FL84
B	191	GLY	-	EXPRESSION TAG	UNP I2FL84
B	192	PRO	-	EXPRESSION TAG	UNP I2FL84
B	193	SER	-	EXPRESSION TAG	UNP I2FL84
B	194	SER	-	EXPRESSION TAG	UNP I2FL84
B	195	GLY	-	EXPRESSION TAG	UNP I2FL84
B	196	GLU	-	EXPRESSION TAG	UNP I2FL84
B	197	ASN	-	EXPRESSION TAG	UNP I2FL84
B	198	LEU	-	EXPRESSION TAG	UNP I2FL84
B	199	TYR	-	EXPRESSION TAG	UNP I2FL84
B	200	PHE	-	EXPRESSION TAG	UNP I2FL84
B	201	GLN	-	EXPRESSION TAG	UNP I2FL84
D	-19	LYS	-	EXPRESSION TAG	UNP I2FL84
D	-18	VAL	-	EXPRESSION TAG	UNP I2FL84
D	-17	THR	-	EXPRESSION TAG	UNP I2FL84
D	-16	VAL	-	EXPRESSION TAG	UNP I2FL84
D	-15	ALA	-	EXPRESSION TAG	UNP I2FL84
D	-14	PHE	-	EXPRESSION TAG	UNP I2FL84

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-13	ASN	-	EXPRESSION TAG	UNP I2FL84
D	-12	GLN	-	EXPRESSION TAG	UNP I2FL84
D	-11	PHE	-	EXPRESSION TAG	UNP I2FL84
D	-10	GLY	-	EXPRESSION TAG	UNP I2FL84
D	-9	GLY	-	EXPRESSION TAG	UNP I2FL84
D	-8	SER	-	EXPRESSION TAG	UNP I2FL84
D	-7	LEU	-	EXPRESSION TAG	UNP I2FL84
D	-6	VAL	-	EXPRESSION TAG	UNP I2FL84
D	-5	PRO	-	EXPRESSION TAG	UNP I2FL84
D	-4	ARG	-	EXPRESSION TAG	UNP I2FL84
D	-3	GLY	-	EXPRESSION TAG	UNP I2FL84
D	-2	SER	-	EXPRESSION TAG	UNP I2FL84
D	-1	GLY	-	EXPRESSION TAG	UNP I2FL84
D	0	GLY	-	EXPRESSION TAG	UNP I2FL84
D	1	GLY	-	EXPRESSION TAG	UNP I2FL84
D	2	GLY	-	EXPRESSION TAG	UNP I2FL84
D	3	SER	-	EXPRESSION TAG	UNP I2FL84
D	4	PRO	-	EXPRESSION TAG	UNP I2FL84
D	5	GLU	-	EXPRESSION TAG	UNP I2FL84
D	190	SER	-	EXPRESSION TAG	UNP I2FL84
D	191	GLY	-	EXPRESSION TAG	UNP I2FL84
D	192	PRO	-	EXPRESSION TAG	UNP I2FL84
D	193	SER	-	EXPRESSION TAG	UNP I2FL84
D	194	SER	-	EXPRESSION TAG	UNP I2FL84
D	195	GLY	-	EXPRESSION TAG	UNP I2FL84
D	196	GLU	-	EXPRESSION TAG	UNP I2FL84
D	197	ASN	-	EXPRESSION TAG	UNP I2FL84
D	198	LEU	-	EXPRESSION TAG	UNP I2FL84
D	199	TYR	-	EXPRESSION TAG	UNP I2FL84
D	200	PHE	-	EXPRESSION TAG	UNP I2FL84
D	201	GLN	-	EXPRESSION TAG	UNP I2FL84
F	-19	LYS	-	EXPRESSION TAG	UNP I2FL84
F	-18	VAL	-	EXPRESSION TAG	UNP I2FL84
F	-17	THR	-	EXPRESSION TAG	UNP I2FL84
F	-16	VAL	-	EXPRESSION TAG	UNP I2FL84
F	-15	ALA	-	EXPRESSION TAG	UNP I2FL84
F	-14	PHE	-	EXPRESSION TAG	UNP I2FL84
F	-13	ASN	-	EXPRESSION TAG	UNP I2FL84
F	-12	GLN	-	EXPRESSION TAG	UNP I2FL84
F	-11	PHE	-	EXPRESSION TAG	UNP I2FL84
F	-10	GLY	-	EXPRESSION TAG	UNP I2FL84
F	-9	GLY	-	EXPRESSION TAG	UNP I2FL84

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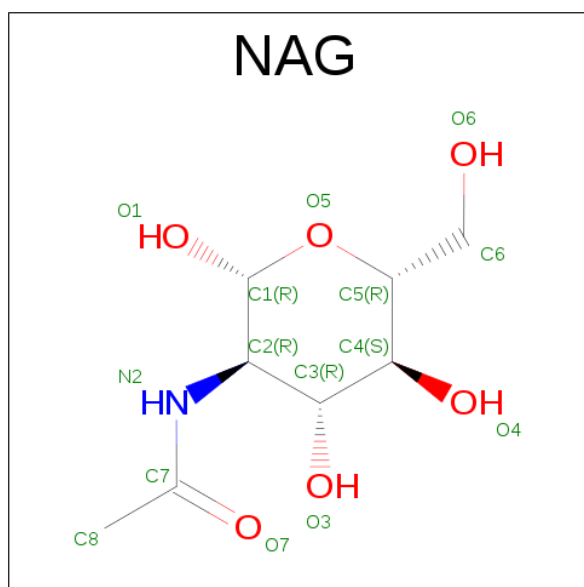
Chain	Residue	Modelled	Actual	Comment	Reference
F	-8	SER	-	EXPRESSION TAG	UNP I2FL84
F	-7	LEU	-	EXPRESSION TAG	UNP I2FL84
F	-6	VAL	-	EXPRESSION TAG	UNP I2FL84
F	-5	PRO	-	EXPRESSION TAG	UNP I2FL84
F	-4	ARG	-	EXPRESSION TAG	UNP I2FL84
F	-3	GLY	-	EXPRESSION TAG	UNP I2FL84
F	-2	SER	-	EXPRESSION TAG	UNP I2FL84
F	-1	GLY	-	EXPRESSION TAG	UNP I2FL84
F	0	GLY	-	EXPRESSION TAG	UNP I2FL84
F	1	GLY	-	EXPRESSION TAG	UNP I2FL84
F	2	GLY	-	EXPRESSION TAG	UNP I2FL84
F	3	SER	-	EXPRESSION TAG	UNP I2FL84
F	4	PRO	-	EXPRESSION TAG	UNP I2FL84
F	5	GLU	-	EXPRESSION TAG	UNP I2FL84
F	190	SER	-	EXPRESSION TAG	UNP I2FL84
F	191	GLY	-	EXPRESSION TAG	UNP I2FL84
F	192	PRO	-	EXPRESSION TAG	UNP I2FL84
F	193	SER	-	EXPRESSION TAG	UNP I2FL84
F	194	SER	-	EXPRESSION TAG	UNP I2FL84
F	195	GLY	-	EXPRESSION TAG	UNP I2FL84
F	196	GLU	-	EXPRESSION TAG	UNP I2FL84
F	197	ASN	-	EXPRESSION TAG	UNP I2FL84
F	198	LEU	-	EXPRESSION TAG	UNP I2FL84
F	199	TYR	-	EXPRESSION TAG	UNP I2FL84
F	200	PHE	-	EXPRESSION TAG	UNP I2FL84
F	201	GLN	-	EXPRESSION TAG	UNP I2FL84
H	-19	LYS	-	EXPRESSION TAG	UNP I2FL84
H	-18	VAL	-	EXPRESSION TAG	UNP I2FL84
H	-17	THR	-	EXPRESSION TAG	UNP I2FL84
H	-16	VAL	-	EXPRESSION TAG	UNP I2FL84
H	-15	ALA	-	EXPRESSION TAG	UNP I2FL84
H	-14	PHE	-	EXPRESSION TAG	UNP I2FL84
H	-13	ASN	-	EXPRESSION TAG	UNP I2FL84
H	-12	GLN	-	EXPRESSION TAG	UNP I2FL84
H	-11	PHE	-	EXPRESSION TAG	UNP I2FL84
H	-10	GLY	-	EXPRESSION TAG	UNP I2FL84
H	-9	GLY	-	EXPRESSION TAG	UNP I2FL84
H	-8	SER	-	EXPRESSION TAG	UNP I2FL84
H	-7	LEU	-	EXPRESSION TAG	UNP I2FL84
H	-6	VAL	-	EXPRESSION TAG	UNP I2FL84
H	-5	PRO	-	EXPRESSION TAG	UNP I2FL84
H	-4	ARG	-	EXPRESSION TAG	UNP I2FL84

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Chain	Residue	Modelled	Actual	Comment	Reference
H	-3	GLY	-	EXPRESSION TAG	UNP I2FL84
H	-2	SER	-	EXPRESSION TAG	UNP I2FL84
H	-1	GLY	-	EXPRESSION TAG	UNP I2FL84
H	0	GLY	-	EXPRESSION TAG	UNP I2FL84
H	1	GLY	-	EXPRESSION TAG	UNP I2FL84
H	2	GLY	-	EXPRESSION TAG	UNP I2FL84
H	3	SER	-	EXPRESSION TAG	UNP I2FL84
H	4	PRO	-	EXPRESSION TAG	UNP I2FL84
H	5	GLU	-	EXPRESSION TAG	UNP I2FL84
H	190	SER	-	EXPRESSION TAG	UNP I2FL84
H	191	GLY	-	EXPRESSION TAG	UNP I2FL84
H	192	PRO	-	EXPRESSION TAG	UNP I2FL84
H	193	SER	-	EXPRESSION TAG	UNP I2FL84
H	194	SER	-	EXPRESSION TAG	UNP I2FL84
H	195	GLY	-	EXPRESSION TAG	UNP I2FL84
H	196	GLU	-	EXPRESSION TAG	UNP I2FL84
H	197	ASN	-	EXPRESSION TAG	UNP I2FL84
H	198	LEU	-	EXPRESSION TAG	UNP I2FL84
H	199	TYR	-	EXPRESSION TAG	UNP I2FL84
H	200	PHE	-	EXPRESSION TAG	UNP I2FL84
H	201	GLN	-	EXPRESSION TAG	UNP I2FL84

- Molecule 3 is N-ACETYL-D-GLUCOSAMINE (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			14	8	1	5		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	C	1	Total	C	N	O	0	0
			14	8	1	5		
3	E	1	Total	C	N	O	0	0
			14	8	1	5		
3	G	1	Total	C	N	O	0	0
			14	8	1	5		


- Molecule 4 is water.

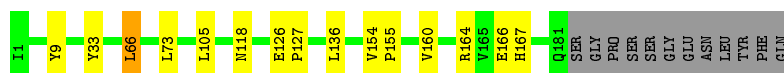
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	34	Total	O	0	0
			34	34		
4	B	32	Total	O	0	0
			32	32		
4	C	35	Total	O	0	0
			35	35		
4	D	37	Total	O	0	0
			37	37		
4	E	32	Total	O	0	0
			32	32		
4	F	23	Total	O	0	0
			23	23		
4	G	39	Total	O	0	0
			39	39		
4	H	35	Total	O	0	0
			35	35		

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 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: MHC class II antigen

Chain A: 




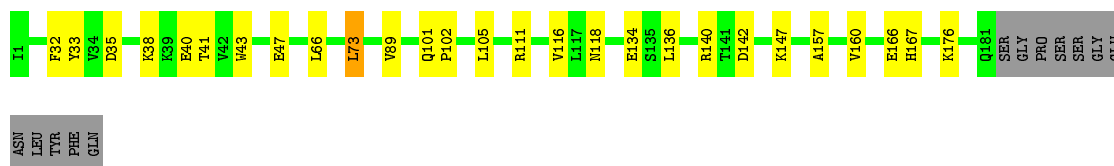
- Molecule 1: MHC class II antigen

Chain C: 




- Molecule 1: MHC class II antigen

Chain E: 




- Molecule 1: MHC class II antigen

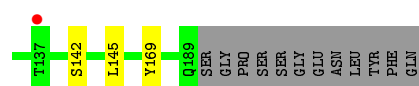
Chain G: 



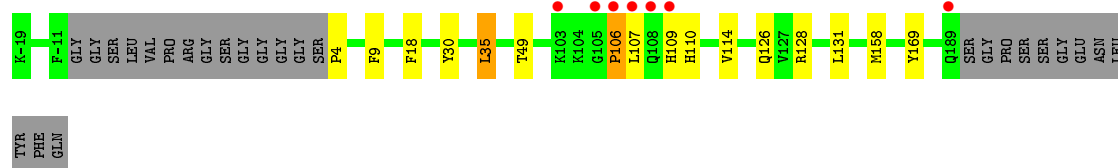
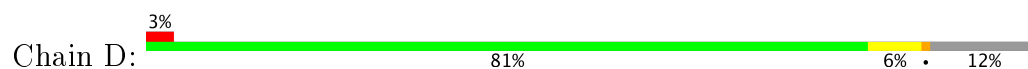
- Molecule 2: MHC class II antigen

Chain B: 

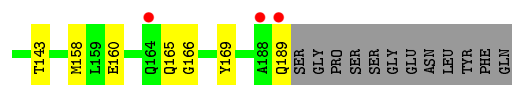
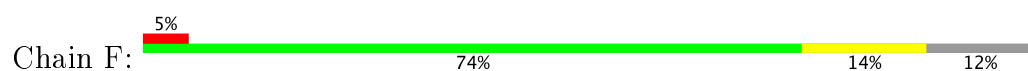




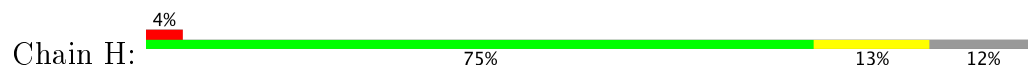
- Molecule 2: MHC class II antigen



- Molecule 2: MHC class II antigen



- Molecule 2: MHC class II antigen



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	61.15Å 64.37Å 130.36Å 92.99° 97.53° 109.41°	Depositor
Resolution (Å)	64.28 – 2.40 64.28 – 2.40	Depositor EDS
% Data completeness (in resolution range)	97.3 (64.28-2.40) 94.0 (64.28-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.20	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.88 (at 2.40Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.1_1168)	Depositor
R, R_{free}	0.202 , 0.237 0.200 , 0.237	Depositor DCC
R_{free} test set	3519 reflections (5.00%)	DCC
Wilson B-factor (Å ²)	28.0	Xtriage
Anisotropy	0.020	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 45.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	12695	wwPDB-VP
Average B, all atoms (Å ²)	29.0	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 38.06 % 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 3.9040e-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, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NAG

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	0.32	0/1531	0.51	0/2090
1	C	0.32	0/1531	0.52	1/2090 (0.0%)
1	E	0.33	0/1531	0.50	0/2090
1	G	0.31	0/1531	0.49	0/2090
2	B	0.33	0/1648	0.55	0/2234
2	D	0.31	0/1648	0.52	0/2234
2	F	0.33	0/1648	0.59	2/2234 (0.1%)
2	H	0.32	0/1648	0.53	0/2234
All	All	0.32	0/12716	0.53	3/17296 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	73	LEU	CA-CB-CG	-5.72	102.14	115.30
2	F	143	THR	N-CA-C	-5.10	97.23	111.00
2	F	107	LEU	CA-CB-CG	5.03	126.88	115.30

There are no chirality outliers.

There are no planarity outliers.

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	1483	0	1386	10	0
1	C	1483	0	1386	28	0
1	E	1483	0	1386	21	0
1	G	1483	0	1386	10	0
2	B	1610	0	1546	18	0
2	D	1610	0	1546	16	1
2	F	1610	0	1546	16	1
2	H	1610	0	1546	21	0
3	A	14	0	13	1	0
3	C	14	0	13	0	0
3	E	14	0	13	2	0
3	G	14	0	13	0	0
4	A	34	0	0	0	0
4	B	32	0	0	0	0
4	C	35	0	0	1	0
4	D	37	0	0	0	0
4	E	32	0	0	0	0
4	F	23	0	0	0	0
4	G	39	0	0	0	0
4	H	35	0	0	0	0
All	All	12695	0	11780	113	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 5.

All (113) 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:H:44:GLU:OE2	2:H:46:ARG:NH1	1.96	0.98
2:F:74:ASP:OD1	2:F:78:ARG:NH1	2.02	0.93
2:B:128:ARG:HG3	2:B:128:ARG:HH11	1.33	0.92
2:B:50:GLU:OE1	2:B:53:ARG:NH2	2.04	0.90
2:F:19:ASN:OD1	2:F:20:GLY:N	2.04	0.90
2:D:126:GLN:NE2	2:D:128:ARG:HH12	1.74	0.85
2:B:8:LEU:HD21	2:B:10:GLN:HG3	1.61	0.82
2:F:8:LEU:HD21	2:F:10:GLN:HG3	1.61	0.81
1:C:164:ARG:NH1	1:C:166:GLU:OE2	2.17	0.77
2:F:53:ARG:NE	2:F:57:GLU:OE2	2.17	0.77
2:F:114:VAL:HG22	2:F:158:MET:HG2	1.70	0.72
1:A:118:ASN:HB2	1:A:166:GLU:HB2	1.73	0.70
2:B:128:ARG:HG3	2:B:128:ARG:NH1	2.08	0.69
2:B:33:GLU:OE2	2:B:49:THR:HG21	1.92	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:111:ARG:HD3	1:E:111:ARG:NH1	2.09	0.67
2:D:126:GLN:CD	2:D:128:ARG:HH12	1.98	0.66
1:A:164:ARG:NH1	1:A:166:GLU:OE2	2.26	0.66
1:C:47:GLU:N	1:C:47:GLU:OE1	2.24	0.66
2:H:114:VAL:HG22	2:H:158:MET:HG2	1.77	0.65
2:H:74:ASP:O	2:H:78:ARG:NH1	2.29	0.65
1:C:16:HIS:CE1	2:D:4:PRO:HD2	2.33	0.64
2:D:114:VAL:HG22	2:D:158:MET:HG2	1.78	0.64
1:E:47:GLU:OE1	1:E:47:GLU:N	2.27	0.64
1:E:73:LEU:HD13	2:F:9:PHE:HE1	1.64	0.62
2:F:189:GLN:N	2:F:189:GLN:OE1	2.32	0.62
1:C:46:GLU:HG2	1:C:50:ARG:CZ	2.30	0.61
1:C:39:LYS:NZ	1:C:60:LEU:HD22	2.17	0.59
1:C:111:ARG:NH1	1:E:111:ARG:HH11	2.01	0.58
1:E:167:HIS:HA	3:E:201:NAG:H81	1.84	0.58
1:C:111:ARG:HH11	1:E:111:ARG:HH11	1.52	0.57
2:B:126:GLN:OE1	2:B:128:ARG:NH1	2.38	0.57
2:D:110:HIS:HE1	1:E:157:ALA:O	1.89	0.56
1:C:66:LEU:HD13	2:D:9:PHE:HD2	1.71	0.56
2:H:128:ARG:NH1	2:H:174:GLU:OE2	2.40	0.55
2:H:131:LEU:HD13	2:H:169:TYR:CE2	2.42	0.55
1:A:167:HIS:HA	3:A:201:NAG:H81	1.89	0.54
2:B:71:ALA:HB1	2:B:75:ARG:NH1	2.23	0.54
2:H:108:GLN:O	2:H:108:GLN:HG3	2.08	0.54
1:C:16:HIS:HE1	2:D:4:PRO:HD2	1.71	0.54
1:A:66:LEU:HD13	2:B:9:PHE:HD2	1.74	0.53
1:C:111:ARG:HH11	1:E:111:ARG:NH1	2.07	0.53
2:H:74:ASP:HB3	2:H:78:ARG:HH12	1.73	0.53
1:E:118:ASN:HB2	1:E:166:GLU:HB2	1.90	0.53
2:D:106:PRO:HG2	2:D:109:HIS:CD2	2.44	0.53
1:C:162:ASP:OD2	4:C:335:HOH:O	2.19	0.52
1:G:66:LEU:HD13	2:H:9:PHE:HD2	1.76	0.51
1:C:118:ASN:HB2	1:C:166:GLU:HB2	1.91	0.51
1:G:66:LEU:HD13	2:H:9:PHE:CD2	2.46	0.50
1:C:73:LEU:HD22	2:D:35:LEU:HD21	1.94	0.49
1:A:66:LEU:HD13	2:B:9:PHE:CD2	2.49	0.48
2:D:131:LEU:HD13	2:D:169:TYR:CE2	2.49	0.48
2:B:71:ALA:HB1	2:B:75:ARG:HH11	1.77	0.48
1:C:134:GLU:OE1	1:C:147:LYS:NZ	2.46	0.48
1:E:73:LEU:HD22	2:F:30:TYR:HB2	1.96	0.47
1:C:46:GLU:HG2	1:C:50:ARG:NH2	2.30	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:88:GLU:OE2	1:E:111:ARG:NE	2.48	0.47
2:B:64:ASP:O	2:B:68:GLU:HG3	2.14	0.47
2:F:131:LEU:HD13	2:F:169:TYR:CE1	2.50	0.47
1:G:105:LEU:HD23	1:G:105:LEU:HA	1.74	0.47
2:D:126:GLN:CD	2:D:128:ARG:NH1	2.65	0.47
1:G:73:LEU:HD22	2:H:35:LEU:HD21	1.96	0.47
1:C:177:HIS:CD2	2:F:112:LEU:HD21	2.49	0.46
1:G:118:ASN:HB2	1:G:166:GLU:HB2	1.97	0.45
1:A:33:TYR:CG	1:A:136:LEU:HD11	2.51	0.45
2:F:165:GLN:CG	2:F:166:GLY:HA2	2.47	0.45
1:A:9:TYR:HB2	2:B:13:GLN:HB2	1.97	0.45
2:B:4:PRO:HB2	2:B:5:GLU:H	1.57	0.45
1:G:85:ASP:OD1	2:H:32:ARG:NH1	2.48	0.45
1:G:76:ARG:NH1	2:H:54:PRO:HG2	2.31	0.45
1:E:33:TYR:CG	1:E:136:LEU:HD11	2.51	0.45
1:C:98:GLU:HB3	2:H:61:SER:HB2	1.99	0.44
2:H:106:PRO:C	2:H:107:LEU:HD12	2.38	0.44
2:H:131:LEU:HD13	2:H:169:TYR:CZ	2.53	0.44
1:C:39:LYS:HZ2	1:C:60:LEU:HD22	1.80	0.44
1:C:66:LEU:HD13	2:D:9:PHE:CD2	2.50	0.44
1:C:2:LYS:HD2	1:C:3:ALA:H	1.83	0.44
1:E:116:VAL:HG12	3:E:201:NAG:H83	2.00	0.44
1:E:40:GLU:CD	1:E:41:THR:H	2.21	0.44
2:B:96:LYS:N	2:B:96:LYS:HD2	2.32	0.43
1:G:176:LYS:HA	1:G:176:LYS:HD3	1.83	0.43
1:G:160:VAL:HG13	1:G:177:HIS:CE1	2.53	0.43
2:H:161:MET:SD	2:H:163:PRO:HG3	2.58	0.43
1:A:105:LEU:HD23	1:A:105:LEU:HA	1.84	0.43
1:C:3:ALA:HA	2:D:18:PHE:HB2	2.01	0.43
1:E:35:ASP:OD2	1:E:38:LYS:HD2	2.18	0.43
2:H:115:CYS:HB2	2:H:129:TRP:CZ2	2.53	0.43
2:B:131:LEU:HD13	2:B:169:TYR:CE2	2.54	0.43
2:F:106:PRO:HB2	2:F:108:GLN:H	1.84	0.43
2:H:74:ASP:O	2:H:78:ARG:CZ	2.67	0.43
1:C:105:LEU:HA	1:C:105:LEU:HD23	1.76	0.43
1:G:110:ASP:OD1	1:G:146:HIS:HB3	2.19	0.43
2:F:94:GLN:HA	2:F:95:PRO:HD3	1.78	0.42
2:B:145:LEU:HD12	2:B:145:LEU:HA	1.91	0.42
1:E:134:GLU:OE1	1:E:147:LYS:NZ	2.52	0.42
1:E:140:ARG:NH1	1:E:142:ASP:OD2	2.52	0.42
2:B:14:GLU:OE2	2:B:27:ARG:NH1	2.51	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:5:GLU:OE1	2:F:7:TYR:OH	2.16	0.42
1:C:73:LEU:HD22	2:D:30:TYR:HB2	2.02	0.42
1:C:39:LYS:HZ3	1:C:60:LEU:HD22	1.83	0.42
1:C:57:GLN:HA	1:C:60:LEU:HD12	2.01	0.42
2:H:145:LEU:HD23	2:H:145:LEU:HA	1.80	0.42
1:A:126:GLU:HA	1:A:127:PRO:HD3	1.88	0.41
1:A:154:VAL:HA	1:A:155:PRO:HD2	1.87	0.41
2:B:113:LEU:HA	2:B:113:LEU:HD12	1.94	0.41
2:H:25:LEU:HA	2:H:38:PHE:O	2.20	0.41
1:C:162:ASP:OD2	2:F:160:GLU:OE1	2.39	0.41
2:H:26:GLU:CD	2:H:69:LYS:HZ2	2.23	0.41
1:E:105:LEU:HD23	1:E:105:LEU:HA	1.81	0.41
1:E:101:GLN:HA	1:E:102:PRO:HD3	1.97	0.41
2:D:107:LEU:HA	2:D:107:LEU:HD12	1.94	0.41
1:E:32:PHE:HB3	1:E:43:TRP:CE3	2.56	0.41
2:D:49:THR:HG22	2:F:48:VAL:O	2.20	0.40
1:E:89:VAL:O	1:E:176:LYS:HE2	2.21	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:D:128:ARG:NH2	2:F:132:ASN:O[1_455]	2.01	0.19

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	179/193 (93%)	179 (100%)	0	0	100	100
1	C	179/193 (93%)	179 (100%)	0	0	100	100
1	E	179/193 (93%)	179 (100%)	0	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	G	179/193 (93%)	179 (100%)	0	0	100	100
2	B	191/221 (86%)	185 (97%)	6 (3%)	0	100	100
2	D	191/221 (86%)	185 (97%)	5 (3%)	1 (0%)	32	46
2	F	191/221 (86%)	182 (95%)	9 (5%)	0	100	100
2	H	191/221 (86%)	186 (97%)	5 (3%)	0	100	100
All	All	1480/1656 (89%)	1454 (98%)	25 (2%)	1 (0%)	55	72

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	106	PRO

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	162/172 (94%)	159 (98%)	3 (2%)	62	80
1	C	162/172 (94%)	158 (98%)	4 (2%)	53	73
1	E	162/172 (94%)	159 (98%)	3 (2%)	62	80
1	G	162/172 (94%)	160 (99%)	2 (1%)	75	88
2	B	177/194 (91%)	174 (98%)	3 (2%)	66	82
2	D	177/194 (91%)	176 (99%)	1 (1%)	89	95
2	F	177/194 (91%)	175 (99%)	2 (1%)	78	90
2	H	177/194 (91%)	174 (98%)	3 (2%)	66	82
All	All	1356/1464 (93%)	1335 (98%)	21 (2%)	70	85

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	66	LEU
1	A	73	LEU

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Mol	Chain	Res	Type
1	A	160	VAL
2	B	35	LEU
2	B	109	HIS
2	B	142	SER
1	C	46	GLU
1	C	66	LEU
1	C	73	LEU
1	C	129	THR
2	D	35	LEU
1	E	66	LEU
1	E	73	LEU
1	E	160	VAL
2	F	-19	LYS
2	F	137	THR
1	G	66	LEU
1	G	73	LEU
2	H	10	GLN
2	H	35	LEU
2	H	102	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	16	HIS
2	D	109	HIS
2	D	110	HIS
2	D	126	GLN
2	H	-12	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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry

4 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
3	NAG	A	201	1	14,14,15	0.42	0	15,19,21	1.18	2 (13%)
3	NAG	C	201	1	14,14,15	0.49	0	15,19,21	1.21	2 (13%)
3	NAG	E	201	1	14,14,15	0.50	0	15,19,21	1.00	1 (6%)
3	NAG	G	201	1	14,14,15	0.54	0	15,19,21	0.76	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	A	201	1	-	0/6/23/26	0/1/1/1
3	NAG	C	201	1	-	0/6/23/26	0/1/1/1
3	NAG	E	201	1	-	0/6/23/26	0/1/1/1
3	NAG	G	201	1	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	201	NAG	C4-C3-C2	-2.02	108.06	111.02
3	C	201	NAG	O5-C1-C2	2.03	114.30	111.47
3	E	201	NAG	C1-O5-C5	2.27	115.30	112.17
3	A	201	NAG	C1-O5-C5	3.21	116.60	112.17
3	C	201	NAG	C1-O5-C5	3.42	116.88	112.17

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	201	NAG	1	0
3	E	201	NAG	2	0

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	181/193 (93%)	-0.36	0 100 100	16, 25, 42, 54	0
1	C	181/193 (93%)	-0.30	0 100 100	16, 25, 42, 55	0
1	E	181/193 (93%)	-0.28	0 100 100	16, 25, 42, 54	0
1	G	181/193 (93%)	-0.35	1 (0%) 89 87	16, 25, 42, 54	0
2	B	195/221 (88%)	-0.12	5 (2%) 56 54	17, 28, 53, 90	0
2	D	195/221 (88%)	-0.04	7 (3%) 43 42	17, 28, 53, 90	0
2	F	195/221 (88%)	0.09	10 (5%) 29 27	18, 28, 53, 91	0
2	H	195/221 (88%)	-0.08	8 (4%) 38 36	18, 28, 53, 90	0
All	All	1504/1656 (90%)	-0.17	31 (2%) 64 61	16, 26, 48, 91	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	F	106	PRO	8.7
2	F	107	LEU	8.2
2	D	108	GLN	7.3
2	F	108	GLN	7.3
2	F	105	GLY	5.5
2	D	106	PRO	5.4
2	H	105	GLY	4.3
2	H	108	GLN	4.3
2	H	106	PRO	4.0
2	H	104	LYS	3.9
2	F	104	LYS	3.8
2	H	107	LEU	3.7
2	B	105	GLY	3.7
2	D	109	HIS	3.6
2	F	109	HIS	3.4
2	F	164	GLN	3.4

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Mol	Chain	Res	Type	RSRZ
2	F	134	GLN	3.0
2	F	188	ALA	2.9
2	D	105	GLY	2.8
2	B	109	HIS	2.7
2	D	107	LEU	2.7
2	D	103	LYS	2.6
2	H	165	GLN	2.5
2	D	189	GLN	2.4
2	H	109	HIS	2.3
2	F	189	GLN	2.3
2	B	137	THR	2.1
2	B	108	GLN	2.1
2	H	189	GLN	2.1
1	G	181	GLN	2.1
2	B	4	PRO	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](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	NAG	C	201	14/15	0.87	0.22	4.28	30,54,65,67	0
3	NAG	G	201	14/15	0.88	0.18	3.21	26,48,68,76	0
3	NAG	E	201	14/15	0.89	0.16	0.70	22,42,52,82	0
3	NAG	A	201	14/15	0.90	0.15	-0.32	32,45,58,59	0

6.5 Other polymers [i](#)

There are no such residues in this entry.