



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 12:52 am GMT

PDB ID : 3PWV
Title : An immunodominant CTL epitope from rinderpest virus presented by cattle MHC class I molecule N*01801 (BoLA-A11)
Authors : Li, X.; Liu, J.; Qi, J.; Gao, F.; Li, Q.; Li, X.; Zhang, N.; Xia, C.; Gao, G.F.
Deposited on : 2010-12-09
Resolution : 2.70 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
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
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) : recalc28949

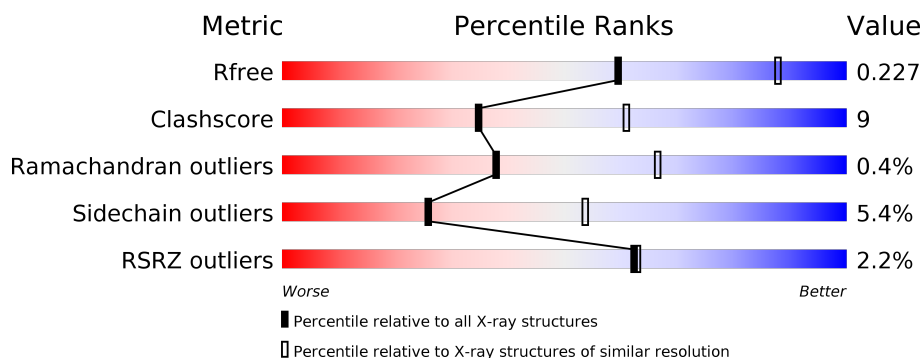
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	2259 (2.70-2.70)
Clashscore	112137	2590 (2.70-2.70)
Ramachandran outliers	110173	2550 (2.70-2.70)
Sidechain outliers	110143	2550 (2.70-2.70)
RSRZ outliers	101464	2275 (2.70-2.70)

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	274	<div> <div>%</div> <div> <div></div> <div>80%</div> <div>18%</div> <div>.</div> </div> </div>
1	D	274	<div> <div>3%</div> <div> <div></div> <div>78%</div> <div>19%</div> <div>.</div> </div> </div>
2	B	98	<div> <div></div> <div> <div>83%</div> <div>16%</div> <div>.</div> </div> </div>
2	E	98	<div> <div>6%</div> <div> <div></div> <div>78%</div> <div>20%</div> <div>.</div> </div> </div>
3	C	9	<div> <div></div> <div> <div>67%</div> <div>33%</div> </div> </div>
3	F	9	<div> <div></div> <div> <div>33%</div> <div>22%</div> <div>33%</div> <div>11%</div> </div> </div>

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 6381 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 I antigen.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	274	Total	C	N	O	S	0	0	0
			2243	1405	400	430	8			
1	D	274	Total	C	N	O	S	0	0	0
			2243	1405	400	430	8			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	98	Total	C	N	O	S	0	0	0
			823	528	142	151	2			
2	E	98	Total	C	N	O	S	0	0	0
			823	528	142	151	2			

- Molecule 3 is a protein called 9-mer peptide from Hemagglutinin.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	9	Total	C	N	O	0	0	0
			67	46	9	12			
3	F	9	Total	C	N	O	0	0	0
			67	46	9	12			

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	62	Total	O	0	0
			62	62		
4	B	26	Total	O	0	0
			26	26		
4	C	1	Total	O	0	0
			1	1		
4	D	19	Total	O	0	0
			19	19		

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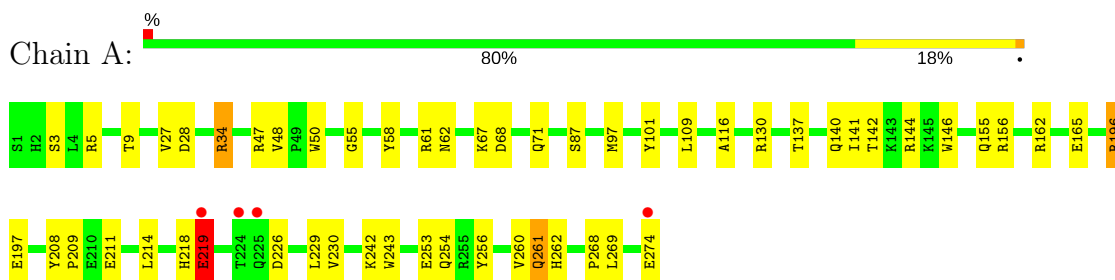
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	E	5	Total	O	0	0
			5	5		
4	F	2	Total	O	0	0
			2	2		

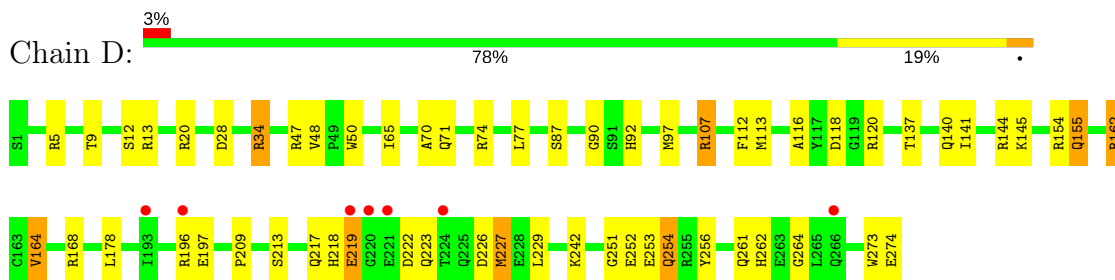
3 Residue-property plots [i](#)

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.

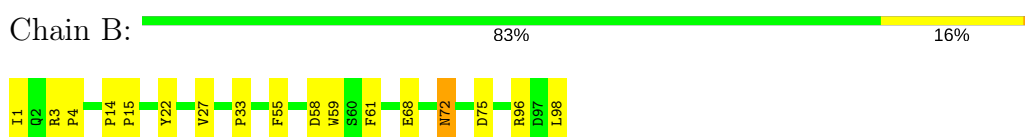
- Molecule 1: MHC class I antigen



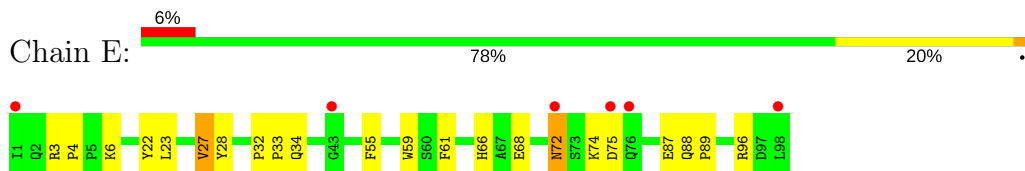
- Molecule 1: MHC class I antigen



- Molecule 2: Beta-2-microglobulin



- Molecule 2: Beta-2-microglobulin



- Molecule 3: 9-mer peptide from Hemagglutinin





- Molecule 3: 9-mer peptide from Hemagglutinin

Chain F:
33% 22% 33% 11%

A horizontal quality scale bar for Chain F. It is divided into four segments: green (33%), yellow (22%), orange (33%), and red (11%).



4 Data and refinement statistics

Property	Value	Source
Space group	P 41	Depositor
Cell constants a, b, c, α , β , γ	83.91Å 83.91Å 153.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.23 – 2.70 38.68 – 2.70	Depositor EDS
% Data completeness (in resolution range)	97.5 (30.23-2.70) 99.4 (38.68-2.70)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	0.14	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.63 (at 2.69Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.5_2)	Depositor
R, R_{free}	0.188 , 0.232 0.181 , 0.227	Depositor DCC
R_{free} test set	1471 reflections (5.08%)	DCC
Wilson B-factor (Å ²)	42.4	Xtriage
Anisotropy	0.113	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 44.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.054 for h,-k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6381	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.79% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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	0.28	0/2305	0.47	0/3131
1	D	0.30	0/2305	0.46	0/3131
2	B	0.28	0/849	0.47	0/1151
2	E	0.25	0/849	0.43	0/1151
3	C	0.86	0/68	0.61	0/92
3	F	1.19	0/68	1.23	1/92 (1.1%)
All	All	0.32	0/6444	0.48	1/8748 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	8	THR	N-CA-CB	-5.24	100.34	110.30

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	2243	0	2095	35	1
1	D	2243	0	2095	55	0
2	B	823	0	801	13	0
2	E	823	0	801	17	0
3	C	67	0	75	2	0
3	F	67	0	75	15	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	62	0	0	0	1
4	B	26	0	0	0	0
4	C	1	0	0	0	0
4	D	19	0	0	0	1
4	E	5	0	0	0	0
4	F	2	0	0	1	0
All	All	6381	0	5942	115	2

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 9.

All (115) 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)
3:F:6:VAL:HG12	3:F:7:LEU:N	1.73	1.01
2:E:72:ASN:HD21	2:E:75:ASP:H	1.02	0.98
1:D:107:ARG:CG	1:D:107:ARG:HH11	1.89	0.84
1:D:262:HIS:CD2	1:D:264:GLY:H	1.97	0.83
2:B:72:ASN:HD21	2:B:75:ASP:H	1.28	0.82
1:D:113:MET:SD	1:D:155:GLN:HG3	2.20	0.81
3:F:8:THR:HG23	4:F:10:HOH:O	1.81	0.79
3:F:6:VAL:HG12	3:F:7:LEU:H	1.45	0.78
2:E:72:ASN:ND2	2:E:75:ASP:H	1.81	0.78
1:A:137:THR:HG23	1:D:252:GLU:OE2	1.85	0.77
1:D:145:LYS:HZ1	3:F:8:THR:HG22	1.50	0.76
1:D:229:LEU:HD22	1:D:242:LYS:HE3	1.69	0.73
1:A:219:GLU:HG2	1:A:219:GLU:O	1.88	0.72
1:A:146:TRP:CZ2	3:C:7:LEU:HB3	2.26	0.71
1:D:262:HIS:HD2	1:D:264:GLY:H	1.36	0.71
1:D:34:ARG:HD2	1:D:47:ARG:HD2	1.74	0.70
1:A:261:GLN:HG3	1:A:268:PRO:HA	1.77	0.67
1:D:107:ARG:HH11	1:D:107:ARG:HG2	1.59	0.66
1:D:217:GLN:HA	1:D:222:ASP:HA	1.78	0.66
1:D:162:ARG:NH2	3:F:4:TYR:HE2	1.93	0.65
1:D:137:THR:HA	1:D:140:GLN:HG3	1.79	0.64
1:D:162:ARG:NE	3:F:1:ILE:HD11	2.12	0.63
1:D:107:ARG:HG3	1:D:107:ARG:HH11	1.60	0.63
1:D:145:LYS:NZ	3:F:8:THR:HG22	2.13	0.62
2:B:33:PRO:HG3	2:B:61:PHE:CZ	2.35	0.61
1:D:137:THR:O	1:D:140:GLN:HB2	2.02	0.60
1:A:218:HIS:HD2	1:A:256:TYR:CZ	2.19	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:219:GLU:HG2	1:D:219:GLU:O	2.01	0.59
1:D:209:PRO:O	1:D:262:HIS:HE1	1.85	0.59
1:D:90:GLY:O	1:D:92:HIS:HD2	1.85	0.59
2:E:72:ASN:ND2	2:E:74:LYS:H	2.01	0.59
1:A:34:ARG:HD2	1:A:47:ARG:HD2	1.84	0.58
2:E:22:TYR:CE1	2:E:68:GLU:HG2	2.39	0.57
1:D:5:ARG:NH2	1:D:28:ASP:O	2.37	0.57
1:A:197:GLU:HG2	1:D:120:ARG:NH2	2.20	0.57
1:A:34:ARG:HD3	1:A:47:ARG:HH11	1.69	0.57
1:A:97:MET:HG3	2:B:55:PHE:HE1	1.71	0.56
1:A:137:THR:HA	1:A:140:GLN:HG3	1.88	0.56
1:D:140:GLN:O	1:D:144:ARG:HG3	2.06	0.55
1:A:209:PRO:O	1:A:262:HIS:HE1	1.90	0.55
1:D:162:ARG:NH2	3:F:4:TYR:CE2	2.73	0.55
1:D:162:ARG:HE	3:F:1:ILE:HD11	1.73	0.54
1:D:155:GLN:HA	1:D:155:GLN:NE2	2.22	0.54
1:A:97:MET:HG3	2:B:55:PHE:CE1	2.42	0.54
2:B:22:TYR:CE1	2:B:68:GLU:HG2	2.43	0.53
2:E:33:PRO:HG3	2:E:61:PHE:CZ	2.43	0.53
1:A:142:THR:HG23	3:C:9:ILE:HA	1.89	0.53
2:E:72:ASN:C	2:E:72:ASN:HD22	2.12	0.53
1:A:34:ARG:CD	1:A:47:ARG:HH11	2.22	0.53
1:D:34:ARG:CD	1:D:47:ARG:HH11	2.21	0.52
1:A:61:ARG:HH11	1:A:62:ASN:ND2	2.09	0.51
1:D:116:ALA:HB2	2:E:59:TRP:CE2	2.46	0.51
1:D:97:MET:O	1:D:113:MET:HA	2.11	0.50
1:D:217:GLN:HG2	1:D:222:ASP:HB3	1.94	0.50
1:D:5:ARG:HG2	1:D:112:PHE:HZ	1.77	0.50
1:A:116:ALA:HB2	2:B:59:TRP:CE2	2.48	0.49
3:F:6:VAL:CG1	3:F:7:LEU:H	2.06	0.49
1:A:229:LEU:HD22	1:A:242:LYS:HE3	1.94	0.49
1:D:48:VAL:HG11	1:D:50:TRP:CE2	2.47	0.49
1:A:230:VAL:HG11	1:A:243:TRP:CZ2	2.48	0.49
2:B:14:PRO:HA	2:B:15:PRO:HD3	1.75	0.48
2:E:23:LEU:O	2:E:66:HIS:HA	2.13	0.48
1:A:141:ILE:HD13	1:A:144:ARG:HH12	1.79	0.48
3:F:1:ILE:HB	3:F:2:PRO:HD2	1.95	0.47
2:B:1:ILE:H1	1:D:197:GLU:CD	2.17	0.47
1:A:219:GLU:O	1:A:219:GLU:CG	2.62	0.47
2:B:72:ASN:H	2:B:72:ASN:HD22	1.62	0.47
1:D:164:VAL:O	1:D:168:ARG:HG3	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:196:ARG:NH2	1:D:118:ASP:HB2	2.31	0.46
1:A:67:LYS:O	1:A:71:GLN:HG3	2.16	0.46
1:D:162:ARG:CZ	3:F:1:ILE:HD11	2.45	0.46
1:A:3:SER:HB3	1:A:101:TYR:CD1	2.52	0.45
1:D:218:HIS:HD2	1:D:256:TYR:CE1	2.34	0.45
1:A:55:GLY:O	1:A:58:TYR:HB3	2.15	0.45
1:D:223:GLN:O	1:D:227:MET:HG2	2.16	0.45
1:D:70:ALA:O	1:D:74:ARG:HG3	2.17	0.45
1:D:107:ARG:CG	1:D:107:ARG:NH1	2.59	0.45
1:D:50:TRP:CZ2	1:D:178:LEU:HD11	2.51	0.45
2:E:88:GLN:HB2	2:E:89:PRO:HD2	1.99	0.45
2:B:72:ASN:HD22	2:B:72:ASN:N	2.15	0.44
1:D:5:ARG:HG2	1:D:112:PHE:CZ	2.51	0.44
1:D:196:ARG:HG3	1:D:197:GLU:HG3	1.99	0.44
1:A:218:HIS:HD2	1:A:256:TYR:CE1	2.36	0.44
1:A:214:LEU:CD2	1:A:260:VAL:HG22	2.47	0.44
2:E:3:ARG:HA	2:E:4:PRO:HD3	1.82	0.43
1:A:5:ARG:NH2	1:A:28:ASP:O	2.51	0.43
2:B:3:ARG:HA	2:B:4:PRO:HD3	1.82	0.43
2:B:58:ASP:O	2:B:59:TRP:HB2	2.17	0.43
2:E:87:GLU:HA	2:E:87:GLU:OE1	2.19	0.43
1:A:208:TYR:CG	1:A:209:PRO:HA	2.54	0.42
1:A:48:VAL:HG11	1:A:50:TRP:CE2	2.54	0.42
3:F:6:VAL:CG1	3:F:7:LEU:N	2.45	0.42
1:A:130:ARG:HG2	1:A:156:ARG:HH21	1.84	0.42
2:E:32:PRO:HB2	2:E:34:GLN:OE1	2.19	0.42
1:A:196:ARG:HH22	1:D:118:ASP:HB2	1.84	0.42
2:E:6:LYS:O	2:E:27:VAL:HA	2.20	0.42
1:D:107:ARG:NH1	1:D:107:ARG:HG3	2.28	0.42
1:D:141:ILE:HD13	1:D:144:ARG:HH12	1.84	0.42
2:E:28:TYR:CD2	2:E:28:TYR:N	2.88	0.42
1:D:65:ILE:HG21	3:F:2:PRO:HB2	2.02	0.41
2:B:33:PRO:HG3	2:B:61:PHE:CE2	2.54	0.41
1:D:154:ARG:NH2	3:F:6:VAL:O	2.53	0.41
1:A:261:GLN:HG3	1:A:268:PRO:CA	2.48	0.41
1:D:254:GLN:HE22	1:D:274:GLU:HA	1.86	0.41
2:E:72:ASN:HD22	2:E:74:LYS:H	1.65	0.41
1:D:13:ARG:CZ	1:D:20:ARG:HB2	2.50	0.41
1:D:213:SER:HB3	1:D:261:GLN:HB2	2.02	0.41
1:D:97:MET:HG3	2:E:55:PHE:HE1	1.86	0.41
2:E:33:PRO:HG3	2:E:61:PHE:CE2	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:254:GLN:HE22	1:A:274:GLU:HA	1.85	0.41
1:D:12:SER:HB3	1:D:77:LEU:HD13	2.03	0.41
1:D:251:GLY:N	1:D:253:GLU:OE2	2.49	0.41
1:A:27:VAL:O	1:A:28:ASP:HB2	2.20	0.41
1:D:254:GLN:NE2	1:D:273:TRP:O	2.54	0.40
1:A:196:ARG:HH22	1:D:118:ASP:CB	2.34	0.40

All (2) 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)
1:A:68:ASP:OD2	3:F:8:THR:CG2[1_565]	2.12	0.08
4:A:335:HOH:O	4:D:285:HOH:O[1_565]	2.17	0.03

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	272/274 (99%)	263 (97%)	8 (3%)	1 (0%)	38	66
1	D	272/274 (99%)	255 (94%)	16 (6%)	1 (0%)	38	66
2	B	96/98 (98%)	93 (97%)	3 (3%)	0	100	100
2	E	96/98 (98%)	90 (94%)	6 (6%)	0	100	100
3	C	7/9 (78%)	7 (100%)	0	0	100	100
3	F	7/9 (78%)	4 (57%)	2 (29%)	1 (14%)	0	0
All	All	750/762 (98%)	712 (95%)	35 (5%)	3 (0%)	38	66

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	219	GLU

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Mol	Chain	Res	Type
1	A	219	GLU
3	F	6	VAL

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	232/232 (100%)	218 (94%)	14 (6%)	22	48
1	D	232/232 (100%)	221 (95%)	11 (5%)	30	60
2	B	94/94 (100%)	90 (96%)	4 (4%)	33	64
2	E	94/94 (100%)	91 (97%)	3 (3%)	44	75
3	C	7/7 (100%)	6 (86%)	1 (14%)	4	9
3	F	7/7 (100%)	4 (57%)	3 (43%)	0	0
All	All	666/666 (100%)	630 (95%)	36 (5%)	26	54

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	THR
1	A	34	ARG
1	A	87	SER
1	A	109	LEU
1	A	155	GLN
1	A	162	ARG
1	A	165	GLU
1	A	196	ARG
1	A	211	GLU
1	A	219	GLU
1	A	226	ASP
1	A	253	GLU
1	A	261	GLN
1	A	269	LEU
2	B	27	VAL
2	B	72	ASN

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Mol	Chain	Res	Type
2	B	96	ARG
2	B	98	LEU
3	C	6	VAL
1	D	9	THR
1	D	34	ARG
1	D	71	GLN
1	D	87	SER
1	D	107	ARG
1	D	155	GLN
1	D	162	ARG
1	D	164	VAL
1	D	226	ASP
1	D	227	MET
1	D	254	GLN
2	E	27	VAL
2	E	72	ASN
2	E	96	ARG
3	F	4	TYR
3	F	7	LEU
3	F	8	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	62	ASN
1	A	71	GLN
1	A	218	HIS
1	A	254	GLN
2	B	72	ASN
2	B	88	GLN
1	D	62	ASN
1	D	71	GLN
1	D	155	GLN
1	D	218	HIS
1	D	254	GLN
1	D	261	GLN
1	D	262	HIS
2	E	72	ASN
2	E	88	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	274/274 (100%)	-0.18	4 (1%) 74 75	17, 32, 62, 121	0
1	D	274/274 (100%)	-0.02	7 (2%) 56 56	26, 47, 82, 129	0
2	B	98/98 (100%)	-0.25	0 100 100	20, 31, 53, 70	0
2	E	98/98 (100%)	0.18	6 (6%) 22 20	34, 55, 99, 119	0
3	C	9/9 (100%)	0.04	0 100 100	25, 30, 63, 74	0
3	F	9/9 (100%)	0.35	0 100 100	36, 49, 92, 111	0
All	All	762/762 (100%)	-0.08	17 (2%) 62 63	17, 40, 83, 129	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	E	98	LEU	4.7
2	E	76	GLN	3.7
1	A	225	GLN	3.6
1	A	224	THR	3.4
1	D	196	ARG	3.1
1	D	220	GLY	2.7
1	A	219	GLU	2.6
1	D	224	THR	2.6
1	D	219	GLU	2.5
2	E	1	ILE	2.5
1	D	266	GLN	2.4
1	A	274	GLU	2.4
2	E	75	ASP	2.3
1	D	221	GLU	2.3
1	D	193	ILE	2.2
2	E	43	GLY	2.1
2	E	72	ASN	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 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.