Full wwPDB X-ray Structure Validation Report ⓘ

May 3, 2025 – 09:24 AM EDT

PDB ID	:	3QQ3 / pdb_00003qq3
Title	:	Crystal structure of swine major histocompatibility complex class I SLA-1 0401 and identification of 2009 pandemic swine-origin influenza A H1N1 virus cytotoxic T lymphocyte epitope peptides
Authors	:	Zhang, N.; Qi, J.; Gao, F.; Pan, X.; Chen, R.; Li, Q.; Chen, Z.; Li, X.; Xia, C.; Gao, G.F.
Deposited on	:	2011-02-15
Resolution	:	2.59 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
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references ①) were used in the production of this report:

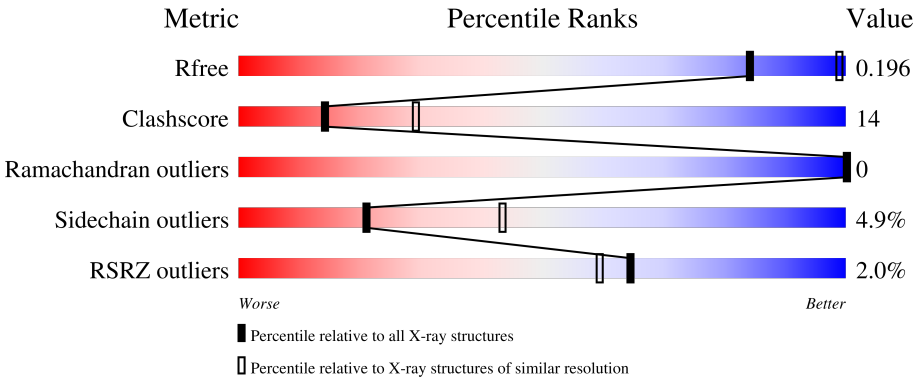
MolProbity	:	4-5-2 with Phenix2.0rc1
Xtriage (Phenix)	:	2.0rc1
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.006 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.43.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.59 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}	164625	3775 (2.60-2.60)
Clashscore	180529	4181 (2.60-2.60)
Ramachandran outliers	177936	4129 (2.60-2.60)
Sidechain outliers	177891	4129 (2.60-2.60)
RSRZ outliers	164620	3775 (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 of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	2% 72% 25% ·
1	D	275	2% 66% 30% ·
2	B	100	% 69% 29% ··
2	E	100	3% 73% 23% · ·
3	C	9	89% 11%

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Mol	Chain	Length	Quality of chain
3	F	9	

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 6593 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	275	Total	C	N	O	S	0	0	0
			2233	1394	398	432	9			
1	D	275	Total	C	N	O	S	0	0	0
			2233	1394	398	432	9			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	99	Total	C	N	O	S	0	0	0
			817	524	142	148	3			
2	E	99	Total	C	N	O	S	0	0	0
			817	524	142	148	3			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	1	GLU	-	expression tag	UNP Q07717
B	2	PHE	-	expression tag	UNP Q07717
E	1	GLU	-	expression tag	UNP Q07717
E	2	PHE	-	expression tag	UNP Q07717

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

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	9	Total	C	N	O	0	0	0
			75	47	12	16			
3	F	9	Total	C	N	O	0	0	0
			75	47	12	16			

- Molecule 4 is water.

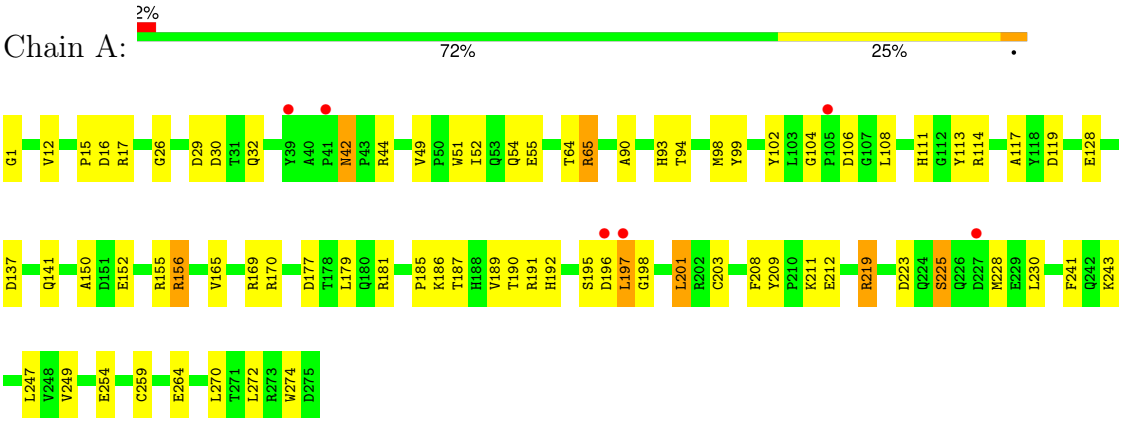
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	112	Total O 112 112	0	0
4	B	72	Total O 72 72	0	0
4	C	6	Total O 6 6	0	0
4	D	102	Total O 102 102	0	0
4	E	43	Total O 43 43	0	0
4	F	8	Total O 8 8	0	0

3 Residue-property plots (i)

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

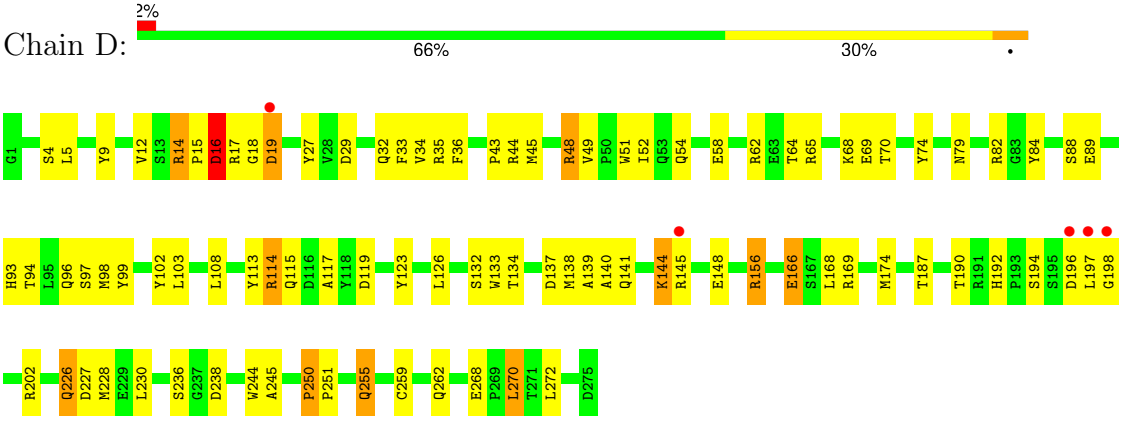
• Molecule 1: MHC class I antigen

Chain A:



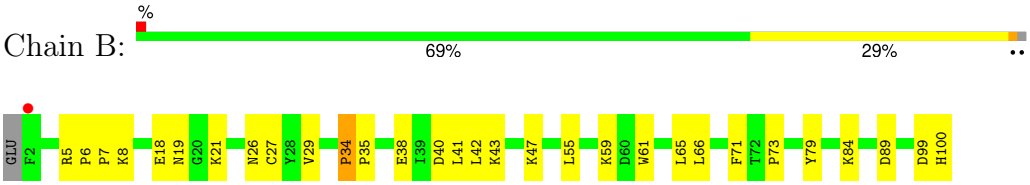
• Molecule 1: MHC class I antigen

Chain D:



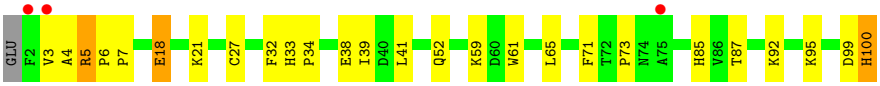
• Molecule 2: Beta-2-microglobulin

Chain B:



• Molecule 2: Beta-2-microglobulin

Chain E:

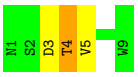


• Molecule 3: 9-mer peptide from Neuraminidase

Chain C:

• Molecule 3: 9-mer peptide from Neuraminidase

Chain F:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	96.73Å 37.65Å 111.43Å 90.00° 113.06° 90.00°	Depositor
Resolution (Å)	28.31 – 2.59 28.31 – 2.59	Depositor EDS
% Data completeness (in resolution range)	99.7 (28.31-2.59) 99.8 (28.31-2.59)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.03 (at 2.51Å)	Xtriage
Refinement program	PHENIX (phenix.refine)	Depositor
R, R_{free}	0.200 , 0.262 0.195 , 0.196	Depositor DCC
R_{free} test set	1212 reflections (5.13%)	wwPDB-VP
Wilson B-factor (Å^2)	27.4	Xtriage
Anisotropy	0.205	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 36.8	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	0.006 for h,-k,-h-l	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	6593	wwPDB-VP
Average B, all atoms (Å^2)	27.0	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 26.63 % 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 2.5453e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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	0.40	0/2293	0.80	5/3114 (0.2%)
1	D	0.47	0/2293	0.81	7/3114 (0.2%)
2	B	0.27	0/843	0.74	4/1145 (0.3%)
2	E	0.29	0/843	0.78	4/1145 (0.3%)
3	C	0.20	0/78	0.41	0/106
3	F	0.22	0/78	0.50	0/106
All	All	0.39	0/6428	0.79	20/8730 (0.2%)

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	A	0	1
1	D	0	1
All	All	0	2

There are no bond length outliers.

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	17	ARG	N-CA-C	-9.59	100.06	114.64
1	A	16	ASP	N-CA-C	8.74	124.00	113.16
1	D	144	LYS	N-CA-C	8.44	122.13	111.24
1	D	18	GLY	CA-C-N	-6.97	108.23	121.54
1	D	18	GLY	C-N-CA	-6.97	108.23	121.54
1	D	19	ASP	CB-CA-C	-6.15	98.18	110.42
2	E	34	PRO	CA-C-N	5.88	125.58	119.05
2	E	34	PRO	C-N-CA	5.88	125.58	119.05
2	E	5	ARG	CA-C-N	5.70	123.82	119.66
2	E	5	ARG	C-N-CA	5.70	123.82	119.66

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	D	250	PRO	CA-C-N	5.59	125.30	119.82
1	D	250	PRO	C-N-CA	5.59	125.30	119.82
1	A	104	GLY	CA-C-N	5.57	125.41	119.28
1	A	104	GLY	C-N-CA	5.57	125.41	119.28
1	A	29	ASP	CB-CA-C	-5.45	110.28	116.54
2	B	5	ARG	CA-C-N	5.32	123.55	119.66
2	B	5	ARG	C-N-CA	5.32	123.55	119.66
2	B	34	PRO	CA-C-N	5.04	124.83	119.28
2	B	34	PRO	C-N-CA	5.04	124.83	119.28
1	D	29	ASP	CB-CA-C	-5.02	110.81	116.63

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	15	PRO	Mainchain
1	D	16	ASP	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	2233	0	2096	60	1
1	D	2233	0	2096	76	0
2	B	817	0	788	19	0
2	E	817	0	788	22	0
3	C	75	0	61	1	0
3	F	75	0	61	2	0
4	A	112	0	0	5	0
4	B	72	0	0	2	1
4	C	6	0	0	0	0
4	D	102	0	0	3	0
4	E	43	0	0	3	0
4	F	8	0	0	0	0
All	All	6593	0	5890	170	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 14.

All (170) 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:D:14:ARG:HH21	1:D:19:ASP:CG	1.57	1.12
2:E:6:PRO:HA	2:E:87:THR:HG21	1.37	1.04
1:D:14:ARG:NH2	1:D:19:ASP:OD1	1.91	1.02
1:D:14:ARG:NH2	1:D:19:ASP:CG	2.27	0.91
2:E:3:VAL:HG12	4:E:331:HOH:O	1.69	0.91
1:D:196:ASP:C	1:D:197:LEU:HD23	1.98	0.88
1:A:196:ASP:O	1:A:197:LEU:HG	1.84	0.77
1:D:196:ASP:O	1:D:197:LEU:HD23	1.85	0.76
1:D:197:LEU:O	1:D:197:LEU:HG	1.87	0.74
1:D:194:SER:N	1:D:198:GLY:O	2.21	0.74
1:D:14:ARG:NH2	1:D:19:ASP:OD2	2.22	0.73
1:D:19:ASP:OD1	1:D:19:ASP:C	2.31	0.73
1:A:259:CYS:HB3	1:A:272:LEU:HB3	1.72	0.71
1:D:244:TRP:HE1	2:E:100:HIS:HD2	1.38	0.71
1:D:244:TRP:NE1	2:E:100:HIS:HD2	1.89	0.69
1:D:69:GLU:HB3	3:F:5:VAL:HG21	1.74	0.69
1:A:223:ASP:OD2	1:A:225:SER:HB2	1.94	0.68
1:A:152:GLU:OE1	3:C:7:TRP:HB2	1.95	0.66
1:D:54:GLN:HG3	1:D:174:MET:HE3	1.75	0.65
1:D:16:ASP:OD2	1:D:16:ASP:N	2.30	0.65
1:D:114:ARG:HD3	1:D:114:ARG:C	2.22	0.65
2:E:4:ALA:O	4:E:331:HOH:O	2.14	0.64
1:D:35:ARG:HG2	1:D:48:ARG:NE	2.14	0.63
1:D:123:TYR:CZ	1:D:140:ALA:HA	2.33	0.63
2:B:40:ASP:OD1	2:B:47:LYS:HE3	1.98	0.62
1:D:244:TRP:HE1	2:E:100:HIS:CD2	2.17	0.62
1:D:194:SER:OG	1:D:198:GLY:C	2.43	0.62
2:E:27:CYS:HB2	2:E:41:LEU:HD21	1.82	0.61
1:A:93:HIS:HD2	1:A:119:ASP:OD2	1.84	0.61
1:A:152:GLU:HA	1:A:155:ARG:NH1	2.17	0.60
1:A:102:TYR:HB2	1:A:111:HIS:HB2	1.83	0.60
1:A:150:ALA:O	1:A:155:ARG:NH2	2.35	0.59
1:A:191:ARG:HD3	1:A:274:TRP:CZ2	2.37	0.59
1:A:12:VAL:HG22	1:A:94:THR:HG22	1.83	0.59
1:D:228:MET:HE3	1:D:230:LEU:HD13	1.83	0.59
1:A:65:ARG:HH11	1:A:65:ARG:HB3	1.69	0.58
1:A:196:ASP:O	1:A:197:LEU:CG	2.50	0.58
1:D:44:ARG:HD2	1:D:64:THR:HG21	1.86	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:202:ARG:HD3	1:D:244:TRP:CE3	2.39	0.57
2:E:71:PHE:CE2	2:E:73:PRO:HG3	2.39	0.57
1:D:58:GLU:O	1:D:62:ARG:HD3	2.04	0.56
1:A:230:LEU:HD11	1:A:243:LYS:HE3	1.88	0.55
1:A:117:ALA:HB2	2:B:61:TRP:CE2	2.42	0.55
1:D:196:ASP:O	1:D:197:LEU:CD2	2.55	0.55
1:D:99:TYR:HB3	1:D:114:ARG:HG2	1.90	0.54
1:A:114:ARG:HD3	1:A:114:ARG:C	2.32	0.54
1:D:255:GLN:HA	4:D:324:HOH:O	2.07	0.54
1:A:197:LEU:HG	1:A:197:LEU:O	2.08	0.54
2:B:19:ASN:HB3	4:B:322:HOH:O	2.08	0.53
1:D:98:MET:HG2	1:D:113:TYR:HE1	1.72	0.53
1:A:191:ARG:HG3	1:A:201:LEU:CD1	2.38	0.53
1:D:226:GLN:H	1:D:226:GLN:CD	2.16	0.53
2:E:7:PRO:HD3	2:E:87:THR:HG21	1.91	0.53
2:B:27:CYS:HB2	2:B:41:LEU:HD21	1.91	0.53
1:D:270:LEU:HD13	1:D:272:LEU:HD11	1.92	0.52
1:A:189:VAL:HG23	1:A:272:LEU:HG	1.91	0.52
1:A:185:PRO:HA	1:A:208:PHE:HB3	1.92	0.52
2:B:43:LYS:HG3	2:B:79:TYR:CE2	2.44	0.52
1:D:70:THR:O	1:D:74:TYR:HD2	1.92	0.52
2:E:7:PRO:HD3	2:E:87:THR:CG2	2.40	0.51
2:B:55:LEU:HA	2:B:65:LEU:HD21	1.93	0.51
2:B:99:ASP:O	2:B:100:HIS:HB2	2.10	0.51
1:D:230:LEU:HD12	1:D:245:ALA:HB2	1.92	0.51
2:B:18:GLU:HB2	2:B:21:LYS:HG3	1.92	0.51
1:A:198:GLY:HA2	4:A:325:HOH:O	2.11	0.51
1:D:145:ARG:HA	1:D:148:GLU:HB2	1.93	0.50
1:A:106:ASP:OD2	1:A:106:ASP:N	2.44	0.50
1:A:211:LYS:HG3	1:A:241:PHE:CE1	2.46	0.50
1:D:103:LEU:HB3	4:D:335:HOH:O	2.11	0.50
1:A:201:LEU:HD22	1:A:249:VAL:HG21	1.93	0.50
2:B:42:LEU:HD23	2:B:47:LYS:HA	1.94	0.50
1:D:226:GLN:H	1:D:226:GLN:NE2	2.10	0.49
1:D:93:HIS:HD2	1:D:119:ASP:OD2	1.95	0.49
1:D:270:LEU:HD13	1:D:272:LEU:CD1	2.42	0.49
1:D:141:GLN:O	1:D:144:LYS:N	2.42	0.49
1:D:15:PRO:O	1:D:16:ASP:C	2.51	0.49
1:D:49:VAL:O	1:D:52:ILE:HG22	2.13	0.49
1:D:96:GLN:NE2	2:E:33:HIS:NE2	2.61	0.48
2:E:85:HIS:ND1	2:E:87:THR:HG22	2.28	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:126:LEU:HD12	1:D:132:SER:O	2.13	0.48
1:D:64:THR:O	1:D:68:LYS:HG3	2.14	0.48
1:A:55:GLU:OE1	1:A:170:ARG:NH2	2.46	0.48
1:A:219:ARG:O	1:A:219:ARG:HG3	2.13	0.48
1:A:111:HIS:HD2	1:A:128:GLU:OE2	1.96	0.48
1:A:195:SER:C	1:A:197:LEU:H	2.21	0.48
1:A:51:TRP:O	1:A:54:GLN:HG2	2.14	0.47
1:D:36:PHE:CE1	1:D:43:PRO:HB2	2.49	0.47
1:D:98:MET:HE3	1:D:115:GLN:NE2	2.29	0.47
1:D:84:TYR:HB3	1:D:139:ALA:HB1	1.97	0.47
1:D:255:GLN:HB3	4:D:354:HOH:O	2.14	0.47
1:A:106:ASP:OD1	1:A:108:LEU:HD12	2.15	0.47
1:D:259:CYS:HB3	1:D:272:LEU:HB2	1.96	0.46
1:A:186:LYS:HE3	4:A:384:HOH:O	2.14	0.46
1:D:138:MET:O	1:D:141:GLN:HB2	2.16	0.46
1:D:156:ARG:HD3	1:D:156:ARG:HA	1.55	0.46
1:D:117:ALA:HB2	2:E:61:TRP:CE2	2.50	0.46
1:A:203:CYS:SG	1:A:272:LEU:HD23	2.55	0.46
1:A:201:LEU:HD22	1:A:249:VAL:HG11	1.97	0.46
1:D:137:ASP:O	1:D:141:GLN:HG3	2.15	0.46
1:A:247:LEU:HA	4:A:333:HOH:O	2.15	0.46
1:A:98:MET:HG2	1:A:113:TYR:HE1	1.81	0.45
1:D:9:TYR:HB2	1:D:97:SER:HB2	1.98	0.45
1:A:156:ARG:HD3	1:A:156:ARG:HA	1.44	0.45
1:D:33:PHE:CD2	1:D:34:VAL:HG13	2.51	0.45
1:A:51:TRP:CZ2	1:A:179:LEU:HD11	2.52	0.45
1:A:185:PRO:CA	1:A:208:PHE:HB3	2.46	0.45
1:D:32:GLN:NE2	1:D:48:ARG:HD2	2.32	0.45
1:D:27:TYR:CE2	1:D:32:GLN:HB2	2.52	0.45
3:F:3:ASP:OD1	3:F:4:THR:N	2.49	0.45
2:B:99:ASP:C	2:B:100:HIS:HD1	2.25	0.45
1:A:197:LEU:CG	1:A:197:LEU:O	2.65	0.44
2:B:8:LYS:HE2	4:B:286:HOH:O	2.17	0.44
2:E:71:PHE:CZ	2:E:73:PRO:HG3	2.52	0.44
2:B:34:PRO:CB	2:B:35:PRO:HD2	2.48	0.44
1:A:196:ASP:C	1:A:197:LEU:HD23	2.42	0.44
2:E:18:GLU:O	2:E:21:LYS:HB2	2.18	0.44
1:A:49:VAL:O	1:A:52:ILE:HG22	2.18	0.44
1:D:98:MET:HE2	2:E:61:TRP:CH2	2.53	0.44
2:B:71:PHE:CZ	2:B:73:PRO:HG3	2.53	0.44
1:D:166:GLU:HG3	1:D:169:ARG:HH22	1.83	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:42:ASN:HD22	1:A:42:ASN:HA	1.63	0.43
1:A:99:TYR:HB3	1:A:114:ARG:HG2	1.98	0.43
1:A:212:GLU:HB2	4:A:372:HOH:O	2.18	0.43
1:D:197:LEU:N	1:D:198:GLY:HA3	2.33	0.43
1:D:236:SER:HB2	1:D:238:ASP:OD1	2.18	0.43
1:D:79:ASN:ND2	1:D:82:ARG:NH1	2.65	0.43
2:E:92:LYS:NZ	4:E:110:HOH:O	2.51	0.43
1:D:34:VAL:CG2	1:D:45:MET:HE3	2.48	0.43
2:E:6:PRO:HA	2:E:7:PRO:HD3	1.80	0.43
1:A:165:VAL:O	1:A:169:ARG:HG3	2.19	0.43
1:D:262:GLN:HE21	1:D:262:GLN:HB2	1.63	0.43
2:B:89:ASP:OD2	2:B:89:ASP:N	2.52	0.42
1:D:5:LEU:HB2	1:D:168:LEU:HD13	2.00	0.42
1:D:202:ARG:HD3	1:D:244:TRP:CD2	2.54	0.42
1:A:195:SER:C	1:A:197:LEU:N	2.77	0.42
1:D:187:THR:HB	1:D:272:LEU:HD21	2.01	0.42
1:D:49:VAL:HG21	1:D:51:TRP:CE2	2.54	0.42
1:D:133:TRP:HB2	1:D:144:LYS:HG3	2.01	0.42
1:A:65:ARG:HB3	1:A:65:ARG:NH1	2.34	0.42
2:E:6:PRO:CA	2:E:87:THR:HG21	2.27	0.42
1:A:1:GLY:HA2	4:A:309:HOH:O	2.20	0.42
1:A:187:THR:HB	1:A:272:LEU:CD2	2.49	0.42
1:A:197:LEU:HA	1:A:198:GLY:HA3	1.70	0.42
1:D:14:ARG:HE	1:D:14:ARG:HB2	1.23	0.42
1:D:166:GLU:HG3	1:D:169:ARG:NH2	2.34	0.42
2:E:38:GLU:C	2:E:39:ILE:HG13	2.43	0.42
1:D:34:VAL:HB	1:D:45:MET:HE3	2.02	0.42
2:B:8:LYS:O	2:B:29:VAL:HA	2.19	0.41
1:D:12:VAL:HG22	1:D:94:THR:HG22	2.02	0.41
1:A:177:ASP:O	1:A:181:ARG:HG2	2.20	0.41
1:A:191:ARG:NH1	1:A:254:GLU:OE2	2.53	0.41
1:A:44:ARG:HD2	1:A:64:THR:HG21	2.02	0.41
1:D:250:PRO:HA	1:D:251:PRO:HD3	1.77	0.41
2:E:5:ARG:O	2:E:32:PHE:HA	2.20	0.41
2:B:34:PRO:HB2	2:B:35:PRO:HD2	2.03	0.41
1:D:144:LYS:O	1:D:148:GLU:HG3	2.21	0.41
1:A:30:ASP:HB2	1:A:209:TYR:CE1	2.56	0.41
1:A:90:ALA:O	1:D:108:LEU:HD13	2.21	0.41
2:B:6:PRO:HA	2:B:7:PRO:HD3	1.80	0.41
2:B:26:ASN:HB3	2:B:66:LEU:HD11	2.02	0.41
1:D:190:THR:OG1	1:D:192:HIS:CE1	2.74	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:190:THR:OG1	1:A:192:HIS:CE1	2.75	0.40
1:A:191:ARG:HD3	1:A:274:TRP:CE2	2.56	0.40
1:D:4:SER:HB3	1:D:102:TYR:CD1	2.56	0.40
2:E:52:GLN:OE1	2:E:65:LEU:HD13	2.20	0.40
1:A:49:VAL:HG21	1:A:51:TRP:CE2	2.56	0.40
1:A:137:ASP:O	1:A:141:GLN:HG3	2.21	0.40
2:B:38:GLU:HB2	2:B:84:LYS:HB2	2.04	0.40
1:A:26:GLY:O	1:A:32:GLN:HA	2.20	0.40
1:A:152:GLU:OE1	1:A:155:ARG:NH1	2.55	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)
1:A:196:ASP:OD2	4:B:129:HOH:O[1_565]	2.10	0.10

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	273/275 (99%)	263 (96%)	10 (4%)	0	100	100
1	D	273/275 (99%)	259 (95%)	14 (5%)	0	100	100
2	B	97/100 (97%)	91 (94%)	6 (6%)	0	100	100
2	E	97/100 (97%)	92 (95%)	5 (5%)	0	100	100
3	C	7/9 (78%)	7 (100%)	0	0	100	100
3	F	7/9 (78%)	6 (86%)	1 (14%)	0	100	100
All	All	754/768 (98%)	718 (95%)	36 (5%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains (i)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	234/234 (100%)	224 (96%)	10 (4%)	25 49
1	D	234/234 (100%)	218 (93%)	16 (7%)	13 28
2	B	92/93 (99%)	91 (99%)	1 (1%)	70 86
2	E	92/93 (99%)	87 (95%)	5 (5%)	18 39
3	C	8/8 (100%)	8 (100%)	0	100 100
3	F	8/8 (100%)	7 (88%)	1 (12%)	3 7
All	All	668/670 (100%)	635 (95%)	33 (5%)	21 43

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	42	ASN
1	A	65	ARG
1	A	156	ARG
1	A	197	LEU
1	A	201	LEU
1	A	219	ARG
1	A	225	SER
1	A	228	MET
1	A	264	GLU
1	A	270	LEU
2	B	59	LYS
1	D	14	ARG
1	D	16	ASP
1	D	17	ARG
1	D	48	ARG
1	D	65	ARG
1	D	88	SER
1	D	89	GLU
1	D	114	ARG
1	D	134	THR
1	D	156	ARG
1	D	166	GLU

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Mol	Chain	Res	Type
1	D	226	GLN
1	D	227	ASP
1	D	255	GLN
1	D	268	GLU
1	D	270	LEU
2	E	18	GLU
2	E	59	LYS
2	E	95	LYS
2	E	99	ASP
2	E	100	HIS
3	F	4	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	42	ASN
1	A	54	GLN
1	A	72	GLN
1	A	87	GLN
1	A	93	HIS
1	A	96	GLN
1	A	111	HIS
1	A	192	HIS
1	A	255	GLN
1	A	262	GLN
2	B	26	ASN
2	B	49	ASN
3	C	1	ASN
1	D	42	ASN
1	D	54	GLN
1	D	79	ASN
1	D	87	GLN
1	D	96	GLN
1	D	115	GLN
1	D	141	GLN
1	D	192	HIS
1	D	255	GLN
1	D	262	GLN
2	E	26	ASN
2	E	49	ASN
3	F	1	ASN

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 oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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, 95^{th} 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%)	-0.25	6 (2%) 62 57	11, 23, 49, 63	0
1	D	275/275 (100%)	-0.16	5 (1%) 67 62	16, 27, 47, 66	0
2	B	99/100 (99%)	-0.53	1 (1%) 79 75	13, 18, 30, 81	0
2	E	99/100 (99%)	-0.10	3 (3%) 52 46	19, 31, 52, 83	0
3	C	9/9 (100%)	-0.34	0 100 100	17, 22, 25, 27	0
3	F	9/9 (100%)	-0.19	0 100 100	21, 26, 35, 36	0
All	All	766/768 (99%)	-0.23	15 (1%) 64 59	11, 25, 48, 83	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	145	ARG	5.0
1	A	197	LEU	4.5
2	B	2	PHE	3.7
2	E	2	PHE	3.2
1	D	198	GLY	3.2
1	D	197	LEU	2.6
1	A	196	ASP	2.6
1	A	39	TYR	2.5
2	E	3	VAL	2.4
1	D	196	ASP	2.3
1	A	105	PRO	2.3
1	D	19	ASP	2.2
2	E	75	ALA	2.2
1	A	227	ASP	2.2
1	A	41	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 monosaccharides in this entry.

6.4 Ligands (i)

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.