



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 03:46 PM UTC

PDB ID : 8KB0 / pdb_00008kb0
Title : Crystal structure of 01JD-AEAIIVAMV
Authors : Tang, Z.; Zhang, N.
Deposited on : 2023-08-03
Resolution : 2.48 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 ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

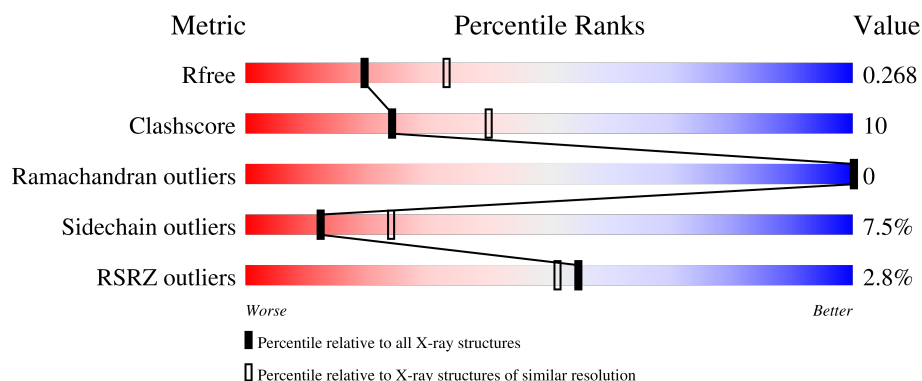
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.48 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}	180053	7589 (2.50-2.46)
Clashscore	190562	8295 (2.50-2.46)
Ramachandran outliers	187476	8164 (2.50-2.46)
Sidechain outliers	187428	8166 (2.50-2.46)
RSRZ outliers	180081	7593 (2.50-2.46)

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 $\leq 5\%$. 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	271	3% 70% 25% ..
1	D	271	4% 76% 21% ..
2	B	103	% 86% 10% .
2	E	103	% 73% 22% ..
3	C	9	67% 22% 11%

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Mol	Chain	Length	Quality of chain
3	F	9	 A horizontal bar chart showing the quality of chain F. The bar is divided into three segments: green (56%), yellow (33%), and orange (11%).

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6391 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 alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	271	Total	C	N	O	S	0	0	0
			2221	1390	394	430	7			
1	D	271	Total	C	N	O	S	0	0	0
			2221	1390	394	430	7			

- Molecule 2 is a protein called Beta2-microglobulin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	99	Total	C	N	O	S	0	0	0
			798	511	134	149	4			
2	E	99	Total	C	N	O	S	0	0	0
			798	511	134	149	4			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-1	GLU	-	expression tag	UNP Q14U75
B	0	PHE	-	expression tag	UNP Q14U75
E	-1	GLU	-	expression tag	UNP Q14U75
E	0	PHE	-	expression tag	UNP Q14U75

- Molecule 3 is a protein called peptide of AIV.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	9	Total	C	N	O	S	0	0	0
			62	41	9	11	1			
3	F	9	Total	C	N	O	S	0	0	0
			62	41	9	11	1			

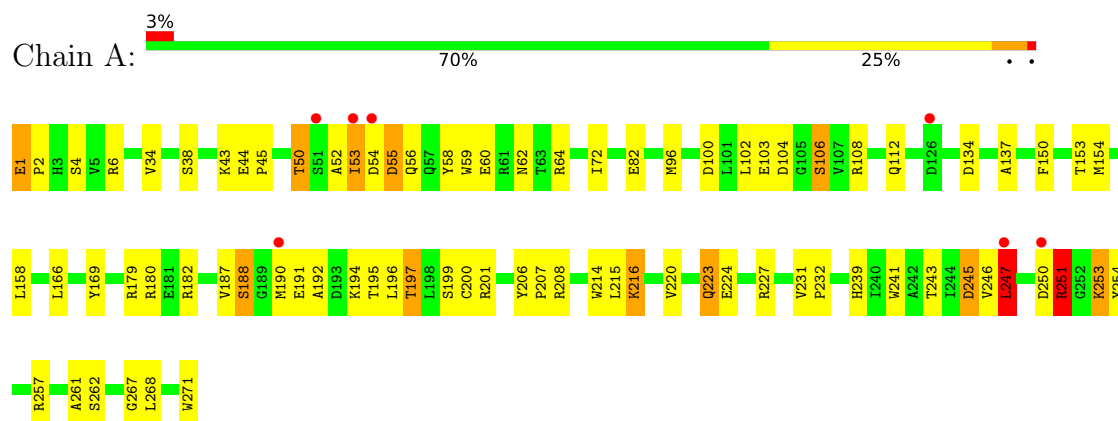
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	83	Total 83	O 83	0	0
4	B	50	Total 50	O 50	0	0
4	C	5	Total 5	O 5	0	0
4	D	60	Total 60	O 60	0	0
4	E	28	Total 28	O 28	0	0
4	F	3	Total 3	O 3	0	0

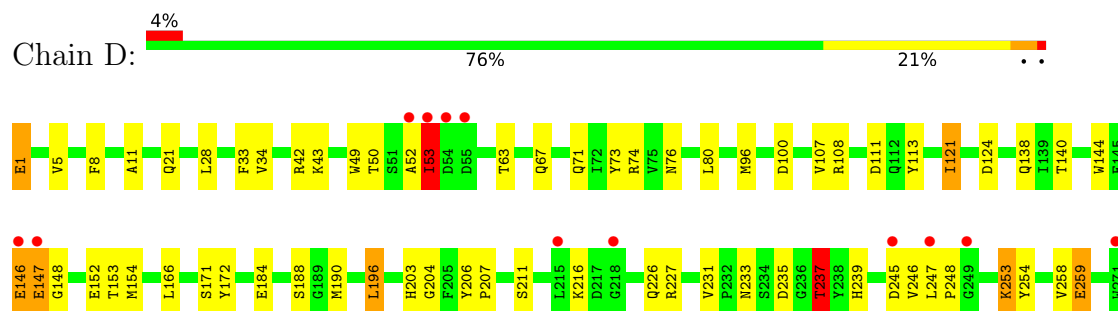
3 Residue-property plots [i](#)

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.

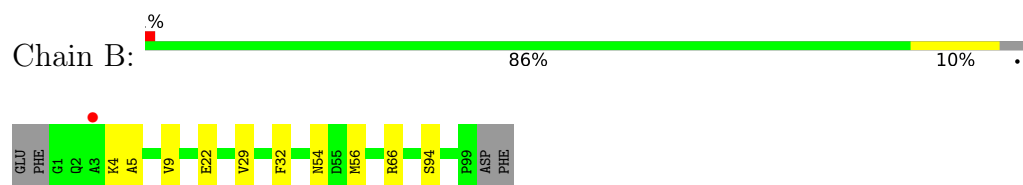
- Molecule 1: MHC class I antigen alpha chain



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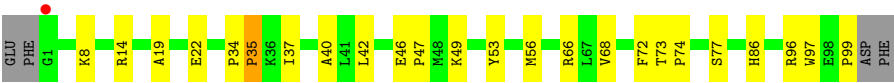


- Molecule 2: Beta2-microglobulin



- Molecule 2: Beta2-microglobulin

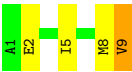




● Molecule 3: peptide of AIV



● Molecule 3: peptide of AIV



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	85.71Å 89.24Å 117.85Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.99 – 2.48 19.99 – 2.48	Depositor EDS
% Data completeness (in resolution range)	87.5 (19.99-2.48) 87.3 (19.99-2.48)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.08 (at 2.48Å)	Xtriage
Refinement program	PHENIX (1.20.1_4487: ???)	Depositor
R, R_{free}	0.203 , 0.275 0.212 , 0.268	Depositor DCC
R_{free} test set	1902 reflections (5.79%)	wwPDB-VP
Wilson B-factor (Å ²)	24.7	Xtriage
Anisotropy	0.167	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 31.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.018 for k,h,-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	6391	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.05% 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

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	0.85	2/2280 (0.1%)	0.95	8/3091 (0.3%)
1	D	0.46	0/2280	0.77	10/3091 (0.3%)
2	B	0.59	0/824	0.71	0/1122
2	E	1.13	1/824 (0.1%)	1.13	0/1122
3	C	0.91	0/61	0.98	0/82
3	F	0.97	0/61	0.90	0/82
All	All	0.75	3/6330 (0.0%)	0.89	18/8590 (0.2%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	232	PRO	C-O	-7.84	1.15	1.23
2	E	35	PRO	C-O	-6.08	1.16	1.24
1	A	208	ARG	C-O	-5.05	1.19	1.24

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	55	ASP	N-CA-C	-11.57	99.87	114.56
1	D	237	THR	CB-CA-C	8.14	124.85	110.36
1	D	253	LYS	N-CA-C	-6.25	105.43	114.12
1	D	52	ALA	CA-C-N	-5.97	112.86	120.98
1	D	52	ALA	C-N-CA	-5.97	112.86	120.98
1	A	247	LEU	N-CA-C	-5.94	102.30	109.72
1	A	52	ALA	CA-C-N	-5.83	110.27	120.29
1	A	52	ALA	C-N-CA	-5.83	110.27	120.29
1	D	53	ILE	N-CA-CB	5.72	117.34	110.31
1	D	147	GLU	N-CA-C	-5.66	101.08	109.11
1	D	248	PRO	CA-C-N	5.62	127.21	120.13
1	D	248	PRO	C-N-CA	5.62	127.21	120.13
1	D	146	GLU	CB-CA-C	5.51	119.53	110.88
1	D	247	LEU	N-CA-C	-5.26	101.90	109.48
1	A	232	PRO	N-CA-C	5.22	120.41	111.68

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	245	ASP	CA-C-N	-5.09	115.78	123.27
1	A	245	ASP	C-N-CA	-5.09	115.78	123.27
1	A	251	ARG	N-CA-C	-5.06	107.18	113.55

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	2221	0	2084	62	0
1	D	2221	0	2084	41	0
2	B	798	0	758	8	0
2	E	798	0	758	10	0
3	C	62	0	72	4	0
3	F	62	0	72	5	0
4	A	83	0	0	3	0
4	B	50	0	0	2	0
4	C	5	0	0	0	0
4	D	60	0	0	5	0
4	E	28	0	0	1	0
4	F	3	0	0	0	0
All	All	6391	0	5828	120	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 10.

All (120) 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:247:LEU:HD12	1:A:247:LEU:H	1.16	1.10
2:B:54:ASN:HB3	4:B:205:HOH:O	1.54	1.06
1:A:45:PRO:HB3	1:A:50:THR:HG21	1.37	1.05
1:D:53:ILE:HG22	1:D:172:TYR:HE2	1.29	0.93
1:D:53:ILE:HG22	1:D:172:TYR:CE2	2.03	0.92

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:104:ASP:OD1	1:A:106:SER:HB2	1.74	0.87
1:D:204:GLY:HA2	1:D:237:THR:HG23	1.56	0.87
1:A:247:LEU:HD13	1:A:250:ASP:HB2	1.57	0.84
1:D:5:VAL:HG23	1:D:166:LEU:HD13	1.61	0.82
1:A:53:ILE:HD11	1:A:169:TYR:HE1	1.44	0.80
2:B:56:MET:HG3	2:B:66:ARG:HG2	1.68	0.75
1:A:247:LEU:H	1:A:247:LEU:CD1	1.95	0.75
1:A:45:PRO:CB	1:A:50:THR:HG21	2.15	0.74
2:B:9:VAL:HG22	2:B:29:VAL:HG22	1.71	0.73
1:A:6:ARG:HH11	1:A:6:ARG:HG3	1.55	0.72
1:A:60:GLU:OE2	1:A:64:ARG:NE	2.23	0.71
1:D:42:ARG:HD3	1:D:63:THR:HG21	1.73	0.70
1:D:233:ASN:HB2	1:D:237:THR:HG22	1.75	0.69
1:D:1:GLU:N	4:D:302:HOH:O	2.26	0.67
1:A:247:LEU:HD12	1:A:247:LEU:N	2.01	0.67
1:A:102:LEU:HD11	1:A:108:ARG:HE	1.59	0.67
1:D:235:ASP:OD1	1:D:235:ASP:N	2.24	0.66
1:D:53:ILE:CG2	1:D:172:TYR:CE2	2.79	0.65
1:A:45:PRO:HB3	1:A:50:THR:CG2	2.21	0.65
1:D:153:THR:HG21	3:F:5:ILE:HG13	1.80	0.64
1:A:196:LEU:HD11	1:A:271:TRP:CZ3	2.32	0.64
1:A:55:ASP:HB3	1:A:58:TYR:HB2	1.80	0.63
1:D:1:GLU:N	4:D:304:HOH:O	2.31	0.63
1:A:43:LYS:HE2	1:A:62:ASN:HB3	1.80	0.62
1:A:53:ILE:HD11	1:A:169:TYR:CE1	2.32	0.62
2:B:56:MET:HE3	2:B:66:ARG:HH11	1.64	0.61
1:A:196:LEU:HD11	1:A:271:TRP:HZ3	1.64	0.61
1:D:50:THR:O	1:D:53:ILE:HG12	2.00	0.60
2:B:56:MET:HG3	2:B:66:ARG:CG	2.31	0.60
1:D:43:LYS:NZ	3:F:2:GLU:OE2	2.36	0.59
2:E:22:GLU:H	2:E:22:GLU:CD	2.12	0.57
2:E:37:ILE:HG13	2:E:86:HIS:HB2	1.87	0.56
1:A:251:ARG:HH21	1:A:271:TRP:HB3	1.69	0.56
1:A:103:GLU:H	1:A:103:GLU:CD	2.14	0.56
2:E:40:ALA:HA	4:E:206:HOH:O	2.06	0.55
1:A:196:LEU:HB3	1:A:246:VAL:HG23	1.89	0.55
1:A:251:ARG:NH2	1:A:271:TRP:C	2.65	0.55
1:A:153:THR:HG21	3:C:5:ILE:HG21	1.89	0.54
1:A:257:ARG:NH1	4:A:304:HOH:O	2.40	0.53
1:D:211:SER:HB3	1:D:259:GLU:HG2	1.91	0.53
1:A:169:TYR:OH	3:C:1:ALA:N	2.40	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:231:VAL:HG23	1:D:239:HIS:HB3	1.92	0.51
1:A:223:GLN:OE1	1:A:223:GLN:HA	2.09	0.51
1:D:216:LYS:HA	1:D:253:LYS:O	2.10	0.51
1:A:6:ARG:HG3	1:A:6:ARG:NH1	2.24	0.51
2:B:56:MET:HE3	2:B:66:ARG:NH1	2.26	0.51
1:D:63:THR:O	1:D:67:GLN:HG3	2.11	0.51
1:D:8:PHE:CE2	1:D:96:MET:HG3	2.46	0.51
1:A:4:SER:HB2	1:A:6:ARG:NH1	2.26	0.51
1:A:250:ASP:HA	1:A:253:LYS:HD3	1.93	0.50
1:D:107:VAL:O	1:D:108:ARG:HD3	2.12	0.49
1:D:246:VAL:HG11	1:D:254:TYR:CE2	2.47	0.49
1:A:201:ARG:HG3	1:A:241:TRP:HB3	1.95	0.49
1:A:102:LEU:HD12	1:A:108:ARG:HH21	1.77	0.48
1:A:251:ARG:HH21	1:A:271:TRP:C	2.21	0.48
1:D:43:LYS:NZ	4:D:314:HOH:O	2.46	0.48
2:E:14:ARG:O	2:E:14:ARG:NH1	2.42	0.48
1:A:134:ASP:H	1:A:137:ALA:HB3	1.78	0.48
1:A:6:ARG:HD3	1:A:96:MET:SD	2.54	0.48
2:E:34:PRO:HB2	2:E:35:PRO:HD2	1.95	0.48
2:B:54:ASN:ND2	4:B:205:HOH:O	2.47	0.47
1:A:43:LYS:CE	1:A:62:ASN:HB3	2.43	0.47
1:D:42:ARG:NH2	4:D:315:HOH:O	2.47	0.47
1:A:206:TYR:CD1	1:A:207:PRO:HA	2.50	0.46
1:D:140:THR:HG21	3:F:9:VAL:HG22	1.98	0.46
1:D:153:THR:CG2	3:F:5:ILE:HG13	2.46	0.46
1:D:144:TRP:O	1:D:148:GLY:HA2	2.15	0.46
1:A:247:LEU:CD1	1:A:247:LEU:N	2.71	0.46
1:A:215:LEU:HD23	1:A:220:VAL:HA	1.98	0.45
1:A:231:VAL:HG23	1:A:239:HIS:HB3	1.98	0.45
1:A:192:ALA:HB3	1:A:195:THR:CG2	2.47	0.45
1:D:11:ALA:HA	1:D:21:GLN:O	2.16	0.45
1:A:108:ARG:CZ	1:D:226:GLN:HE22	2.29	0.45
1:A:179:ARG:O	1:A:180:ARG:HG3	2.17	0.45
1:A:1:GLU:N	1:A:2:PRO:HD2	2.32	0.44
1:A:180:ARG:HH21	1:A:261:ALA:HB1	1.82	0.44
1:D:73:TYR:HA	1:D:76:ASN:HB2	1.99	0.44
1:A:166:LEU:C	1:A:166:LEU:HD23	2.42	0.44
2:E:19:ALA:HA	2:E:74:PRO:O	2.17	0.44
1:A:188:SER:OG	4:A:301:HOH:O	2.21	0.44
1:D:33:PHE:CD2	1:D:34:VAL:HG13	2.52	0.44
1:A:72:ILE:HD12	3:C:8:MET:HG3	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:150:PHE:CE1	3:C:5:ILE:HD12	2.54	0.43
1:D:71:GLN:OE1	1:D:74:ARG:NH1	2.52	0.43
1:A:200:CYS:HB2	1:A:214:TRP:CZ2	2.53	0.43
1:D:206:TYR:CD1	1:D:207:PRO:HA	2.54	0.43
1:D:28:LEU:HD12	1:D:49:TRP:CZ2	2.54	0.43
1:D:190:MET:O	1:D:196:LEU:HA	2.18	0.43
1:A:4:SER:HB2	1:A:6:ARG:HH12	1.83	0.43
2:B:5:ALA:O	2:B:32:PHE:HA	2.19	0.43
1:A:246:VAL:HG12	1:A:254:TYR:CZ	2.54	0.43
1:D:80:LEU:HD21	3:F:9:VAL:HG11	2.01	0.42
1:A:216:LYS:HB2	1:A:254:TYR:CE1	2.54	0.42
1:A:187:VAL:HG13	1:A:214:TRP:HH2	1.83	0.42
2:E:53:TYR:CD2	2:E:68:VAL:HG22	2.54	0.42
1:D:111:ASP:OD2	1:D:154:MET:HE2	2.18	0.42
2:E:56:MET:HE3	2:E:66:ARG:NH1	2.33	0.42
1:A:192:ALA:HB3	1:A:195:THR:HG23	2.02	0.42
1:A:6:ARG:NH1	1:A:100:ASP:OD1	2.52	0.42
1:A:50:THR:HG23	1:A:59:TRP:HZ2	1.85	0.42
1:A:182:ARG:HA	1:A:262:SER:OG	2.19	0.42
1:A:267:GLY:C	1:A:268:LEU:HD22	2.44	0.42
1:A:268:LEU:HD13	1:A:268:LEU:HA	1.91	0.41
2:E:46:GLU:HA	2:E:47:PRO:HD3	1.99	0.41
1:A:182:ARG:O	4:A:302:HOH:O	2.22	0.41
1:A:154:MET:O	1:A:158:LEU:HG	2.21	0.41
2:E:97:TRP:CH2	2:E:99:PRO:HB3	2.55	0.41
1:D:113:TYR:HB2	1:D:121:ILE:HG22	2.03	0.41
1:A:197:THR:HG23	1:A:245:ASP:OD1	2.21	0.40
1:D:184:GLU:HB2	1:D:203:HIS:HB2	2.02	0.40
1:D:100:ASP:HB2	1:D:108:ARG:HB2	2.03	0.40
1:D:152:GLU:HG2	4:D:306:HOH:O	2.20	0.40
1:A:199:SER:HA	1:A:243:THR:HG22	2.04	0.40
1:D:211:SER:O	1:D:258:VAL:HA	2.21	0.40
1:D:227:ARG:HE	1:D:227:ARG:HB3	1.79	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	269/271 (99%)	259 (96%)	10 (4%)	0	100	100
1	D	269/271 (99%)	255 (95%)	14 (5%)	0	100	100
2	B	97/103 (94%)	93 (96%)	4 (4%)	0	100	100
2	E	97/103 (94%)	93 (96%)	4 (4%)	0	100	100
3	C	7/9 (78%)	7 (100%)	0	0	100	100
3	F	7/9 (78%)	7 (100%)	0	0	100	100
All	All	746/766 (97%)	714 (96%)	32 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	235/235 (100%)	212 (90%)	23 (10%)	7	14
1	D	235/235 (100%)	222 (94%)	13 (6%)	19	37
2	B	86/90 (96%)	83 (96%)	3 (4%)	32	55
2	E	86/90 (96%)	79 (92%)	7 (8%)	11	21
3	C	6/6 (100%)	5 (83%)	1 (17%)	2	3
3	F	6/6 (100%)	4 (67%)	2 (33%)	0	0
All	All	654/662 (99%)	605 (92%)	49 (8%)	12	24

All (49) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	GLU
1	A	34	VAL
1	A	38	SER
1	A	44	GLU
1	A	50	THR
1	A	53	ILE
1	A	54	ASP
1	A	56	GLN
1	A	82	GLU
1	A	106	SER
1	A	112	GLN
1	A	188	SER
1	A	190	MET
1	A	191	GLU
1	A	194	LYS
1	A	197	THR
1	A	216	LYS
1	A	223	GLN
1	A	224	GLU
1	A	227	ARG
1	A	247	LEU
1	A	251	ARG
1	A	253	LYS
2	B	4	LYS
2	B	22	GLU
2	B	94	SER
3	C	8	MET
1	D	1	GLU
1	D	53	ILE
1	D	121	ILE
1	D	124	ASP
1	D	138	GLN
1	D	146	GLU
1	D	147	GLU
1	D	171	SER
1	D	188	SER
1	D	196	LEU
1	D	237	THR
1	D	245	ASP
1	D	259	GLU
2	E	8	LYS
2	E	42	LEU

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Mol	Chain	Res	Type
2	E	49	LYS
2	E	72	PHE
2	E	73	THR
2	E	77	SER
2	E	96	ARG
3	F	8	MET
3	F	9	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	160	ASN
1	A	203	HIS
1	A	226	GLN
2	B	2	GLN
2	B	26	ASN
1	D	57	GLN
1	D	62	ASN
1	D	86	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 oligosaccharides 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	271/271 (100%)	-0.05	7 (2%) 57 53	12, 24, 54, 77	0
1	D	271/271 (100%)	0.23	12 (4%) 39 35	16, 32, 67, 83	0
2	B	99/103 (96%)	-0.35	1 (1%) 79 77	12, 19, 35, 48	0
2	E	99/103 (96%)	0.07	1 (1%) 79 77	21, 28, 50, 59	0
3	C	9/9 (100%)	-0.01	0 100 100	16, 21, 30, 31	0
3	F	9/9 (100%)	-0.31	0 100 100	16, 18, 32, 33	0
All	All	758/766 (98%)	0.02	21 (2%) 55 51	12, 27, 60, 83	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	247	LEU	4.8
1	D	247	LEU	3.7
1	A	53	ILE	3.4
1	D	53	ILE	2.8
1	D	147	GLU	2.8
1	A	190	MET	2.7
1	D	52	ALA	2.6
1	D	218	GLY	2.5
1	D	249	GLY	2.4
1	A	54	ASP	2.4
1	D	271	TRP	2.4
1	D	55	ASP	2.3
1	D	215	LEU	2.2
2	E	1	GLY	2.2
2	B	3	ALA	2.1
1	D	146	GLU	2.1
1	A	250	ASP	2.1
1	A	126	ASP	2.1
1	D	54	ASP	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	245	ASP	2.1
1	A	51	SER	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 oligosaccharides 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.