



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 02:02 PM UTC

PDB ID : 3P39 / pdb_00003p39
Title : Crystal structure of the NS1 effector domain W182A mutant from influenza A/Vietnam/1203/2004 (H5N1) virus
Authors : Chen, S.; Xiao, Y.B.; Bricogne, G.; Sharff, A.J.; Hilgenfeld, R.
Deposited on : 2010-10-04
Resolution : 3.14 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

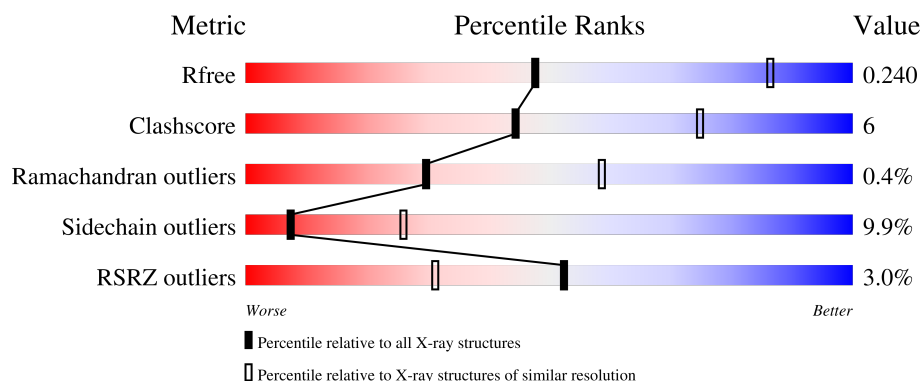
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.14 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	2351 (3.18-3.10)
Clashscore	190562	2452 (3.18-3.10)
Ramachandran outliers	187476	2324 (3.18-3.10)
Sidechain outliers	187428	2324 (3.18-3.10)
RSRZ outliers	180081	2351 (3.18-3.10)

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	147	0% 59% 22% •• 16%
1	B	147	3% 66% 14% • 16%
1	C	147	4% 65% 18% • 16%
1	D	147	3% 68% 16% • 14%
1	E	147	3% 67% 16% • 16%

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Mol	Chain	Length	Quality of chain
1	F	147	% 59% 22% • 16%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5900 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 Nonstructural protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	123	Total	C	N	O	S	0	0	0
			963	613	163	179	8			
1	B	123	Total	C	N	O	S	0	0	0
			963	613	163	179	8			
1	C	123	Total	C	N	O	S	0	0	0
			963	613	163	179	8			
1	D	126	Total	C	N	O	S	0	0	0
			985	628	167	182	8			
1	E	123	Total	C	N	O	S	0	0	0
			963	613	163	179	8			
1	F	123	Total	C	N	O	S	0	0	0
			963	613	163	179	8			

There are 30 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	69	GLY	-	expression tag	UNP D1LP63
A	70	PRO	-	expression tag	UNP D1LP63
A	71	LEU	-	expression tag	UNP D1LP63
A	72	GLY	-	expression tag	UNP D1LP63
A	182	ALA	TRP	engineered mutation	UNP D1LP63
B	69	GLY	-	expression tag	UNP D1LP63
B	70	PRO	-	expression tag	UNP D1LP63
B	71	LEU	-	expression tag	UNP D1LP63
B	72	GLY	-	expression tag	UNP D1LP63
B	182	ALA	TRP	engineered mutation	UNP D1LP63
C	69	GLY	-	expression tag	UNP D1LP63
C	70	PRO	-	expression tag	UNP D1LP63
C	71	LEU	-	expression tag	UNP D1LP63
C	72	GLY	-	expression tag	UNP D1LP63
C	182	ALA	TRP	engineered mutation	UNP D1LP63
D	69	GLY	-	expression tag	UNP D1LP63
D	70	PRO	-	expression tag	UNP D1LP63

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Chain	Residue	Modelled	Actual	Comment	Reference
D	71	LEU	-	expression tag	UNP D1LP63
D	72	GLY	-	expression tag	UNP D1LP63
D	182	ALA	TRP	engineered mutation	UNP D1LP63
E	69	GLY	-	expression tag	UNP D1LP63
E	70	PRO	-	expression tag	UNP D1LP63
E	71	LEU	-	expression tag	UNP D1LP63
E	72	GLY	-	expression tag	UNP D1LP63
E	182	ALA	TRP	engineered mutation	UNP D1LP63
F	69	GLY	-	expression tag	UNP D1LP63
F	70	PRO	-	expression tag	UNP D1LP63
F	71	LEU	-	expression tag	UNP D1LP63
F	72	GLY	-	expression tag	UNP D1LP63
F	182	ALA	TRP	engineered mutation	UNP D1LP63

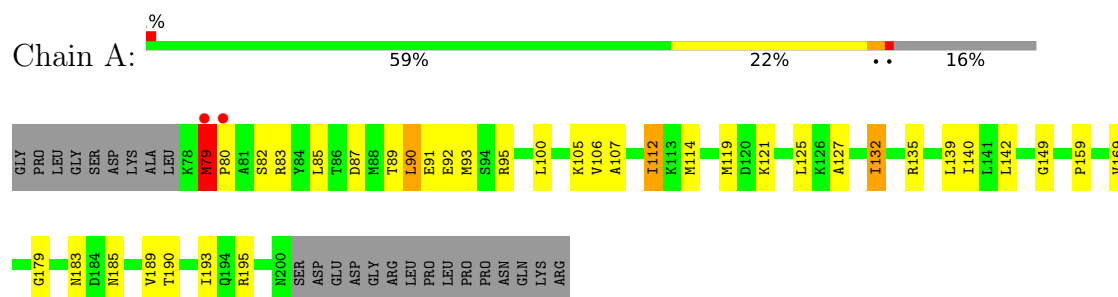
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	15	Total O 15 15	0	0
2	B	16	Total O 16 16	0	0
2	C	15	Total O 15 15	0	0
2	D	23	Total O 23 23	0	0
2	E	17	Total O 17 17	0	0
2	F	14	Total O 14 14	0	0

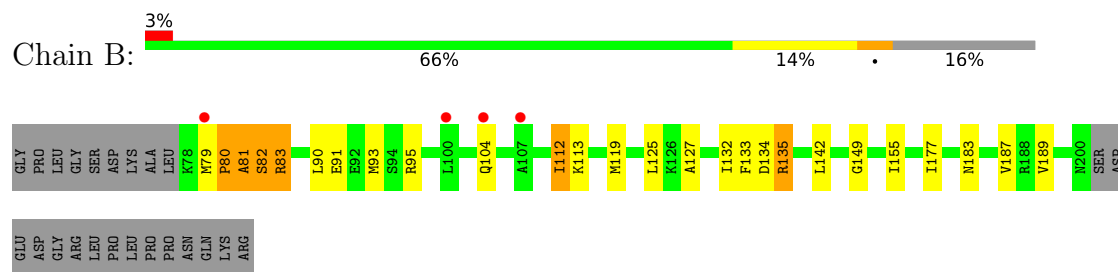
3 Residue-property plots [i](#)

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.

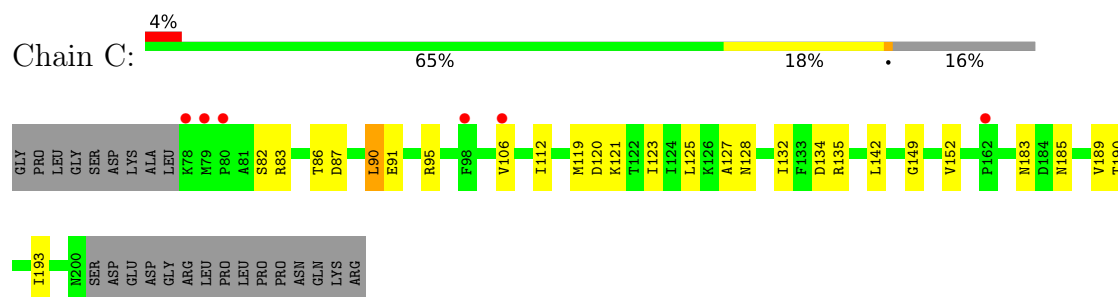
- Molecule 1: Nonstructural protein 1



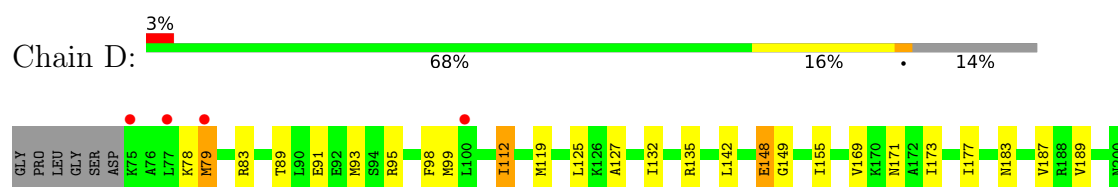
- Molecule 1: Nonstructural protein 1



- Molecule 1: Nonstructural protein 1



- Molecule 1: Nonstructural protein 1



SER
ASP
GLU
ASP
GLY
ASP
GLY
ARG
LEU
PRO
PRO
PRO
ASN
GLN
LYS
ARG

● Molecule 1: Nonstructural protein 1



GLY
PRO
LEU
GLY
SER
ASP
LYS
ALA
LEU
K78
M79
P80
A81
S82
R83
D87
L90
E91
M101
Q104
K105
V106
I112
K113
M114
M119
D120
K121
L125
K126
A127
D134
R135
L136
E137
L142
M183
D184
M185
V189
T190
I193
K200
SER
ASP
GLU
ASP

GLY
ARG
LEU
PRO
GLY
LEU
PRO
ASN
LYS
ARG

● Molecule 1: Nonstructural protein 1



GLY
PRO
LEU
GLY
SER
ASP
LYS
ALA
LEU
K78
M79
P80
A81
S82
R83
T86
D87
L90
E91
R95
Q104
K105
V106
I112
K113
M114
M119
D120
K121
T122
I123
I124
L125
K126
A127
M128
I132
F133
D134
R135
L142
R143
E147
E148
G149
V152
G153
E154
I155

L176
G179
N183
D184
N185
V189
T190
K200
SER
ASP
GLU
ASP
GLY
ARG
LEU
PRO
LEU
PRO
PRO
ASN
GLN
LYS
ARG

4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	113.49Å 113.49Å 197.44Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	37.30 – 3.14 37.30 – 3.14	Depositor EDS
% Data completeness (in resolution range)	94.6 (37.30-3.14) 94.7 (37.30-3.14)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.22 (at 3.12Å)	Xtriage
Refinement program	BUSTER-TNT BUSTER 2.11.0, BUSTER 2.11.0	Depositor
R, R_{free}	0.207 , 0.237 0.213 , 0.240	Depositor DCC
R_{free} test set	1128 reflections (5.13%)	wwPDB-VP
Wilson B-factor (Å ²)	47.7	Xtriage
Anisotropy	0.551	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 50.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	5900	wwPDB-VP
Average B, all atoms (Å ²)	47.0	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 27.26 % 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 2.2702e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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	1/978 (0.1%)	1.34	3/1320 (0.2%)
1	B	0.77	0/978	1.34	5/1320 (0.4%)
1	C	0.82	0/978	1.30	2/1320 (0.2%)
1	D	0.78	0/1000	1.31	2/1349 (0.1%)
1	E	0.87	0/978	1.34	2/1320 (0.2%)
1	F	0.82	1/978 (0.1%)	1.30	1/1320 (0.1%)
All	All	0.82	2/5890 (0.0%)	1.32	15/7949 (0.2%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	132	ILE	CG1-CD1	5.58	1.73	1.51
1	A	132	ILE	CG1-CD1	5.03	1.71	1.51

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	82	SER	N-CA-C	7.69	120.24	110.24
1	C	120	ASP	CA-CB-CG	6.80	119.40	112.60
1	B	81	ALA	CA-C-N	6.22	129.96	120.82
1	B	81	ALA	C-N-CA	6.22	129.96	120.82
1	A	127	ALA	N-CA-C	6.00	118.08	109.14
1	C	127	ALA	N-CA-C	5.95	117.82	108.96
1	D	127	ALA	N-CA-C	5.87	117.70	108.96
1	B	80	PRO	N-CA-C	5.86	124.54	112.47
1	A	140	ILE	N-CA-C	-5.83	106.79	111.81
1	E	127	ALA	N-CA-C	5.70	117.46	108.96
1	B	127	ALA	N-CA-C	5.67	117.40	108.96
1	F	127	ALA	N-CA-C	5.43	117.05	108.96
1	E	106	VAL	CB-CA-C	5.42	116.71	111.23
1	A	79	MET	N-CA-C	5.16	121.20	109.81
1	D	148	GLU	CB-CG-CD	5.07	121.22	112.60

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	963	0	990	19	0
1	B	963	0	990	13	0
1	C	963	0	990	10	0
1	D	985	0	1019	12	0
1	E	963	0	990	8	0
1	F	963	0	990	14	0
2	A	15	0	0	1	0
2	B	16	0	0	1	0
2	C	15	0	0	0	0
2	D	23	0	0	0	0
2	E	17	0	0	0	0
2	F	14	0	0	0	0
All	All	5900	0	5969	71	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 6.

All (71) 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:E:119:MET:HG2	1:E:183:ASN:HB3	1.74	0.68
1:F:119:MET:HG2	1:F:183:ASN:HB3	1.76	0.68
1:D:119:MET:HG2	1:D:183:ASN:HB3	1.77	0.67
1:A:90:LEU:HD11	1:B:93:MET:HE2	1.76	0.66
1:A:119:MET:HG2	1:A:183:ASN:HB3	1.77	0.66
1:B:119:MET:HG2	1:B:183:ASN:HB3	1.76	0.66
1:C:119:MET:HG2	1:C:183:ASN:HB3	1.77	0.65
1:A:85:LEU:HD21	1:B:83:ARG:HH21	1.66	0.61
1:B:112:ILE:HG23	1:B:155:ILE:HG12	1.82	0.60
1:A:159:PRO:HD2	1:D:98:PHE:CZ	2.39	0.57
1:D:169:VAL:O	1:D:173:ILE:HG12	2.05	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:112:ILE:HG23	1:D:155:ILE:HG12	1.86	0.56
1:C:87:ASP:OD2	1:C:190:THR:HG21	2.06	0.54
1:B:135:ARG:HD3	2:B:63:HOH:O	2.08	0.54
1:A:87:ASP:OD2	1:A:190:THR:HG21	2.08	0.54
1:A:190:THR:HG22	1:A:193:ILE:HD12	1.88	0.54
1:F:114:MET:HE1	1:F:179:GLY:HA3	1.89	0.54
1:C:90:LEU:HD11	1:D:93:MET:HE2	1.90	0.53
1:E:87:ASP:OD2	1:E:190:THR:HG21	2.08	0.53
1:D:173:ILE:HG22	1:D:177:ILE:HD12	1.91	0.52
1:F:87:ASP:OD2	1:F:190:THR:HG21	2.09	0.52
1:E:121:LYS:H	1:E:185:ASN:ND2	2.08	0.52
1:E:121:LYS:H	1:E:185:ASN:HD22	1.58	0.51
1:A:114:MET:HE1	1:A:179:GLY:HA3	1.91	0.51
1:B:125:LEU:HG	1:B:142:LEU:HD21	1.92	0.51
1:C:125:LEU:HG	1:C:142:LEU:HD21	1.93	0.50
1:A:125:LEU:HG	1:A:142:LEU:HD21	1.94	0.50
1:D:125:LEU:HG	1:D:142:LEU:HD21	1.93	0.50
1:D:177:ILE:HG23	1:D:187:VAL:HG11	1.94	0.50
1:F:125:LEU:HG	1:F:142:LEU:HD21	1.93	0.49
1:B:177:ILE:HG23	1:B:187:VAL:HG11	1.94	0.49
1:C:121:LYS:H	1:C:185:ASN:HD22	1.60	0.48
1:D:177:ILE:HG23	1:D:187:VAL:CG1	2.44	0.48
1:A:121:LYS:H	1:A:185:ASN:ND2	2.12	0.47
1:B:133:PHE:O	1:B:135:ARG:HG2	2.15	0.47
1:C:190:THR:HG22	1:C:193:ILE:HD12	1.96	0.47
1:E:125:LEU:HG	1:E:142:LEU:HD21	1.96	0.47
1:F:121:LYS:H	1:F:185:ASN:HD22	1.62	0.47
1:B:177:ILE:HG23	1:B:187:VAL:CG1	2.45	0.47
1:E:190:THR:HG22	1:E:193:ILE:HD12	1.97	0.47
1:A:121:LYS:H	1:A:185:ASN:HD22	1.62	0.47
1:C:121:LYS:H	1:C:185:ASN:ND2	2.12	0.47
1:F:121:LYS:H	1:F:185:ASN:ND2	2.13	0.46
1:A:139:LEU:HD22	1:A:169:VAL:HG13	1.98	0.46
1:A:89:THR:OG1	1:A:92:GLU:HG3	2.16	0.45
1:A:195:ARG:HD3	2:A:47:HOH:O	2.17	0.45
1:E:104:GLN:HG2	1:E:113:LYS:HG2	1.98	0.44
1:E:190:THR:HG23	1:E:193:ILE:H	1.83	0.44
1:F:142:LEU:HD22	1:F:155:ILE:HD12	2.00	0.44
1:D:79:MET:HE3	1:D:79:MET:HA	2.00	0.43
1:A:90:LEU:HD11	1:B:93:MET:CE	2.46	0.43
1:A:95:ARG:NH2	1:A:149:GLY:O	2.51	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:95:ARG:NH2	1:D:149:GLY:O	2.51	0.43
1:A:89:THR:O	1:A:93:MET:HG3	2.20	0.42
1:C:95:ARG:NH2	1:C:149:GLY:O	2.52	0.42
1:A:79:MET:HA	1:A:80:PRO:HD3	1.84	0.42
1:B:95:ARG:NH2	1:B:149:GLY:O	2.54	0.41
1:B:79:MET:HA	1:B:80:PRO:HD3	1.85	0.41
1:A:105:LYS:HE3	1:A:107:ALA:HB2	2.03	0.41
1:C:123:ILE:HD12	1:C:152:VAL:HG21	2.02	0.41
1:D:89:THR:O	1:D:93:MET:HG3	2.21	0.41
1:F:79:MET:HA	1:F:80:PRO:HD3	1.99	0.41
1:F:104:GLN:HG2	1:F:113:LYS:HG2	2.02	0.41
1:A:105:LYS:HB3	1:A:112:ILE:HG23	2.02	0.41
1:F:95:ARG:NH2	1:F:149:GLY:O	2.54	0.41
1:F:143:ARG:HG2	1:F:154:GLU:HG3	2.02	0.41
1:B:104:GLN:HG2	1:B:113:LYS:HG2	2.02	0.41
1:F:112:ILE:CD1	1:F:176:LEU:HB2	2.51	0.41
1:F:123:ILE:HD12	1:F:152:VAL:HG21	2.03	0.41
1:C:86:THR:HB	1:C:128:ASN:OD1	2.21	0.40
1:F:86:THR:HB	1:F:128:ASN:OD1	2.20	0.40

There are no symmetry-related clashes.

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	121/147 (82%)	116 (96%)	4 (3%)	1 (1%)	16	44
1	B	121/147 (82%)	117 (97%)	3 (2%)	1 (1%)	16	44
1	C	121/147 (82%)	116 (96%)	5 (4%)	0	100	100
1	D	124/147 (84%)	121 (98%)	3 (2%)	0	100	100
1	E	121/147 (82%)	116 (96%)	4 (3%)	1 (1%)	16	44

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	F	121/147 (82%)	118 (98%)	3 (2%)	0	100	100
All	All	729/882 (83%)	704 (97%)	22 (3%)	3 (0%)	30	59

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	81	ALA
1	E	79	MET
1	A	79	MET

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	106/126 (84%)	96 (91%)	10 (9%)	8	28
1	B	106/126 (84%)	97 (92%)	9 (8%)	10	31
1	C	106/126 (84%)	96 (91%)	10 (9%)	8	28
1	D	108/126 (86%)	97 (90%)	11 (10%)	7	25
1	E	106/126 (84%)	95 (90%)	11 (10%)	7	24
1	F	106/126 (84%)	94 (89%)	12 (11%)	5	21
All	All	638/756 (84%)	575 (90%)	63 (10%)	7	26

All (63) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	82	SER
1	A	83	ARG
1	A	90	LEU
1	A	91	GLU
1	A	100	LEU
1	A	106	VAL
1	A	112	ILE
1	A	132	ILE

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Mol	Chain	Res	Type
1	A	135	ARG
1	A	189	VAL
1	B	82	SER
1	B	83	ARG
1	B	90	LEU
1	B	91	GLU
1	B	112	ILE
1	B	132	ILE
1	B	134	ASP
1	B	135	ARG
1	B	189	VAL
1	C	82	SER
1	C	83	ARG
1	C	90	LEU
1	C	91	GLU
1	C	106	VAL
1	C	112	ILE
1	C	132	ILE
1	C	134	ASP
1	C	135	ARG
1	C	189	VAL
1	D	78	LYS
1	D	79	MET
1	D	83	ARG
1	D	91	GLU
1	D	99	MET
1	D	112	ILE
1	D	132	ILE
1	D	135	ARG
1	D	148	GLU
1	D	171	ASN
1	D	189	VAL
1	E	79	MET
1	E	82	SER
1	E	83	ARG
1	E	90	LEU
1	E	91	GLU
1	E	101	MET
1	E	112	ILE
1	E	134	ASP
1	E	135	ARG
1	E	137	GLU

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Mol	Chain	Res	Type
1	E	189	VAL
1	F	82	SER
1	F	83	ARG
1	F	90	LEU
1	F	91	GLU
1	F	106	VAL
1	F	112	ILE
1	F	121	LYS
1	F	132	ILE
1	F	134	ASP
1	F	135	ARG
1	F	147	GLU
1	F	189	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	116	GLN
1	A	185	ASN
1	C	116	GLN
1	C	185	ASN
1	D	104	GLN
1	E	116	GLN
1	E	185	ASN
1	F	185	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, 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	123/147 (83%)	-0.01	2 (1%) 70 49	23, 40, 62, 79	0
1	B	123/147 (83%)	0.03	4 (3%) 49 28	26, 50, 74, 82	0
1	C	123/147 (83%)	0.12	6 (4%) 35 18	22, 44, 72, 82	0
1	D	126/147 (85%)	0.04	4 (3%) 50 29	24, 43, 68, 81	0
1	E	123/147 (83%)	0.18	5 (4%) 41 22	22, 46, 71, 81	0
1	F	123/147 (83%)	0.20	1 (0%) 82 65	28, 53, 78, 84	0
All	All	741/882 (84%)	0.09	22 (2%) 52 31	22, 46, 74, 84	0

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	78	LYS	3.4
1	E	79	MET	3.0
1	A	80	PRO	2.8
1	B	107	ALA	2.8
1	E	114	MET	2.6
1	D	75	LYS	2.5
1	D	100	LEU	2.4
1	A	79	MET	2.3
1	D	77	LEU	2.3
1	E	80	PRO	2.3
1	C	79	MET	2.3
1	D	79	MET	2.3
1	C	80	PRO	2.3
1	B	104	GLN	2.2
1	C	78	LYS	2.2
1	C	106	VAL	2.2
1	F	79	MET	2.2
1	C	98	PHE	2.1
1	B	79	MET	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	162	PRO	2.1
1	E	81	ALA	2.0
1	B	100	LEU	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.