



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 02:44 PM UTC

PDB ID : 2YGC / pdb_00002ygc
Title : Structure of vaccinia virus D13 scaffolding protein
Authors : Bahar, M.W.; Graham, S.C.; Stuart, D.I.; Grimes, J.M.
Deposited on : 2011-04-13
Resolution : 3.02 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
Mogul	:	2022.3.0, CSD as543be (2022)
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

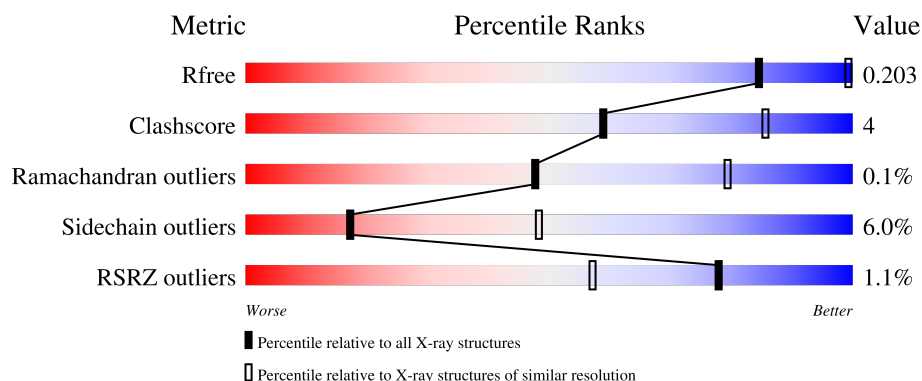
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.02 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	3131 (3.04-3.00)
Clashscore	190562	3444 (3.04-3.00)
Ramachandran outliers	187476	3319 (3.04-3.00)
Sidechain outliers	187428	3322 (3.04-3.00)
RSRZ outliers	180081	3130 (3.04-3.00)

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	569	% 80% 15% • 5%
1	B	569	% 80% 13% • 5%
1	C	569	% 77% 17% • 5%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 12992 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 RIFAMPICIN RESISTANCE PROTEIN.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	542	Total	C	N	O	S	Se	0	1	0
			4308	2753	703	840	4	8			
1	B	542	Total	C	N	O	S	Se	0	0	0
			4300	2748	700	840	4	8			
1	C	542	Total	C	N	O	S	Se	0	1	0
			4305	2752	700	840	4	9			

There are 60 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-17	MSE	-	expression tag	UNP P68440
A	-16	ALA	-	expression tag	UNP P68440
A	-15	HIS	-	expression tag	UNP P68440
A	-14	HIS	-	expression tag	UNP P68440
A	-13	HIS	-	expression tag	UNP P68440
A	-12	HIS	-	expression tag	UNP P68440
A	-11	HIS	-	expression tag	UNP P68440
A	-10	HIS	-	expression tag	UNP P68440
A	-9	SER	-	expression tag	UNP P68440
A	-8	SER	-	expression tag	UNP P68440
A	-7	GLY	-	expression tag	UNP P68440
A	-6	LEU	-	expression tag	UNP P68440
A	-5	GLU	-	expression tag	UNP P68440
A	-4	VAL	-	expression tag	UNP P68440
A	-3	LEU	-	expression tag	UNP P68440
A	-2	PHE	-	expression tag	UNP P68440
A	-1	GLN	-	expression tag	UNP P68440
A	0	GLY	-	expression tag	UNP P68440
A	1	PRO	-	expression tag	UNP P68440
A	513	GLY	ASP	engineered mutation	UNP P68440
B	-17	MSE	-	expression tag	UNP P68440
B	-16	ALA	-	expression tag	UNP P68440
B	-15	HIS	-	expression tag	UNP P68440

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-14	HIS	-	expression tag	UNP P68440
B	-13	HIS	-	expression tag	UNP P68440
B	-12	HIS	-	expression tag	UNP P68440
B	-11	HIS	-	expression tag	UNP P68440
B	-10	HIS	-	expression tag	UNP P68440
B	-9	SER	-	expression tag	UNP P68440
B	-8	SER	-	expression tag	UNP P68440
B	-7	GLY	-	expression tag	UNP P68440
B	-6	LEU	-	expression tag	UNP P68440
B	-5	GLU	-	expression tag	UNP P68440
B	-4	VAL	-	expression tag	UNP P68440
B	-3	LEU	-	expression tag	UNP P68440
B	-2	PHE	-	expression tag	UNP P68440
B	-1	GLN	-	expression tag	UNP P68440
B	0	GLY	-	expression tag	UNP P68440
B	1	PRO	-	expression tag	UNP P68440
B	513	GLY	ASP	engineered mutation	UNP P68440
C	-17	MSE	-	expression tag	UNP P68440
C	-16	ALA	-	expression tag	UNP P68440
C	-15	HIS	-	expression tag	UNP P68440
C	-14	HIS	-	expression tag	UNP P68440
C	-13	HIS	-	expression tag	UNP P68440
C	-12	HIS	-	expression tag	UNP P68440
C	-11	HIS	-	expression tag	UNP P68440
C	-10	HIS	-	expression tag	UNP P68440
C	-9	SER	-	expression tag	UNP P68440
C	-8	SER	-	expression tag	UNP P68440
C	-7	GLY	-	expression tag	UNP P68440
C	-6	LEU	-	expression tag	UNP P68440
C	-5	GLU	-	expression tag	UNP P68440
C	-4	VAL	-	expression tag	UNP P68440
C	-3	LEU	-	expression tag	UNP P68440
C	-2	PHE	-	expression tag	UNP P68440
C	-1	GLN	-	expression tag	UNP P68440
C	0	GLY	-	expression tag	UNP P68440
C	1	PRO	-	expression tag	UNP P68440
C	513	GLY	ASP	engineered mutation	UNP P68440

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	34	Total O 34 34	0	0

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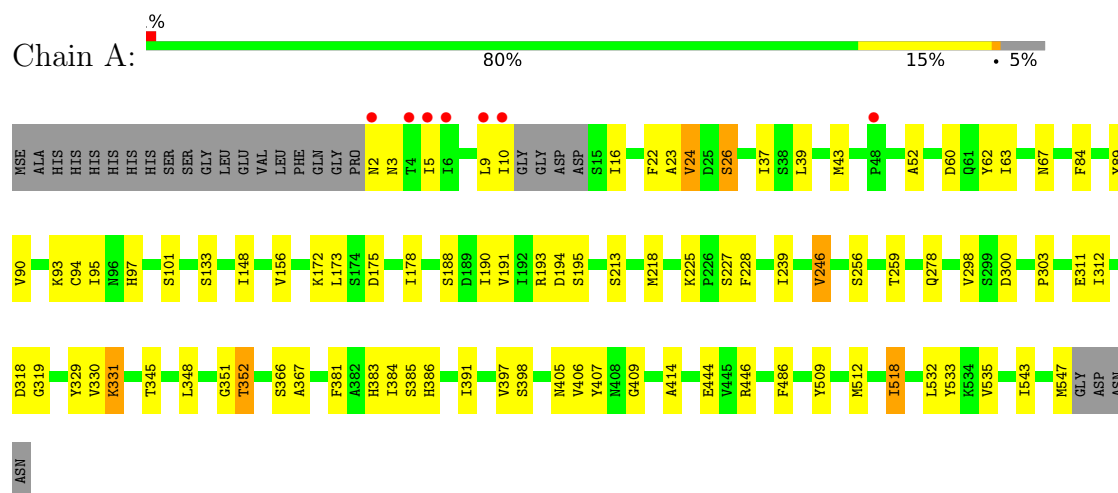
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	28	Total	O	0	0
			28	28		
2	C	17	Total	O	0	0
			17	17		

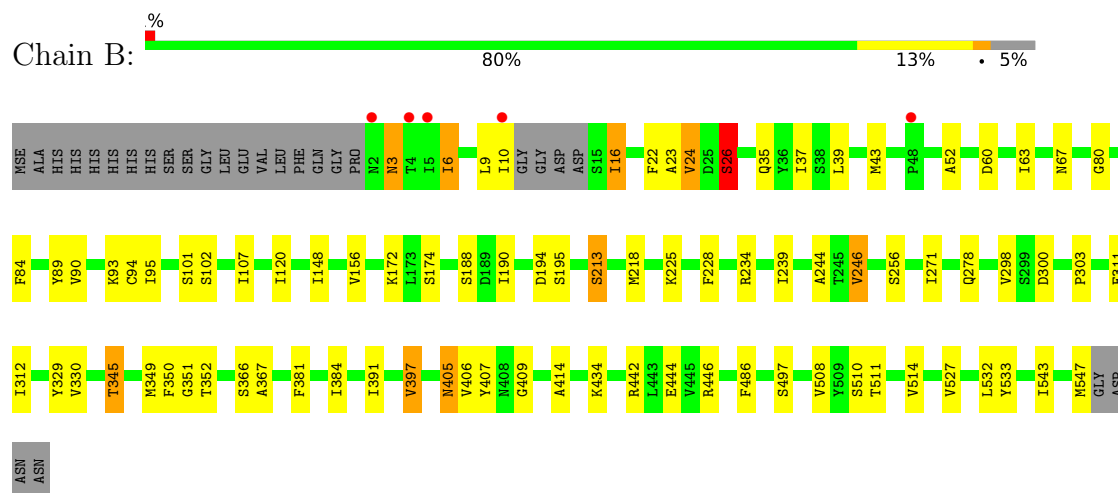
3 Residue-property plots [i](#)

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.

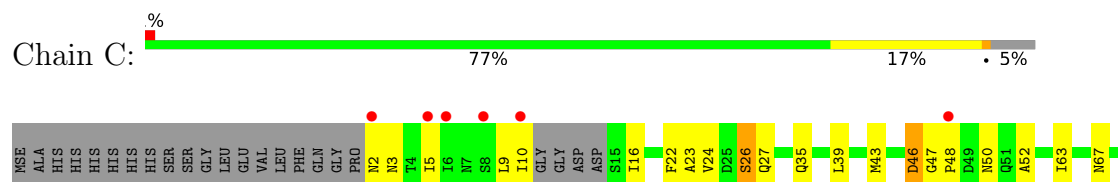
• Molecule 1: RIFAMPICIN RESISTANCE PROTEIN

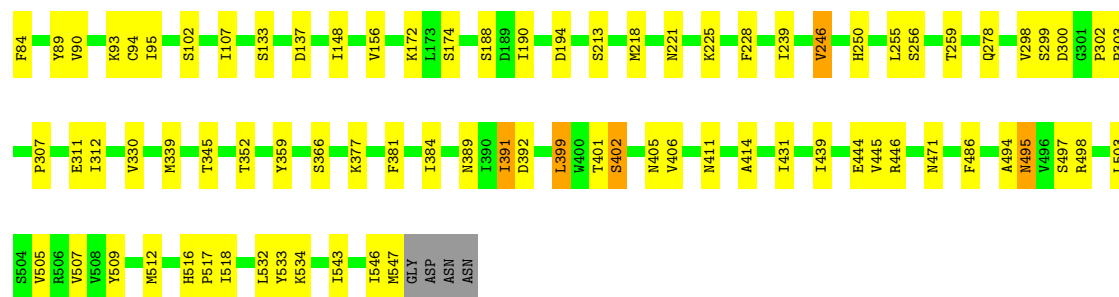


• Molecule 1: RIFAMPICIN RESISTANCE PROTEIN



• Molecule 1: RIFAMPICIN RESISTANCE PROTEIN





4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α , β , γ	190.58Å 190.58Å 252.96Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	39.28 – 3.02 39.28 – 3.02	Depositor EDS
% Data completeness (in resolution range)	99.7 (39.28-3.02) 99.6 (39.28-3.02)	Depositor EDS
R_{merge}	0.21	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.12 (at 3.01Å)	Xtriage
Refinement program	BUSTER 2.11.1	Depositor
R, R_{free}	0.170 , 0.203 0.177 , 0.203	Depositor DCC
R_{free} test set	2717 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	65.0	Xtriage
Anisotropy	0.352	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 70.1	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	0.95	EDS
Total number of atoms	12992	wwPDB-VP
Average B, all atoms (Å ²)	62.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.38% 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	0.77	0/4396	1.20	17/5967 (0.3%)
1	B	0.76	0/4385	1.18	14/5953 (0.2%)
1	C	0.76	0/4393	1.18	16/5963 (0.3%)
All	All	0.77	0/13174	1.19	47/17883 (0.3%)

There are no bond length outliers.

All (47) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	246	VAL	N-CA-CB	6.81	119.17	111.21
1	B	405	ASN	CA-C-N	6.56	131.52	120.64
1	B	405	ASN	C-N-CA	6.56	131.52	120.64
1	B	414	ALA	CA-C-N	6.48	128.86	120.44
1	B	414	ALA	C-N-CA	6.48	128.86	120.44
1	C	414	ALA	CA-C-N	6.23	128.54	120.44
1	C	414	ALA	C-N-CA	6.23	128.54	120.44
1	A	405	ASN	CA-C-N	6.17	130.41	120.30
1	A	405	ASN	C-N-CA	6.17	130.41	120.30
1	A	414	ALA	CA-C-N	6.17	128.46	120.44
1	A	414	ALA	C-N-CA	6.17	128.46	120.44
1	A	300	ASP	CA-CB-CG	5.96	118.56	112.60
1	A	175	ASP	CA-CB-CG	5.95	118.55	112.60
1	C	405	ASN	CA-C-N	5.91	130.45	120.64
1	C	405	ASN	C-N-CA	5.91	130.45	120.64
1	B	300	ASP	CA-CB-CG	5.72	118.32	112.60
1	C	194	ASP	CA-CB-CG	5.62	118.22	112.60
1	C	46	ASP	CA-CB-CG	5.59	118.19	112.60
1	A	67	ASN	CA-CB-CG	5.55	118.15	112.60
1	C	67	ASN	CA-CB-CG	5.55	118.15	112.60
1	B	514	VAL	N-CA-C	-5.54	105.80	113.00
1	B	195	SER	CA-C-N	5.50	127.92	120.38
1	B	195	SER	C-N-CA	5.50	127.92	120.38
1	B	89	TYR	CA-C-N	5.48	128.25	120.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	89	TYR	C-N-CA	5.48	128.25	120.53
1	A	195	SER	CA-C-N	5.41	127.53	120.28
1	A	195	SER	C-N-CA	5.41	127.53	120.28
1	C	300	ASP	CA-CB-CG	5.40	118.00	112.60
1	A	194	ASP	CA-CB-CG	5.39	117.99	112.60
1	A	133	SER	CA-C-N	5.37	127.43	120.44
1	A	133	SER	C-N-CA	5.37	127.43	120.44
1	C	495	ASN	CA-C-N	5.37	128.11	120.53
1	C	495	ASN	C-N-CA	5.37	128.11	120.53
1	B	246	VAL	N-CA-CB	5.34	117.46	111.21
1	B	67	ASN	CA-CB-CG	5.30	117.90	112.60
1	A	246	VAL	N-CA-CB	5.26	117.37	111.21
1	B	194	ASP	CA-CB-CG	5.26	117.86	112.60
1	C	89	TYR	CA-C-N	5.18	127.83	120.53
1	C	89	TYR	C-N-CA	5.18	127.83	120.53
1	A	352	THR	N-CA-C	-5.16	107.16	113.50
1	B	80	GLY	N-CA-C	5.06	116.86	110.38
1	C	221	ASN	CA-CB-CG	5.05	117.65	112.60
1	C	174	SER	CA-C-N	5.05	129.72	122.40
1	C	174	SER	C-N-CA	5.05	129.72	122.40
1	A	318	ASP	CA-CB-CG	5.04	117.64	112.60
1	A	89	TYR	CA-C-N	5.00	127.58	120.53
1	A	89	TYR	C-N-CA	5.00	127.58	120.53

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	4308	0	4272	40	0
1	B	4300	0	4259	43	0
1	C	4305	0	4268	48	0
2	A	34	0	0	1	0
2	B	28	0	0	0	0
2	C	17	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	12992	0	12799	114	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 4.

All (114) 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:C:431:ILE:HD13	1:C:512[B]:MSE:HE3	1.24	1.10
1:C:512[B]:MSE:HE2	1:C:518:ILE:HG23	1.55	0.89
1:C:431:ILE:HD13	1:C:512[B]:MSE:CE	2.04	0.87
1:A:218:MSE:HE3	1:B:6:ILE:HG23	1.59	0.85
1:A:218:MSE:CE	1:B:6:ILE:HG23	2.15	0.75
1:B:311:GLU:O	1:B:345:THR:HB	1.87	0.75
1:A:173:LEU:HD22	1:A:178:ILE:HG13	1.69	0.74
1:C:431:ILE:CD1	1:C:512[B]:MSE:HE3	2.14	0.73
1:C:94:CYS:HB3	1:C:190:ILE:HG21	1.73	0.70
1:A:172:LYS:HE2	1:B:22:PHE:O	1.92	0.69
1:A:218:MSE:SE	1:B:6:ILE:HD12	2.44	0.68
1:B:102:SER:HB2	1:B:107:ILE:HD11	1.76	0.68
1:A:84:PHE:HB3	1:A:148:ILE:HG21	1.78	0.66
1:B:63:ILE:HD13	1:B:218:MSE:SE	2.46	0.66
1:A:311:GLU:O	1:A:345:THR:HB	1.96	0.65
1:A:94:CYS:HB3	1:A:190:ILE:HG21	1.77	0.65
1:B:172:LYS:HE2	1:C:22:PHE:O	1.97	0.65
1:C:24:VAL:HG12	1:C:26:SER:H	1.61	0.64
1:C:311:GLU:O	1:C:345:THR:HB	1.98	0.64
1:A:24:VAL:HG11	1:C:486:PHE:HA	1.80	0.64
1:C:43:MSE:HE1	1:C:52:ALA:HB1	1.80	0.64
1:A:60:ASP:HB3	1:A:63:ILE:HG12	1.80	0.63
1:C:84:PHE:HB3	1:C:148:ILE:HG21	1.81	0.63
1:B:442:ARG:HB2	1:B:508:VAL:HB	1.80	0.63
1:A:43:MSE:HE1	1:A:52:ALA:HB1	1.82	0.62
1:B:84:PHE:HB3	1:B:148:ILE:HG21	1.82	0.61
1:B:94:CYS:HB3	1:B:190:ILE:HG21	1.82	0.61
1:C:63:ILE:HD13	1:C:218:MSE:SE	2.50	0.61
1:B:43:MSE:HE1	1:B:52:ALA:HB1	1.83	0.61
1:A:22:PHE:O	1:C:172:LYS:HE2	2.02	0.59
1:C:533:TYR:CD1	1:C:543:ILE:HD12	2.38	0.58
1:B:486:PHE:HA	1:C:24:VAL:HG11	1.86	0.57
1:A:193[B]:ARG:HG2	1:A:193[B]:ARG:HH11	1.71	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:193[B]:ARG:HH11	1:A:193[B]:ARG:CG	2.20	0.55
1:A:24:VAL:HG12	1:A:26:SER:H	1.71	0.55
1:A:95:ILE:HD12	1:A:156:VAL:HG11	1.90	0.54
1:C:47:GLY:HA2	1:C:50:ASN:H	1.73	0.53
1:A:259:THR:HG21	1:A:509:TYR:HE1	1.74	0.53
1:B:24:VAL:HG12	1:B:26:SER:H	1.73	0.53
1:C:95:ILE:HD12	1:C:156:VAL:HG11	1.91	0.53
1:B:95:ILE:HD12	1:B:156:VAL:HG11	1.91	0.52
1:C:381:PHE:HB3	1:C:384:ILE:HD11	1.91	0.52
1:A:329:TYR:CZ	1:A:331:LYS:HG3	2.44	0.52
1:C:299:SER:HA	1:C:339:MSE:HE2	1.93	0.51
1:A:367:ALA:HB3	1:A:385:SER:HB2	1.94	0.50
1:B:244:ALA:HB3	1:B:527:VAL:HG21	1.93	0.50
1:C:359:TYR:OH	1:C:392:ASP:OD1	2.25	0.49
1:C:259:THR:HG21	1:C:509:TYR:HE1	1.75	0.49
1:C:512[B]:MSE:CE	1:C:518:ILE:HG23	2.35	0.49
1:B:444:GLU:OE2	1:B:446:ARG:NH1	2.46	0.48
1:C:303:PRO:HG2	1:C:312:ILE:HG21	1.95	0.48
1:C:431:ILE:HG12	1:C:439:ILE:HG21	1.95	0.48
1:C:495:ASN:OD1	1:C:498:ARG:HG2	2.13	0.48
1:A:384:ILE:HG22	1:A:386:HIS:HD2	1.78	0.48
1:B:351:GLY:HA3	1:B:407:TYR:HE2	1.79	0.48
1:C:307:PRO:HG2	1:C:411:ASN:HA	1.96	0.48
1:A:62:TYR:CD1	1:B:6:ILE:HG21	2.49	0.47
1:C:35:GLN:CB	1:C:218:MSE:HE1	2.43	0.47
1:A:303:PRO:HG2	1:A:312:ILE:HG21	1.95	0.47
1:B:35:GLN:CB	1:B:218:MSE:HE1	2.45	0.47
1:B:329:TYR:O	1:B:367:ALA:HA	2.15	0.46
1:C:35:GLN:HB3	1:C:218:MSE:HE1	1.98	0.46
1:B:303:PRO:HG2	1:B:312:ILE:HG21	1.97	0.46
1:A:172:LYS:CE	1:B:22:PHE:O	2.62	0.46
1:C:399:LEU:O	1:C:402:SER:HB2	2.16	0.46
1:A:23:ALA:HB1	1:C:225:LYS:HB2	1.97	0.46
1:A:191:VAL:HG12	1:A:193[B]:ARG:HG2	1.98	0.46
1:A:97:HIS:HD2	2:A:2004:HOH:O	1.98	0.45
1:A:381:PHE:HB3	1:A:384:ILE:HD11	1.98	0.45
1:A:228:PHE:HB3	1:A:532:LEU:HD11	1.98	0.45
1:B:228:PHE:HB3	1:B:532:LEU:HD11	1.98	0.45
1:C:444:GLU:OE2	1:C:446:ARG:NH1	2.50	0.45
1:A:486:PHE:HA	1:B:24:VAL:HG11	1.98	0.45
1:A:90:VAL:HG23	1:A:190:ILE:HB	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:533:TYR:CD2	1:A:543:ILE:HD12	2.52	0.45
1:A:218:MSE:HE1	1:B:6:ILE:HG23	1.99	0.45
1:C:228:PHE:HB3	1:C:532:LEU:HD11	1.99	0.45
1:C:516:HIS:ND1	1:C:517:PRO:HD2	2.32	0.44
1:C:534:LYS:HB2	1:C:546:ILE:HD11	2.00	0.44
1:C:389:ASN:OD1	1:C:391:ILE:HB	2.18	0.44
1:A:225:LYS:HB2	1:B:23:ALA:HB1	2.00	0.43
1:C:445:VAL:HG22	1:C:505:VAL:HG22	2.00	0.43
1:B:16:ILE:H	1:B:16:ILE:HG13	1.68	0.43
1:C:439:ILE:HG22	1:C:512[A]:MSE:HE3	2.00	0.43
1:C:255:LEU:HD13	1:C:503:LEU:HD13	2.00	0.43
1:A:2:ASN:HA	1:A:5:ILE:HD12	2.01	0.43
1:B:486:PHE:O	1:C:27:GLN:HG2	2.19	0.43
1:B:35:GLN:HB3	1:B:218:MSE:HE1	2.01	0.43
1:A:351:GLY:HA3	1:A:407:TYR:HE2	1.84	0.42
1:C:102:SER:HB3	1:C:107:ILE:HD11	2.01	0.42
1:A:319:GLY:HA3	1:A:331:LYS:HE3	2.01	0.42
1:C:90:VAL:HG23	1:C:190:ILE:HB	2.02	0.42
1:B:3:ASN:HA	1:B:6:ILE:HG12	2.02	0.42
1:A:348:LEU:HA	1:A:409:GLY:O	2.19	0.42
1:A:512:MSE:HG2	1:A:518:ILE:CG2	2.50	0.42
1:C:133:SER:O	1:C:137:ASP:HB2	2.20	0.42
1:B:350:PHE:CE2	1:B:405:ASN:HB2	2.54	0.42
1:B:349:MSE:O	1:B:409:GLY:HA2	2.20	0.42
1:A:62:TYR:HB3	1:A:218:MSE:HG2	2.02	0.42
1:B:90:VAL:HG23	1:B:190:ILE:HB	2.02	0.41
1:B:381:PHE:HB3	1:B:384:ILE:HD11	2.02	0.41
1:B:60:ASP:HB3	1:B:63:ILE:HG12	2.02	0.41
1:C:2:ASN:HA	1:C:5:ILE:HD12	2.03	0.41
1:C:302:PRO:HA	1:C:303:PRO:HD3	1.99	0.41
1:B:434:LYS:HA	1:B:434:LYS:HD2	1.95	0.41
1:B:225:LYS:HB2	1:C:23:ALA:HB1	2.03	0.41
1:B:172:LYS:CE	1:C:22:PHE:O	2.67	0.41
1:B:244:ALA:HB3	1:B:527:VAL:CG2	2.51	0.41
1:C:250:HIS:HA	1:C:494:ALA:HB2	2.03	0.41
1:A:444:GLU:OE2	1:A:446:ARG:NH1	2.52	0.40
1:B:533:TYR:CD1	1:B:543:ILE:HD12	2.55	0.40
1:B:37:ILE:O	1:B:213:SER:HA	2.21	0.40
1:C:431:ILE:HD12	1:C:518:ILE:HG12	2.03	0.40
1:B:271:ILE:HG13	1:B:397:VAL:HG12	2.03	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	539/569 (95%)	506 (94%)	33 (6%)	0	100	100
1	B	538/569 (95%)	506 (94%)	31 (6%)	1 (0%)	43	75
1	C	539/569 (95%)	507 (94%)	31 (6%)	1 (0%)	43	75
All	All	1616/1707 (95%)	1519 (94%)	95 (6%)	2 (0%)	48	79

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	26	SER
1	C	48	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	499/510 (98%)	469 (94%)	30 (6%)	17	48
1	B	498/510 (98%)	467 (94%)	31 (6%)	16	47
1	C	499/510 (98%)	471 (94%)	28 (6%)	19	51
All	All	1496/1530 (98%)	1407 (94%)	89 (6%)	17	49

All (89) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	ASN

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Mol	Chain	Res	Type
1	A	9	LEU
1	A	10	ILE
1	A	16	ILE
1	A	24	VAL
1	A	26	SER
1	A	37	ILE
1	A	39	LEU
1	A	93	LYS
1	A	101	SER
1	A	188	SER
1	A	213	SER
1	A	227	SER
1	A	239	ILE
1	A	246	VAL
1	A	256	SER
1	A	278	GLN
1	A	298	VAL
1	A	330	VAL
1	A	331	LYS
1	A	352	THR
1	A	366	SER
1	A	383	HIS
1	A	391	ILE
1	A	397	VAL
1	A	398	SER
1	A	406	VAL
1	A	518	ILE
1	A	535	VAL
1	A	547	MSE
1	B	3	ASN
1	B	6	ILE
1	B	9	LEU
1	B	10	ILE
1	B	16	ILE
1	B	24	VAL
1	B	26	SER
1	B	39	LEU
1	B	93	LYS
1	B	101	SER
1	B	120	ILE
1	B	174	SER
1	B	188	SER

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Mol	Chain	Res	Type
1	B	213	SER
1	B	234	ARG
1	B	239	ILE
1	B	246	VAL
1	B	256	SER
1	B	278	GLN
1	B	298	VAL
1	B	330	VAL
1	B	345	THR
1	B	352	THR
1	B	366	SER
1	B	391	ILE
1	B	397	VAL
1	B	406	VAL
1	B	497	SER
1	B	510	SER
1	B	511	THR
1	B	547	MSE
1	C	3	ASN
1	C	9	LEU
1	C	10	ILE
1	C	16	ILE
1	C	26	SER
1	C	39	LEU
1	C	46	ASP
1	C	93	LYS
1	C	188	SER
1	C	213	SER
1	C	239	ILE
1	C	246	VAL
1	C	256	SER
1	C	278	GLN
1	C	298	VAL
1	C	330	VAL
1	C	352	THR
1	C	366	SER
1	C	377	LYS
1	C	391	ILE
1	C	399	LEU
1	C	401	THR
1	C	402	SER
1	C	406	VAL

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Mol	Chain	Res	Type
1	C	471	ASN
1	C	497	SER
1	C	507	VAL
1	C	547	MSE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (30) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	2	ASN
1	A	3	ASN
1	A	45	ASN
1	A	96	ASN
1	A	203	ASN
1	A	221	ASN
1	A	223	GLN
1	A	241	GLN
1	A	268	ASN
1	A	278	GLN
1	A	324	GLN
1	A	383	HIS
1	A	522	GLN
1	B	45	ASN
1	B	96	ASN
1	B	203	ASN
1	B	223	GLN
1	B	268	ASN
1	B	278	GLN
1	B	324	GLN
1	C	35	GLN
1	C	96	ASN
1	C	203	ASN
1	C	223	GLN
1	C	268	ASN
1	C	278	GLN
1	C	324	GLN
1	C	383	HIS
1	C	435	ASN
1	C	530	ASN

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	534/569 (93%)	-0.47	7 (1%)	75 53	28, 53, 83, 169	1 (0%)
1	B	534/569 (93%)	-0.44	5 (0%)	81 61	42, 56, 87, 154	0
1	C	534/569 (93%)	-0.30	6 (1%)	78 57	42, 65, 100, 170	0
All	All	1602/1707 (93%)	-0.40	18 (1%)	78 57	28, 58, 93, 170	1 (0%)

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	6	ILE	6.9
1	B	10	ILE	6.1
1	C	6	ILE	5.8
1	A	5	ILE	5.1
1	C	10	ILE	4.7
1	A	2	ASN	4.3
1	A	10	ILE	3.6
1	C	5	ILE	3.4
1	A	48	PRO	3.3
1	C	2	ASN	3.0
1	A	4	THR	2.9
1	B	4	THR	2.7
1	B	48	PRO	2.6
1	B	5	ILE	2.5
1	B	2	ASN	2.5
1	A	9	LEU	2.2
1	C	8	SER	2.1
1	C	48	PRO	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 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.