



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 03:28 AM UTC

PDB ID : 1M9D / pdb_00001m9d
Title : X-ray crystal structure of Cyclophilin A/HIV-1 CA N-terminal domain (1-146)
O-type chimera Complex.
Authors : Howard, B.R.; Vajdos, F.F.; Li, S.; Sundquist, W.I.; Hill, C.P.
Deposited on : 2002-07-28
Resolution : 1.90 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

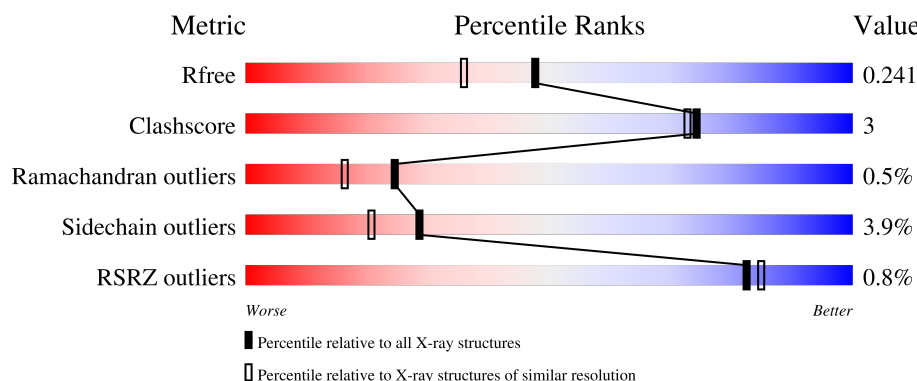
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	165	% 84% 15% .
1	B	165	% 82% 15% ..
2	C	146	% 82% 16% .
2	D	146	% 71% 21% 8%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5363 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 Cyclophilin A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	165	Total	C	N	O	S	0	0	0
			1266	802	218	237	9			
1	B	163	Total	C	N	O	S	0	0	0
			1251	792	216	235	8			

- Molecule 2 is a protein called HIV-1 Capsid.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	146	Total	C	N	O	S	0	0	0
			1131	716	198	209	8			
2	D	146	Total	C	N	O	S	0	0	0
			1131	716	198	209	8			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	83	THR	LEU	engineered mutation	UNP Q72497
C	86	PRO	VAL	engineered mutation	UNP Q72497
C	87	ALA	HIS	engineered mutation	UNP Q72497
C	88	MET	ALA	engineered mutation	UNP Q72497
C	91	LEU	ILE	engineered mutation	UNP Q72497
C	92	PRO	ALA	engineered mutation	UNP Q72497
C	96	ILE	MET	engineered mutation	UNP Q72497
C	120	HIS	ASN	SEE REMARK 999	UNP Q72497
D	83	THR	LEU	engineered mutation	UNP Q72497
D	86	PRO	VAL	engineered mutation	UNP Q72497
D	87	ALA	HIS	engineered mutation	UNP Q72497
D	88	MET	ALA	engineered mutation	UNP Q72497
D	91	LEU	ILE	engineered mutation	UNP Q72497
D	92	PRO	ALA	engineered mutation	UNP Q72497
D	96	ILE	MET	engineered mutation	UNP Q72497
D	120	HIS	ASN	SEE REMARK 999	UNP Q72497

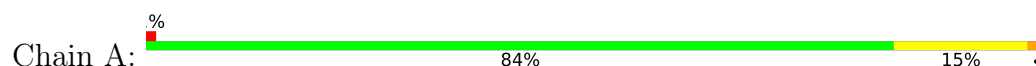
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	172	Total 172	O 172	0	0
3	B	144	Total 144	O 144	0	0
3	C	154	Total 154	O 154	0	0
3	D	114	Total 114	O 114	0	0

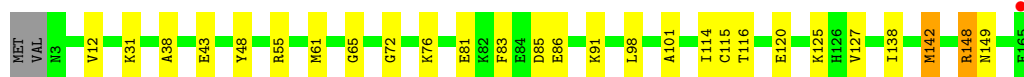
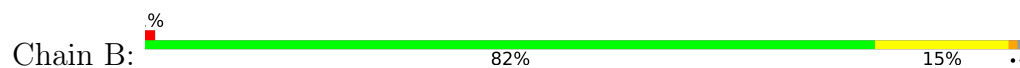
3 Residue-property plots [i](#)

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.

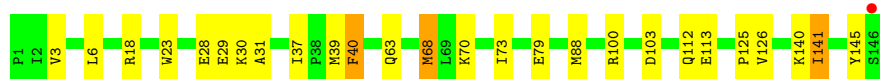
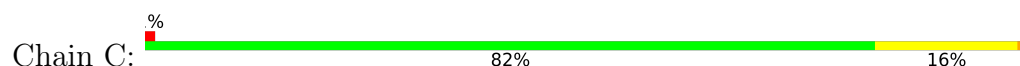
- Molecule 1: Cyclophilin A



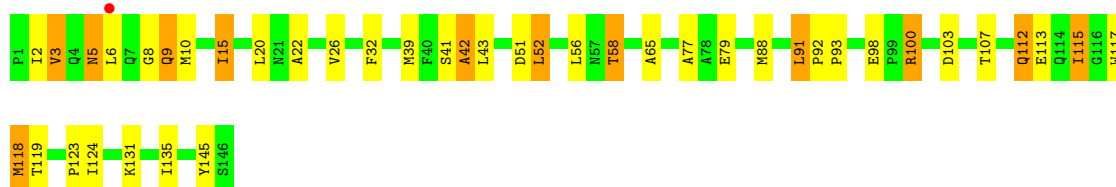
- Molecule 1: Cyclophilin A



- Molecule 2: HIV-1 Capsid



- Molecule 2: HIV-1 Capsid



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	38.90Å 109.30Å 67.70Å 90.00° 99.70° 90.00°	Depositor
Resolution (Å)	67.42 – 1.90 66.73 – 1.90	Depositor EDS
% Data completeness (in resolution range)	94.0 (67.42-1.90) 93.3 (66.73-1.90)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.00 (at 1.89Å)	Xtriage
Refinement program	REFMAC 5.0	Depositor
R, R_{free}	0.163 , 0.232 0.173 , 0.241	Depositor DCC
R_{free} test set	4031 reflections (9.55%)	wwPDB-VP
Wilson B-factor (Å ²)	24.4	Xtriage
Anisotropy	0.062	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 40.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	5363	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.97% 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	1.00	2/1294 (0.2%)	1.72	22/1733 (1.3%)
1	B	0.91	2/1279 (0.2%)	1.74	21/1713 (1.2%)
2	C	1.02	1/1160 (0.1%)	1.75	16/1580 (1.0%)
2	D	1.00	4/1160 (0.3%)	1.98	34/1580 (2.2%)
All	All	0.98	9/4893 (0.2%)	1.80	93/6606 (1.4%)

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
2	C	0	1
2	D	0	2
All	All	0	3

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	68	MET	SD-CE	-10.38	1.53	1.79
1	B	142	MET	SD-CE	-10.35	1.53	1.79
2	D	39	MET	SD-CE	-6.41	1.63	1.79
1	B	38	ALA	CA-CB	-6.14	1.43	1.53
2	D	88	MET	SD-CE	-6.10	1.64	1.79
1	A	61	MET	SD-CE	5.70	1.93	1.79
1	A	89	ILE	CA-CB	-5.33	1.47	1.54
2	D	65	ALA	CA-CB	-5.32	1.45	1.53
2	D	77	ALA	CA-CB	-5.10	1.45	1.53

All (93) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	100	ARG	NE-CZ-NH2	-9.80	110.38	119.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	91	LEU	CA-C-N	9.26	126.33	119.66
2	D	91	LEU	C-N-CA	9.26	126.33	119.66
1	A	72	GLY	N-CA-C	-8.22	103.14	114.64
2	C	126	VAL	N-CA-C	7.74	119.44	110.62
1	B	43	GLU	N-CA-C	7.28	121.26	112.38
2	D	93	PRO	CB-CA-C	-7.09	102.50	111.71
1	B	148	ARG	CA-C-N	7.01	130.97	120.31
1	B	148	ARG	C-N-CA	7.01	130.97	120.31
2	D	115	ILE	N-CA-CB	6.96	118.69	110.55
2	C	112	GLN	CB-CG-CD	6.95	124.42	112.60
1	A	164	LEU	CB-CA-C	-6.81	105.89	114.40
2	C	37	ILE	CA-C-O	6.70	123.18	118.69
2	D	41	SER	CA-C-O	-6.63	113.39	120.42
1	A	65	GLY	N-CA-C	6.63	124.51	115.30
2	D	145	TYR	OH-CZ-CE2	6.62	139.77	119.90
1	A	101	ALA	N-CA-C	-6.61	101.63	110.55
1	B	48	TYR	N-CA-C	6.58	119.45	111.82
2	D	32	PHE	N-CA-C	6.56	120.44	111.39
2	D	88	MET	CA-C-O	6.51	128.30	120.81
2	C	113	GLU	O-C-N	6.49	128.75	122.07
2	D	124	ILE	O-C-N	6.46	125.44	121.37
1	A	68	THR	N-CA-C	-6.41	105.48	113.55
2	D	100	ARG	NE-CZ-NH1	6.34	127.84	121.50
2	D	52	LEU	CB-CG-CD1	6.31	129.64	110.70
2	D	79	GLU	N-CA-C	6.30	117.81	111.07
1	B	98	LEU	CD1-CG-CD2	6.26	124.56	110.80
1	A	70	HIS	N-CA-C	6.25	120.48	112.86
1	A	70	HIS	N-CA-CB	-6.20	102.79	112.41
2	C	145	TYR	CA-C-O	-6.20	114.15	121.16
2	C	73	ILE	N-CA-C	-6.16	104.51	110.42
2	D	103	ASP	CA-C-O	-6.11	114.41	120.70
2	D	58	THR	OG1-CB-CG2	-6.11	97.09	109.30
2	D	43	LEU	N-CA-CB	-6.10	101.47	110.49
1	B	85	ASP	N-CA-C	-6.10	100.63	109.59
2	C	100	ARG	NE-CZ-NH1	5.96	127.46	121.50
2	C	3	VAL	CA-CB-CG2	5.78	120.22	110.40
1	A	164	LEU	CA-C-N	5.73	132.01	121.70
1	A	164	LEU	C-N-CA	5.73	132.01	121.70
1	B	101	ALA	N-CA-C	-5.72	101.23	110.32
2	D	117	TRP	CA-C-O	5.71	126.58	120.70
2	D	41	SER	CA-CB-OG	-5.71	99.69	111.10
2	D	113	GLU	N-CA-C	5.70	118.22	111.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	123	PRO	CA-C-O	5.69	127.83	121.34
2	C	23	TRP	N-CA-C	5.68	117.28	111.14
2	C	125	PRO	N-CA-CB	5.67	106.78	102.36
2	D	3	VAL	CB-CA-C	-5.66	101.63	110.69
1	A	163	GLN	CB-CA-C	5.66	119.06	109.50
2	D	100	ARG	CG-CD-NE	-5.63	99.61	112.00
1	B	149	ASN	CB-CA-C	-5.62	99.89	110.67
1	B	127	VAL	N-CA-C	5.58	116.79	108.54
1	B	115	CYS	CA-C-O	-5.57	114.93	121.16
2	D	135	ILE	N-CA-C	5.56	116.29	110.62
1	B	116	THR	OG1-CB-CG2	-5.56	98.19	109.30
2	C	103	ASP	N-CA-C	5.55	117.41	111.36
1	B	116	THR	N-CA-CB	5.54	119.05	110.80
2	D	123	PRO	CA-C-N	-5.50	116.58	122.85
2	D	123	PRO	C-N-CA	-5.50	116.58	122.85
1	A	148	ARG	N-CA-C	5.50	118.79	111.75
1	B	120	GLU	N-CA-C	5.45	117.97	111.71
1	B	65	GLY	N-CA-C	5.43	122.82	115.32
1	B	72	GLY	CA-C-O	-5.42	113.01	119.02
1	A	48	TYR	N-CA-C	5.36	118.04	111.82
1	A	81	GLU	N-CA-CB	5.35	119.53	110.49
2	C	40	PHE	CA-C-O	5.35	126.44	120.82
2	D	113	GLU	O-C-N	5.34	127.78	122.12
2	D	107	THR	OG1-CB-CG2	-5.32	98.66	109.30
1	A	143	GLU	CB-CA-C	-5.32	99.11	110.32
1	A	131	LYS	CD-CE-NZ	5.29	128.82	111.90
1	A	105	PRO	CA-C-N	5.26	129.83	122.36
1	A	105	PRO	C-N-CA	5.26	129.83	122.36
2	C	30	LYS	N-CA-C	5.26	119.01	112.59
2	D	32	PHE	CA-C-O	5.25	127.93	121.84
1	A	141	ALA	N-CA-C	-5.24	105.65	111.36
1	B	12	VAL	N-CA-C	-5.24	100.02	107.77
1	B	114	ILE	CA-CB-CG2	-5.24	101.59	110.50
1	A	19	ARG	CA-C-N	-5.23	115.72	122.99
1	A	19	ARG	C-N-CA	-5.23	115.72	122.99
1	B	98	LEU	CA-C-O	5.20	125.86	120.30
2	D	98	GLU	N-CA-CB	5.17	117.44	110.14
1	B	86	GLU	O-C-N	5.14	127.64	122.09
1	A	114	ILE	N-CA-CB	5.13	117.92	111.46
1	A	20	VAL	N-CA-C	-5.11	100.76	108.12
1	B	76	LYS	N-CA-CB	-5.11	102.29	110.77
2	D	79	GLU	CG-CD-OE2	-5.09	106.69	118.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	42	ALA	O-C-N	-5.08	116.28	122.22
2	C	141	ILE	CA-C-O	-5.06	115.00	120.47
2	C	39	MET	CG-SD-CE	-5.06	89.77	100.90
1	B	83	PHE	CA-C-O	5.05	127.37	121.46
2	D	112	GLN	CB-CG-CD	5.05	121.18	112.60
2	D	92	PRO	CA-C-N	5.02	125.80	119.98
2	D	92	PRO	C-N-CA	5.02	125.80	119.98
2	C	79	GLU	CG-CD-OE1	-5.01	106.89	118.40

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	C	40	PHE	Sidechain
2	D	118	MET	Mainchain
2	D	42	ALA	Mainchain

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	1266	0	1237	6	0
1	B	1251	0	1216	5	0
2	C	1131	0	1131	8	0
2	D	1131	0	1131	10	0
3	A	172	0	0	2	0
3	B	144	0	0	3	0
3	C	154	0	0	3	0
3	D	114	0	0	0	0
All	All	5363	0	4715	29	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 3.

All (29) 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)
2:C:68:MET:HE2	2:C:141:ILE:HG22	1.40	0.99
1:A:134:GLU:OE1	3:A:312:HOH:O	2.12	0.66
2:D:8:GLY:O	2:D:9:GLN:HG2	1.97	0.64
2:D:2:ILE:HD12	2:D:119:THR:HG22	1.79	0.64
2:D:20:LEU:HD11	2:D:58:THR:HG21	1.80	0.63
2:C:68:MET:HE3	2:C:140:LYS:HG2	1.82	0.61
2:D:15:ILE:HG23	2:D:51:ASP:OD1	2.02	0.60
2:D:2:ILE:HG12	2:D:10:MET:HE3	1.86	0.57
2:C:140:LYS:NZ	3:C:289:HOH:O	2.38	0.56
1:B:125:LYS:NZ	3:B:280:HOH:O	2.39	0.54
1:A:84:GLU:H	1:A:84:GLU:CD	2.16	0.53
1:A:10:ILE:HD12	1:A:10:ILE:N	2.24	0.52
2:C:70:LYS:NZ	3:C:286:HOH:O	2.43	0.52
2:C:18:ARG:NH1	3:C:299:HOH:O	2.35	0.50
1:B:81:GLU:HG3	3:B:182:HOH:O	2.12	0.50
1:A:31:LYS:NZ	1:A:84:GLU:OE2	2.35	0.50
2:C:68:MET:HE3	2:C:140:LYS:CG	2.43	0.48
1:A:149:ASN:ND2	3:A:266:HOH:O	2.30	0.47
1:B:138:ILE:HG22	1:B:142:MET:CE	2.45	0.47
2:C:68:MET:CE	2:C:141:ILE:HG22	2.28	0.46
2:C:6:LEU:H	2:C:6:LEU:HD12	1.83	0.43
2:D:2:ILE:HD11	2:D:10:MET:CB	2.48	0.43
2:D:2:ILE:CD1	2:D:119:THR:HG22	2.46	0.43
1:A:118:LYS:HD3	1:A:120:GLU:HG3	2.01	0.43
1:B:55:ARG:HG2	1:B:61:MET:HE1	2.02	0.42
2:D:22:ALA:O	2:D:26:VAL:HG23	2.20	0.42
2:D:2:ILE:CG2	2:D:118:MET:HE2	2.49	0.42
2:D:5:ASN:OD1	2:D:5:ASN:C	2.64	0.40
1:B:31:LYS:NZ	3:B:241:HOH:O	2.54	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	163/165 (99%)	155 (95%)	7 (4%)	1 (1%)	21	13
1	B	161/165 (98%)	155 (96%)	6 (4%)	0	100	100
2	C	144/146 (99%)	140 (97%)	3 (2%)	1 (1%)	18	10
2	D	144/146 (99%)	140 (97%)	3 (2%)	1 (1%)	18	10
All	All	612/622 (98%)	590 (96%)	19 (3%)	3 (0%)	24	16

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	C	31	ALA
1	A	81	GLU
2	D	9	GLN

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	133/133 (100%)	130 (98%)	3 (2%)	44	40
1	B	131/133 (98%)	129 (98%)	2 (2%)	57	56
2	C	123/124 (99%)	119 (97%)	4 (3%)	33	26
2	D	123/124 (99%)	112 (91%)	11 (9%)	9	4
All	All	510/514 (99%)	490 (96%)	20 (4%)	28	21

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	84	GLU
1	A	131	LYS
1	A	133	LYS
1	B	91	LYS
1	B	148	ARG
2	C	28	GLU
2	C	29	GLU
2	C	63	GLN

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Mol	Chain	Res	Type
2	C	88	MET
2	D	3	VAL
2	D	5	ASN
2	D	6	LEU
2	D	15	ILE
2	D	52	LEU
2	D	56	LEU
2	D	91	LEU
2	D	100	ARG
2	D	112	GLN
2	D	115	ILE
2	D	131	LYS

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	137	ASN
1	B	106	ASN
2	C	57	ASN
2	C	95	GLN
2	D	7	GLN
2	D	62	HIS
2	D	95	GLN
2	D	139	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 [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	165/165 (100%)	-0.43	2 (1%) 76 79	15, 22, 36, 61	0
1	B	163/165 (98%)	-0.27	1 (0%) 85 87	18, 28, 38, 60	0
2	C	146/146 (100%)	-0.20	1 (0%) 84 86	14, 26, 50, 63	0
2	D	146/146 (100%)	-0.01	1 (0%) 84 86	15, 29, 48, 60	0
All	All	620/622 (99%)	-0.24	5 (0%) 82 85	14, 26, 44, 63	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	81	GLU	2.6
1	A	80	GLY	2.3
2	C	146	SER	2.3
2	D	6	LEU	2.2
1	B	165	GLU	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.