



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 06:01 AM UTC

PDB ID : 3EYO / pdb_00003eyo
Title : Crystal structure of anti-human cytomegalovirus antibody 8F9
Authors : Thomson, C.A.; Bryson, S.; McLean, G.R.; Creagh, A.L.; Pai, E.F.; Schrader, J.W.
Deposited on : 2008-10-21
Resolution : 2.50 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

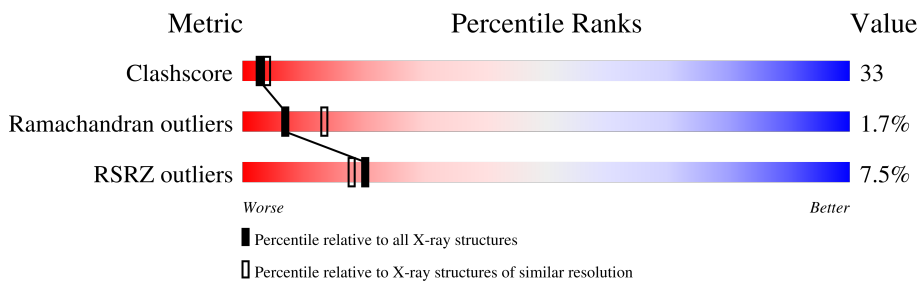
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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))
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	216	7% 58% 31% 8% ..
1	C	216	4% 54% 39% 6% .
2	B	242	8% 45% 38% 5% 12%
2	D	242	9% 46% 36% 8% 10%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6906 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 8f9 Fab.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	213	Total	C	N	O	S	0	0	0
			1643	1034	281	323	5			
1	C	214	Total	C	N	O	S	0	0	0
			1652	1039	282	326	5			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MET	-	expression tag	PDB 3EYO
C	0	MET	-	expression tag	PDB 3EYO

- Molecule 2 is a protein called AD-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	212	Total	C	N	O	S	0	0	0
			1598	1003	274	313	8			
2	D	218	Total	C	N	O	S	0	0	0
			1633	1021	283	319	10			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	235	LEU	-	expression tag	PDB 3EYO
B	236	GLU	-	expression tag	PDB 3EYO
B	237	HIS	-	expression tag	PDB 3EYO
B	238	HIS	-	expression tag	PDB 3EYO
B	239	HIS	-	expression tag	PDB 3EYO
B	240	HIS	-	expression tag	PDB 3EYO
B	241	HIS	-	expression tag	PDB 3EYO
B	242	HIS	-	expression tag	PDB 3EYO
D	235	LEU	-	expression tag	PDB 3EYO
D	236	GLU	-	expression tag	PDB 3EYO

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Chain	Residue	Modelled	Actual	Comment	Reference
D	237	HIS	-	expression tag	PDB 3EYO
D	238	HIS	-	expression tag	PDB 3EYO
D	239	HIS	-	expression tag	PDB 3EYO
D	240	HIS	-	expression tag	PDB 3EYO
D	241	HIS	-	expression tag	PDB 3EYO
D	242	HIS	-	expression tag	PDB 3EYO

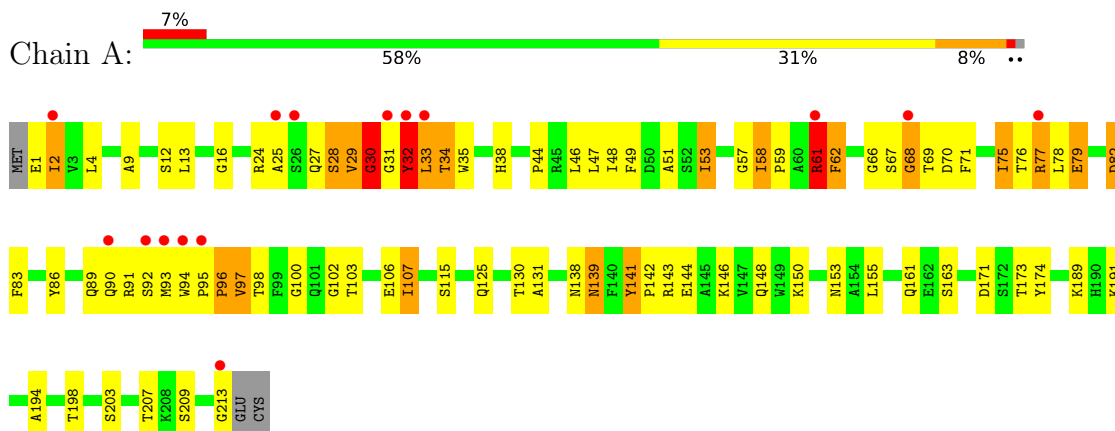
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	114	Total O 114 114	0	0
3	B	96	Total O 96 96	0	0
3	C	89	Total O 89 89	0	0
3	D	81	Total O 81 81	0	0

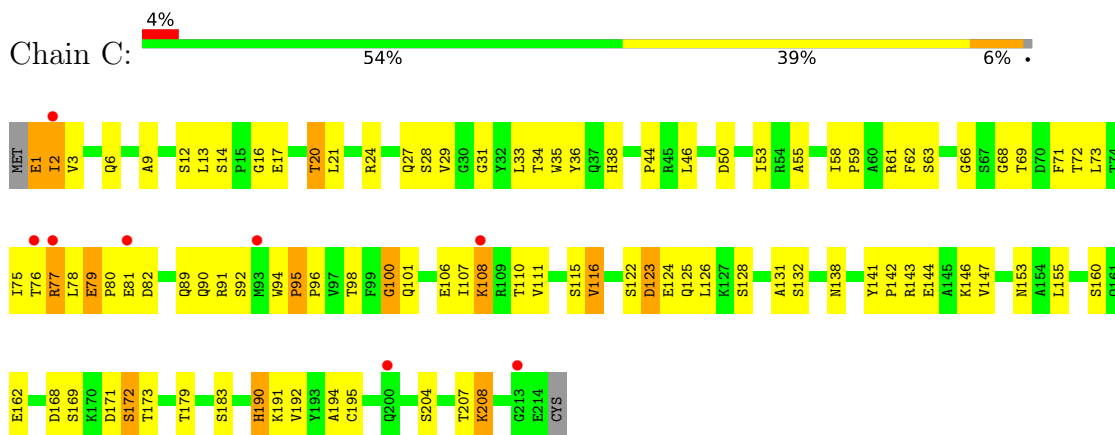
3 Residue-property plots [i](#)

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.

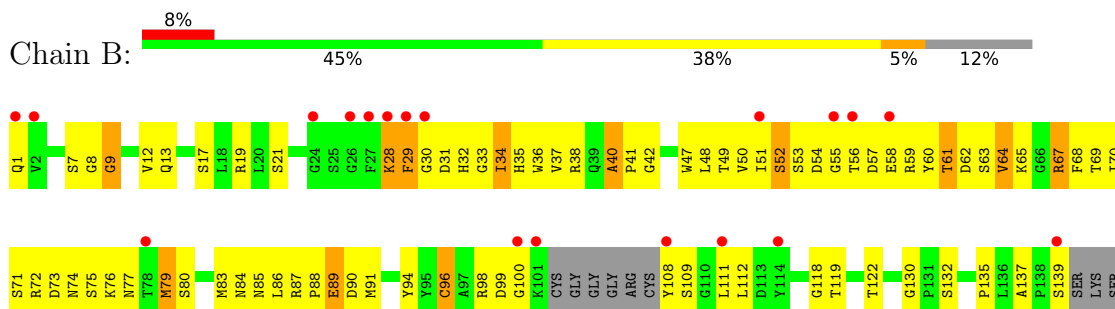
• Molecule 1: 8f9 Fab



• Molecule 1: 8f9 Fab

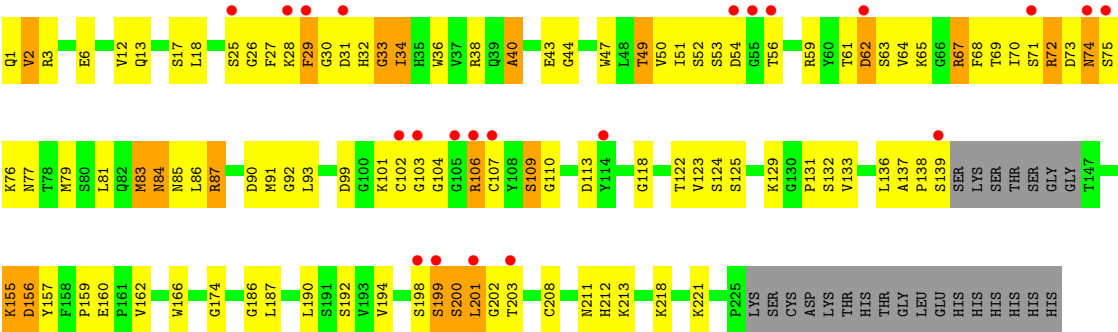


• Molecule 2: AD-2





● Molecule 2: AD-2



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	63.03Å 110.73Å 180.25Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.50 50.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	(Not available) (50.00-2.50) 95.1 (50.00-2.50)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.47 (at 2.39Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.241 , 0.268 (Not available) , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	39.7	Xtriage
Anisotropy	0.329	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 50.9	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	0.91	EDS
Total number of atoms	6906	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.47% 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 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	0.93	4/1681 (0.2%)	1.57	43/2285 (1.9%)
1	C	0.78	0/1690	1.37	34/2297 (1.5%)
2	B	0.81	0/1634	1.51	30/2217 (1.4%)
2	D	0.89	3/1670 (0.2%)	1.62	44/2265 (1.9%)
All	All	0.86	7/6675 (0.1%)	1.52	151/9064 (1.7%)

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

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	82	ASP	CA-C	-12.21	1.36	1.52
1	A	82	ASP	CA-CB	-7.51	1.42	1.54
2	D	162	VAL	N-CA	6.13	1.53	1.46
1	A	83	PHE	N-CA	-5.82	1.39	1.46
2	D	84	ASN	CA-C	-5.76	1.45	1.52
1	A	82	ASP	CB-CG	-5.27	1.38	1.52
2	D	203	THR	CA-C	-5.05	1.49	1.52

All (151) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	29	VAL	N-CA-C	-19.98	93.72	113.10
2	B	89	GLU	N-CA-C	-16.23	93.95	114.56
2	D	103	GLY	N-CA-C	14.00	129.59	112.64
2	B	29	PHE	N-CA-C	13.19	128.48	112.38
2	D	138	PRO	N-CA-C	11.84	128.68	110.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	200	SER	N-CA-C	-11.43	95.59	110.33
2	D	139	SER	N-CA-CB	-10.98	91.84	110.50
2	D	110	GLY	N-CA-C	-10.71	100.33	115.43
2	D	138	PRO	CA-C-N	10.62	140.81	121.70
2	D	138	PRO	C-N-CA	10.62	140.81	121.70
2	B	137	ALA	CA-C-N	9.68	129.79	120.31
2	B	137	ALA	C-N-CA	9.68	129.79	120.31
2	D	138	PRO	CA-C-O	9.67	132.06	121.03
2	B	28	LYS	CA-C-N	9.44	135.70	120.60
2	B	28	LYS	C-N-CA	9.44	135.70	120.60
2	D	106	ARG	N-CA-C	-9.33	99.24	110.44
2	B	63	SER	N-CA-C	9.21	123.62	112.38
2	D	199	SER	N-CA-C	-9.17	100.97	110.97
1	C	3	VAL	N-CA-C	9.15	121.34	107.99
1	C	66	GLY	N-CA-C	9.01	121.18	112.08
2	B	201	LEU	N-CA-C	8.73	121.58	111.11
2	B	79	MET	N-CA-C	-8.67	92.33	110.80
2	D	155	LYS	N-CA-C	8.64	123.50	109.59
1	A	32	TYR	N-CA-C	8.53	128.97	110.80
2	D	113	ASP	N-CA-C	8.49	122.93	112.23
2	B	155	LYS	N-CA-C	8.45	123.17	109.40
1	C	29	VAL	N-CA-C	-8.43	104.67	112.43
2	D	67	ARG	N-CA-C	-8.43	101.72	112.93
2	B	56	THR	N-CA-C	8.42	124.16	112.04
2	D	203	THR	CB-CA-C	-8.17	99.89	111.65
1	C	100	GLY	N-CA-C	-8.01	99.95	112.85
1	A	61	ARG	NE-CZ-NH2	-7.97	112.02	119.20
1	A	53	ILE	N-CA-C	7.94	119.58	107.99
2	D	61	THR	N-CA-C	-7.89	100.71	110.41
1	A	96	PRO	CA-C-N	7.82	132.18	120.77
1	A	96	PRO	C-N-CA	7.82	132.18	120.77
1	A	100	GLY	N-CA-C	-7.76	100.73	112.60
1	A	34	THR	N-CA-C	7.75	122.00	110.52
2	D	139	SER	CB-CA-C	7.70	124.73	110.10
1	C	79	GLU	N-CA-C	-7.67	100.36	110.40
1	A	32	TYR	CA-C-N	-7.59	109.87	121.86
1	A	32	TYR	C-N-CA	-7.59	109.87	121.86
1	A	33	LEU	N-CA-C	7.57	121.14	108.13
2	B	64	VAL	N-CA-C	-7.55	105.65	112.90
2	D	202	GLY	N-CA-C	-7.51	103.21	115.46
1	A	76	THR	N-CA-C	7.43	122.11	112.89
2	D	74	ASN	N-CA-C	-7.43	103.12	111.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	58	ILE	CA-C-N	7.36	128.35	120.11
1	A	58	ILE	C-N-CA	7.36	128.35	120.11
1	A	92	SER	N-CA-C	-7.29	102.98	112.68
1	A	115	SER	N-CA-C	-7.27	97.39	109.24
1	A	77	ARG	CB-CA-C	7.20	121.13	111.70
1	A	97	VAL	N-CA-C	-7.17	99.25	108.93
1	A	107	ILE	N-CA-C	7.16	118.34	108.89
1	A	143	ARG	N-CA-C	7.14	120.11	111.82
2	B	40	ALA	CA-C-N	7.08	128.93	120.23
2	B	40	ALA	C-N-CA	7.08	128.93	120.23
2	B	132	SER	N-CA-C	-6.99	97.85	109.24
1	A	75	ILE	N-CA-C	-6.89	95.76	106.72
1	A	16	GLY	N-CA-C	-6.89	105.69	114.37
1	A	78	LEU	N-CA-C	6.87	120.60	110.30
1	A	31	GLY	N-CA-C	-6.86	96.93	113.18
2	D	138	PRO	O-C-N	6.83	130.75	123.10
2	D	83	MET	CA-C-N	-6.75	111.99	122.59
2	D	83	MET	C-N-CA	-6.75	111.99	122.59
1	A	30	GLY	N-CA-C	6.75	129.17	113.18
1	C	115	SER	N-CA-C	-6.63	98.45	109.46
2	D	132	SER	N-CA-C	-6.57	98.99	109.76
1	A	97	VAL	N-CA-CB	-6.50	103.02	110.49
2	D	87	ARG	CA-C-N	6.48	126.41	119.28
2	D	87	ARG	C-N-CA	6.48	126.41	119.28
1	A	57	GLY	N-CA-C	-6.36	106.36	114.37
1	C	183	SER	N-CA-C	-6.36	100.16	110.20
2	B	34	ILE	N-CA-C	6.34	117.24	108.12
1	C	143	ARG	N-CA-C	6.30	120.23	112.54
2	B	160	GLU	N-CA-C	-6.28	101.41	110.08
2	D	139	SER	CA-CB-OG	6.27	123.64	111.10
2	D	136	LEU	N-CA-C	-6.27	95.28	107.62
2	D	62	ASP	N-CA-C	6.21	117.72	111.07
1	C	204	SER	CA-C-N	6.21	126.23	119.90
1	C	204	SER	C-N-CA	6.21	126.23	119.90
2	D	125	SER	N-CA-C	-6.20	105.36	113.17
1	A	203	SER	N-CA-C	-6.20	104.63	111.82
1	A	62	PHE	N-CA-C	6.16	118.56	108.52
1	A	2	ILE	N-CA-C	6.14	117.56	109.21
1	C	123	ASP	CA-C-N	6.12	131.08	120.68
1	C	123	ASP	C-N-CA	6.12	131.08	120.68
2	B	96	CYS	N-CA-C	-6.05	99.86	109.59
1	C	2	ILE	N-CA-CB	-6.03	101.27	111.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	83	PHE	N-CA-C	-5.97	100.55	109.15
2	B	29	PHE	N-CA-CB	-5.91	100.27	110.32
2	D	156	ASP	N-CA-C	5.89	117.80	110.91
2	B	42	GLY	N-CA-C	-5.87	106.05	115.08
1	C	138	ASN	N-CA-C	5.84	118.48	109.07
2	D	160	GLU	N-CA-C	-5.84	101.63	110.39
1	C	122	SER	CA-C-N	5.83	129.56	120.82
1	C	122	SER	C-N-CA	5.83	129.56	120.82
2	D	33	GLY	N-CA-C	-5.81	103.71	112.89
1	C	116	VAL	N-CA-C	5.80	116.94	108.53
1	C	153	ASN	N-CA-C	5.78	120.14	113.38
1	A	138	ASN	N-CA-C	5.78	118.37	109.07
1	C	95	PRO	CA-C-N	-5.77	113.16	127.00
1	C	95	PRO	C-N-CA	-5.77	113.16	127.00
1	C	9	ALA	N-CA-C	-5.76	104.44	112.45
2	D	92	GLY	N-CA-C	5.71	118.80	110.63
1	A	61	ARG	CG-CD-NE	-5.71	99.44	112.00
2	B	130	GLY	CA-C-N	5.71	125.62	119.85
2	B	130	GLY	C-N-CA	5.71	125.62	119.85
2	D	34	ILE	N-CA-C	5.71	116.09	108.27
2	D	203	THR	N-CA-C	5.70	119.68	111.02
1	C	2	ILE	CB-CA-C	5.70	120.63	111.29
1	A	51	ALA	N-CA-C	5.62	122.78	110.80
1	A	66	GLY	N-CA-C	5.62	117.80	111.85
2	B	28	LYS	CA-C-O	5.59	126.73	120.92
1	C	27	GLN	N-CA-C	-5.53	101.12	109.85
1	A	103	THR	N-CA-C	-5.50	100.06	109.24
1	A	79	GLU	N-CA-C	-5.48	102.17	110.50
1	A	9	ALA	N-CA-C	-5.47	105.23	111.14
1	C	192	VAL	N-CA-C	5.46	115.97	107.99
1	C	124	GLU	N-CA-C	5.44	119.09	112.23
2	B	186	GLY	N-CA-C	-5.44	107.81	115.32
2	D	26	GLY	N-CA-C	-5.44	107.28	114.95
1	C	20	THR	N-CA-C	-5.43	100.05	108.90
1	C	131	ALA	N-CA-C	5.41	117.22	108.34
2	B	74	ASN	N-CA-C	5.32	118.56	111.75
2	D	27	PHE	N-CA-C	5.30	116.34	108.60
2	D	87	ARG	N-CA-C	-5.30	100.29	108.55
1	C	208	LYS	N-CA-C	-5.25	100.84	109.40
1	A	82	ASP	CA-CB-CG	-5.25	107.35	112.60
1	A	77	ARG	NE-CZ-NH1	-5.22	116.28	121.50
1	C	172	SER	N-CA-C	-5.19	105.77	113.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	137	ALA	N-CA-C	5.19	117.82	110.24
2	D	49	THR	N-CA-C	5.18	115.37	108.38
1	C	194	ALA	N-CA-C	5.15	117.80	109.40
2	B	61	THR	N-CA-C	-5.14	101.89	110.17
2	B	56	THR	CA-C-N	-5.12	113.31	122.07
2	B	56	THR	C-N-CA	-5.12	113.31	122.07
1	C	190	HIS	N-CA-C	5.12	117.19	109.41
2	B	67	ARG	N-CA-C	5.10	118.65	112.23
1	A	191	LYS	N-CA-C	5.09	119.11	112.13
2	D	109	SER	N-CA-C	-5.09	104.71	112.04
2	D	186	GLY	N-CA-C	-5.08	107.68	115.00
2	D	40	ALA	N-CA-C	-5.07	103.39	109.83
2	D	162	VAL	CB-CA-C	-5.07	102.58	110.69
1	A	141	TYR	C-N-CD	5.05	131.71	120.60
2	B	119	THR	N-CA-C	-5.05	101.64	109.72
1	C	108	LYS	CD-CE-NZ	-5.05	95.75	111.90
2	D	192	SER	N-CA-C	-5.04	100.24	108.76
1	A	77	ARG	CA-CB-CG	5.03	124.16	114.10
1	C	1	GLU	CA-C-N	-5.01	112.95	121.97
1	C	1	GLU	C-N-CA	-5.01	112.95	121.97

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	61	ARG	Sidechain

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	1643	0	1613	109	0
1	C	1652	0	1619	90	0
2	B	1598	0	1555	110	0
2	D	1633	0	1586	122	0
3	A	114	0	0	12	0
3	B	96	0	0	9	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	89	0	0	8	0
3	D	81	0	0	6	0
All	All	6906	0	6373	423	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 33.

All (423) 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:34:THR:HG21	1:A:46:LEU:HD11	1.22	1.12
1:A:91:ARG:NH1	2:B:109:SER:O	1.83	1.11
2:B:62:ASP:HA	2:B:65:LYS:HE2	1.36	1.07
2:B:30:GLY:O	2:B:53:SER:HB2	1.57	1.04
1:C:28:SER:OG	1:C:68:GLY:HA2	1.56	1.03
1:A:61:ARG:HH22	1:A:82:ASP:CG	1.66	1.02
1:C:171:ASP:HB3	1:C:173:THR:OG1	1.60	1.02
2:B:36:TRP:NE1	2:B:79:MET:HE1	1.75	1.01
1:A:48:ILE:HD12	1:A:48:ILE:O	1.59	1.00
1:A:34:THR:HG21	1:A:46:LEU:CD1	1.93	0.98
2:D:6:GLU:OE1	2:D:118:GLY:N	1.96	0.96
2:D:200:SER:O	2:D:201:LEU:HB2	1.64	0.96
2:B:29:PHE:O	2:B:72:ARG:NH2	2.01	0.94
1:A:34:THR:CG2	1:A:46:LEU:HD11	1.98	0.94
1:A:48:ILE:HD12	1:A:48:ILE:C	1.92	0.93
2:D:71:SER:O	2:D:72:ARG:HB2	1.70	0.91
1:A:30:GLY:HA2	1:A:68:GLY:HA2	1.52	0.91
2:B:17:SER:HB2	3:B:337:HOH:O	1.70	0.90
2:B:69:THR:HG22	3:B:334:HOH:O	1.71	0.90
3:C:229:HOH:O	2:D:155:LYS:HE3	1.70	0.90
1:A:77:ARG:O	1:A:77:ARG:HD3	1.71	0.90
1:C:81:GLU:CD	1:C:81:GLU:H	1.80	0.89
2:D:63:SER:O	2:D:67:ARG:NH1	2.06	0.87
1:C:106:GLU:HG2	1:C:107:ILE:N	1.89	0.86
2:D:62:ASP:HA	2:D:65:LYS:HE2	1.56	0.86
2:D:32:HIS:O	2:D:72:ARG:NH2	2.08	0.86
2:B:62:ASP:HA	2:B:65:LYS:CE	2.06	0.85
1:A:2:ILE:O	1:A:98:THR:HG21	1.75	0.85
2:D:52:SER:O	2:D:72:ARG:NH1	2.10	0.84
1:A:61:ARG:NH1	1:A:82:ASP:OD1	2.10	0.84
1:C:20:THR:C	1:C:21:LEU:HD23	2.02	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:54:ASP:HB2	2:D:56:THR:HG23	1.57	0.84
1:A:28:SER:HA	1:A:68:GLY:O	1.78	0.83
2:D:91:MET:CE	2:D:123:VAL:H	1.93	0.82
1:C:123:ASP:HA	3:C:238:HOH:O	1.79	0.81
1:A:2:ILE:HD11	1:A:93:MET:CB	2.09	0.81
2:D:67:ARG:NH2	2:D:90:ASP:OD2	2.13	0.81
1:A:61:ARG:NH2	1:A:82:ASP:CG	2.37	0.81
2:B:28:LYS:O	2:B:31:ASP:HB2	1.80	0.81
2:D:36:TRP:NE1	2:D:79:MET:HE1	1.96	0.79
1:A:34:THR:HG23	1:A:49:PHE:HA	1.64	0.79
2:D:12:VAL:HG11	2:D:86:LEU:HD13	1.65	0.79
2:D:129:LYS:HE2	3:D:268:HOH:O	1.84	0.78
1:C:171:ASP:CB	1:C:173:THR:OG1	2.32	0.78
1:A:161:GLN:HB3	2:B:181:VAL:HG11	1.66	0.78
2:D:17:SER:OG	2:D:84:ASN:HB3	1.83	0.77
1:C:95:PRO:CB	1:C:96:PRO:HA	2.15	0.77
2:B:36:TRP:CE2	2:B:79:MET:HE1	2.20	0.76
1:A:90:GLN:OE1	1:A:98:THR:OG1	2.04	0.76
1:A:150:LYS:HD3	3:A:219:HOH:O	1.86	0.75
1:C:79:GLU:HB3	1:C:81:GLU:OE1	1.86	0.75
1:A:34:THR:HG23	1:A:48:ILE:O	1.86	0.75
2:D:91:MET:HE2	2:D:122:THR:HA	1.69	0.74
1:A:61:ARG:NH2	1:A:82:ASP:OD2	2.20	0.74
1:A:34:THR:CG2	1:A:48:ILE:O	2.35	0.74
1:C:81:GLU:CD	1:C:81:GLU:N	2.46	0.74
1:A:106:GLU:HG2	1:A:107:ILE:N	2.03	0.74
2:D:36:TRP:CE2	2:D:79:MET:HE1	2.23	0.74
1:A:89:GLN:HE21	1:A:97:VAL:HG12	1.52	0.74
1:C:106:GLU:OE2	1:C:141:TYR:OH	2.05	0.73
1:A:2:ILE:HD11	1:A:93:MET:HB2	1.69	0.73
1:A:35:TRP:HD1	1:A:48:ILE:HD11	1.54	0.73
1:A:77:ARG:O	1:A:77:ARG:CD	2.36	0.72
2:D:91:MET:HE2	2:D:123:VAL:H	1.53	0.72
2:B:30:GLY:O	2:B:53:SER:O	2.07	0.72
1:C:14:SER:O	1:C:17:GLU:HG3	1.88	0.72
1:A:48:ILE:C	1:A:48:ILE:CD1	2.60	0.72
2:D:62:ASP:HA	2:D:65:LYS:CE	2.21	0.71
2:D:51:ILE:HG23	2:D:70:ILE:HD13	1.73	0.71
2:D:211:ASN:OD1	2:D:218:LYS:HE3	1.91	0.70
1:A:2:ILE:HD11	1:A:93:MET:HB3	1.72	0.70
1:A:4:LEU:HD12	1:A:25:ALA:HA	1.72	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:33:GLY:O	2:D:34:ILE:HG12	1.92	0.70
1:A:79:GLU:O	1:A:82:ASP:OD2	2.11	0.69
1:A:34:THR:HG22	1:A:35:TRP:H	1.55	0.69
2:B:91:MET:CE	2:B:122:THR:HG23	2.23	0.69
2:D:54:ASP:CB	2:D:56:THR:HG23	2.22	0.69
2:D:200:SER:O	2:D:201:LEU:CB	2.41	0.69
2:D:71:SER:O	2:D:72:ARG:CB	2.40	0.69
1:A:30:GLY:HA2	1:A:68:GLY:CA	2.22	0.68
1:C:55:ALA:HB3	1:C:58:ILE:HD13	1.75	0.68
1:C:116:VAL:HG12	1:C:208:LYS:HG3	1.75	0.68
2:B:171:LEU:HD21	2:B:194:VAL:HG21	1.76	0.67
1:C:77:ARG:O	1:C:77:ARG:HG2	1.93	0.67
2:B:36:TRP:NE1	2:B:79:MET:CE	2.54	0.67
2:B:67:ARG:NH2	2:B:90:ASP:OD2	2.28	0.67
2:B:174:GLY:O	2:B:194:VAL:HA	1.94	0.67
1:A:30:GLY:CA	1:A:68:GLY:HA2	2.24	0.67
1:A:163:SER:OG	2:B:179:PRO:HD2	1.95	0.67
1:A:62:PHE:CD1	1:A:75:ILE:HG12	2.30	0.66
2:D:73:ASP:OD2	2:D:76:LYS:HB2	1.95	0.66
1:C:6:GLN:HG2	3:C:254:HOH:O	1.95	0.66
2:B:57:ASP:OD2	2:B:59:ARG:NE	2.29	0.66
1:A:27:GLN:HG2	1:A:27:GLN:O	1.94	0.66
1:A:171:ASP:HB3	1:A:173:THR:OG1	1.96	0.66
2:D:52:SER:C	2:D:72:ARG:NH1	2.54	0.65
2:D:54:ASP:HB2	2:D:56:THR:CG2	2.25	0.65
1:A:89:GLN:NE2	1:A:97:VAL:HG12	2.11	0.65
1:C:61:ARG:HH21	1:C:82:ASP:CG	2.04	0.65
2:D:62:ASP:O	2:D:65:LYS:HG2	1.97	0.65
1:C:6:GLN:OE1	1:C:100:GLY:HA3	1.97	0.65
2:D:28:LYS:CB	2:D:31:ASP:OD2	2.45	0.65
2:D:174:GLY:O	2:D:194:VAL:HA	1.98	0.64
2:B:17:SER:HB3	2:B:83:MET:O	1.97	0.64
2:B:112:LEU:N	2:B:112:LEU:HD12	2.11	0.64
1:C:24:ARG:HA	1:C:69:THR:O	1.96	0.64
2:B:73:ASP:OD2	2:B:76:LYS:HG3	1.98	0.64
1:C:110:THR:HA	3:C:295:HOH:O	1.98	0.64
1:A:34:THR:HG23	1:A:49:PHE:CA	2.27	0.64
2:B:72:ARG:HA	2:B:79:MET:HA	1.79	0.64
2:D:6:GLU:OE1	2:D:118:GLY:CA	2.45	0.64
1:A:62:PHE:CE1	1:A:75:ILE:HG12	2.33	0.64
1:A:24:ARG:HD3	3:A:327:HOH:O	1.97	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:17:GLU:O	1:C:76:THR:O	2.15	0.63
1:C:106:GLU:CG	1:C:107:ILE:N	2.61	0.63
2:D:28:LYS:O	2:D:31:ASP:OD2	2.16	0.63
1:A:2:ILE:HG13	1:A:90:GLN:OE1	1.99	0.62
2:B:224:GLU:HB2	2:B:225:PRO:HD2	1.80	0.62
2:B:62:ASP:CA	2:B:65:LYS:HE2	2.21	0.62
1:C:13:LEU:HD23	1:C:78:LEU:HD11	1.82	0.62
2:D:51:ILE:HG21	2:D:70:ILE:HG12	1.79	0.62
1:A:48:ILE:O	1:A:48:ILE:CD1	2.43	0.62
1:C:77:ARG:O	1:C:77:ARG:CG	2.47	0.62
2:D:76:LYS:HD2	3:D:294:HOH:O	2.00	0.62
1:C:28:SER:HA	1:C:68:GLY:O	2.00	0.61
2:B:7:SER:OG	2:B:21:SER:HB3	2.01	0.61
1:C:1:GLU:HG2	1:C:2:ILE:H	1.65	0.61
2:B:41:PRO:HA	3:B:314:HOH:O	2.00	0.61
1:C:20:THR:O	1:C:21:LEU:HD23	1.99	0.61
2:B:87:ARG:O	2:B:90:ASP:HB2	2.01	0.60
1:C:162:GLU:HG2	3:C:222:HOH:O	2.00	0.60
2:D:199:SER:C	2:D:200:SER:O	2.41	0.60
2:B:1:GLN:OE1	2:B:1:GLN:N	2.31	0.60
1:A:94:TRP:HB2	1:A:95:PRO:HA	1.84	0.60
2:D:28:LYS:HB2	2:D:31:ASP:OD2	2.02	0.60
2:B:32:HIS:C	2:B:53:SER:HB3	2.27	0.60
1:A:61:ARG:HH12	1:A:82:ASP:CG	2.09	0.59
1:C:55:ALA:HB3	1:C:58:ILE:CD1	2.31	0.59
1:C:16:GLY:H	1:C:78:LEU:HB2	1.68	0.59
1:C:155:LEU:HD12	1:C:155:LEU:N	2.16	0.59
1:C:168:ASP:HB3	1:C:171:ASP:HB2	1.85	0.59
2:B:13:GLN:HB2	3:B:306:HOH:O	2.02	0.59
1:A:27:GLN:O	1:A:28:SER:C	2.46	0.58
2:B:91:MET:HE3	2:B:122:THR:HG23	1.84	0.58
2:B:36:TRP:HE1	2:B:79:MET:HE1	1.64	0.58
2:B:98:ARG:O	2:B:112:LEU:HA	2.03	0.58
2:D:59:ARG:HG2	2:D:59:ARG:HH11	1.69	0.58
1:A:1:GLU:N	3:A:277:HOH:O	2.37	0.58
2:B:34:ILE:O	2:B:50:VAL:HA	2.04	0.58
2:B:112:LEU:N	2:B:112:LEU:CD1	2.67	0.58
2:D:36:TRP:HE1	2:D:79:MET:HE1	1.67	0.58
2:D:64:VAL:O	2:D:67:ARG:HB2	2.04	0.58
2:D:51:ILE:HG23	2:D:70:ILE:CD1	2.34	0.57
2:D:68:PHE:CZ	2:D:83:MET:HG2	2.39	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:52:SER:OG	2:B:54:ASP:OD1	2.21	0.57
2:B:54:ASP:OD1	2:B:55:GLY:N	2.37	0.57
1:A:34:THR:HG22	1:A:35:TRP:N	2.20	0.57
2:D:36:TRP:CZ2	2:D:79:MET:HE1	2.39	0.57
2:D:32:HIS:HD2	2:D:99:ASP:O	1.87	0.57
1:C:89:GLN:HG2	1:C:90:GLN:N	2.20	0.56
1:C:33:LEU:HD22	1:C:71:PHE:CG	2.40	0.56
1:A:141:TYR:CG	1:A:142:PRO:HA	2.41	0.56
1:A:47:LEU:HA	1:A:58:ILE:HG12	1.87	0.56
1:A:34:THR:HG22	1:A:48:ILE:O	2.05	0.56
2:D:40:ALA:HB3	2:D:43:GLU:HG3	1.88	0.56
2:D:51:ILE:CG2	2:D:70:ILE:CG2	2.84	0.56
1:C:21:LEU:HD23	1:C:21:LEU:N	2.21	0.55
1:A:94:TRP:HB2	1:A:97:VAL:HG22	1.87	0.55
1:C:90:GLN:HG2	1:C:92:SER:H	1.69	0.55
2:B:91:MET:HE3	2:B:122:THR:HA	1.88	0.55
2:D:91:MET:HE2	2:D:123:VAL:N	2.22	0.55
2:B:67:ARG:HH22	2:B:90:ASP:CG	2.15	0.55
1:C:61:ARG:NH2	1:C:82:ASP:OD1	2.39	0.55
1:A:34:THR:CG2	1:A:46:LEU:CD1	2.71	0.55
2:B:60:TYR:OH	2:B:70:ILE:HG22	2.07	0.55
2:D:51:ILE:CG2	2:D:70:ILE:HG12	2.36	0.55
2:D:74:ASN:C	2:D:76:LYS:N	2.62	0.54
1:C:79:GLU:CB	1:C:81:GLU:OE1	2.55	0.54
2:B:64:VAL:HB	2:B:68:PHE:CG	2.43	0.54
1:A:91:ARG:HG3	1:A:91:ARG:HH11	1.73	0.54
2:D:74:ASN:O	2:D:77:ASN:N	2.40	0.54
2:D:12:VAL:HG11	2:D:86:LEU:CD1	2.34	0.54
2:D:33:GLY:C	2:D:34:ILE:CG1	2.81	0.54
1:C:20:THR:HG22	1:C:72:THR:CG2	2.38	0.53
1:C:125:GLN:HE22	1:C:132:SER:H	1.54	0.53
2:B:48:LEU:HA	2:B:61:THR:CG2	2.38	0.53
2:B:67:ARG:NH2	2:B:90:ASP:CG	2.67	0.53
1:C:76:THR:O	1:C:78:LEU:N	2.41	0.53
2:D:28:LYS:HB3	2:D:31:ASP:OD2	2.09	0.53
2:D:73:ASP:OD2	2:D:76:LYS:N	2.41	0.53
1:A:90:GLN:HG2	1:A:91:ARG:N	2.24	0.53
2:B:48:LEU:HA	2:B:61:THR:HG22	1.91	0.53
1:A:29:VAL:O	1:A:32:TYR:HB2	2.09	0.53
2:B:36:TRP:HE1	2:B:79:MET:CE	2.21	0.53
1:C:171:ASP:O	1:C:172:SER:HB2	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:84:ASN:O	2:B:85:ASN:C	2.52	0.52
1:C:195:CYS:O	1:C:207:THR:HA	2.10	0.52
2:D:52:SER:OG	2:D:54:ASP:OD1	2.27	0.52
2:D:74:ASN:O	2:D:76:LYS:N	2.43	0.52
1:C:146:LYS:HZ1	1:C:147:VAL:H	1.58	0.52
1:A:61:ARG:HH21	1:A:75:ILE:CG2	2.22	0.52
2:D:2:VAL:C	2:D:3:ARG:CG	2.83	0.52
1:A:29:VAL:HG23	1:A:32:TYR:HB2	1.92	0.51
2:D:190:LEU:C	2:D:190:LEU:HD12	2.35	0.51
1:A:91:ARG:CG	1:A:91:ARG:O	2.58	0.51
1:C:125:GLN:O	1:C:128:SER:HB2	2.11	0.51
2:B:12:VAL:HG13	2:B:13:GLN:N	2.24	0.51
2:D:84:ASN:O	2:D:85:ASN:HB2	2.09	0.51
2:B:7:SER:HG	2:B:21:SER:HB3	1.75	0.51
1:A:61:ARG:CZ	1:A:82:ASP:CG	2.83	0.51
2:B:139:SER:CB	3:B:316:HOH:O	2.59	0.51
2:D:62:ASP:O	2:D:65:LYS:CG	2.58	0.51
1:A:94:TRP:HZ2	2:B:108:TYR:HB3	1.76	0.51
1:A:130:THR:HG22	1:A:131:ALA:N	2.26	0.51
2:B:73:ASP:CG	2:B:76:LYS:HG3	2.36	0.51
1:A:12:SER:O	1:A:13:LEU:HD13	2.11	0.50
2:D:28:LYS:O	2:D:30:GLY:N	2.45	0.50
1:A:69:THR:C	1:A:70:ASP:OD2	2.54	0.50
2:B:64:VAL:HG13	3:B:318:HOH:O	2.11	0.50
1:C:146:LYS:NZ	1:C:147:VAL:H	2.09	0.50
2:B:62:ASP:O	2:B:65:LYS:HB2	2.12	0.50
2:D:102:CYS:SG	2:D:107:CYS:C	2.94	0.50
2:B:139:SER:HB3	3:B:316:HOH:O	2.12	0.50
1:C:94:TRP:HA	1:C:95:PRO:C	2.37	0.50
2:D:1:GLN:CB	3:D:274:HOH:O	2.59	0.50
1:A:61:ARG:NH1	1:A:82:ASP:CG	2.69	0.50
1:A:153:ASN:HB3	3:A:298:HOH:O	2.12	0.50
1:A:91:ARG:NH1	1:A:91:ARG:HG3	2.27	0.49
2:B:79:MET:C	2:B:79:MET:SD	2.95	0.49
1:C:95:PRO:HB2	1:C:96:PRO:HA	1.94	0.49
1:C:144:GLU:CD	1:C:144:GLU:H	2.20	0.49
2:B:34:ILE:HG13	2:B:72:ARG:NH1	2.26	0.49
2:B:99:ASP:OD2	2:B:109:SER:OG	2.13	0.49
1:C:63:SER:HB3	3:C:224:HOH:O	2.11	0.49
2:D:2:VAL:C	2:D:3:ARG:HG3	2.36	0.49
2:D:76:LYS:O	2:D:77:ASN:HB2	2.11	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:131:PRO:HB3	2:D:157:TYR:HB3	1.94	0.49
1:A:146:LYS:HB3	1:A:198:THR:HB	1.94	0.49
2:B:51:ILE:HD12	2:B:71:SER:HA	1.94	0.49
2:D:28:LYS:C	2:D:30:GLY:H	2.20	0.49
2:D:199:SER:OG	2:D:200:SER:N	2.43	0.49
1:A:49:PHE:HB2	2:B:111:LEU:HD11	1.95	0.49
1:C:12:SER:OG	1:C:106:GLU:OE2	2.25	0.49
1:C:12:SER:HB3	1:C:108:LYS:HD2	1.95	0.49
2:B:12:VAL:HG13	2:B:13:GLN:O	2.13	0.49
2:B:37:VAL:HG22	2:B:47:TRP:HA	1.95	0.48
2:D:12:VAL:O	2:D:123:VAL:HA	2.13	0.48
2:D:211:ASN:OD1	2:D:218:LYS:HD2	2.13	0.48
2:B:30:GLY:O	2:B:53:SER:CB	2.45	0.48
1:C:160:SER:HA	1:C:179:THR:O	2.13	0.48
2:D:68:PHE:HB3	2:D:81:LEU:HD11	1.95	0.48
1:C:94:TRP:HB2	1:C:95:PRO:HA	1.96	0.48
2:B:59:ARG:CD	2:B:59:ARG:N	2.76	0.48
2:B:76:LYS:O	2:B:77:ASN:HB2	2.14	0.48
2:B:17:SER:HA	2:B:86:LEU:HD12	1.95	0.48
2:D:87:ARG:O	2:D:90:ASP:HB2	2.14	0.48
1:C:126:LEU:C	1:C:128:SER:N	2.68	0.48
2:D:18:LEU:HD23	2:D:18:LEU:HA	1.73	0.48
2:D:91:MET:HB2	2:D:91:MET:HE3	1.60	0.48
2:B:36:TRP:HB2	2:B:49:THR:HG23	1.96	0.48
1:C:155:LEU:N	1:C:155:LEU:CD1	2.77	0.48
2:D:13:GLN:HA	2:D:124:SER:O	2.14	0.48
2:D:198:SER:C	2:D:200:SER:O	2.57	0.48
2:B:28:LYS:H	2:B:28:LYS:HZ2	1.62	0.47
2:B:53:SER:HA	2:B:72:ARG:NH1	2.29	0.47
2:B:94:TYR:O	2:B:118:GLY:HA2	2.14	0.47
2:D:85:ASN:O	2:D:86:LEU:C	2.51	0.47
2:D:1:GLN:HB3	3:D:274:HOH:O	2.14	0.47
2:B:19:ARG:HH21	2:B:80:SER:HB2	1.80	0.47
2:B:48:LEU:O	2:B:61:THR:HG22	2.15	0.47
2:D:74:ASN:C	2:D:76:LYS:H	2.22	0.47
1:A:30:GLY:N	1:A:68:GLY:HA2	2.29	0.47
1:A:61:ARG:O	1:A:75:ILE:HA	2.14	0.47
1:C:146:LYS:HA	1:C:146:LYS:HZ2	1.80	0.47
2:D:91:MET:CE	2:D:123:VAL:N	2.72	0.47
2:D:159:PRO:O	2:D:212:HIS:NE2	2.40	0.47
1:A:58:ILE:HA	1:A:59:PRO:HD3	1.71	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:80:PRO:N	1:C:81:GLU:OE1	2.48	0.47
1:C:190:HIS:HE1	3:C:285:HOH:O	1.98	0.47
1:A:2:ILE:HG13	1:A:90:GLN:CD	2.41	0.46
2:B:135:PRO:HD3	2:B:221:LYS:HE2	1.96	0.46
2:B:197:PRO:O	2:B:200:SER:OG	2.30	0.46
1:A:30:GLY:H	1:A:68:GLY:HA2	1.80	0.46
2:B:158:PHE:HA	2:B:159:PRO:HA	1.67	0.46
2:D:70:ILE:HD11	2:D:79:MET:SD	2.55	0.46
1:A:94:TRP:CE3	1:A:97:VAL:HG22	2.50	0.46
2:B:51:ILE:O	2:B:52:SER:O	2.33	0.46
2:B:98:ARG:HG2	2:B:99:ASP:N	2.31	0.46
2:D:59:ARG:HG2	2:D:59:ARG:NH1	2.31	0.46
2:B:64:VAL:HB	2:B:68:PHE:CD1	2.51	0.46
1:C:6:GLN:N	1:C:101:GLN:OE1	2.48	0.46
2:D:211:ASN:HD21	2:D:213:LYS:HE3	1.80	0.46
1:A:141:TYR:CD1	1:A:142:PRO:HA	2.50	0.46
1:A:189:LYS:HE2	1:A:189:LYS:HB3	1.37	0.46
2:B:41:PRO:HB3	3:B:314:HOH:O	2.16	0.46
2:D:38:ARG:HA	2:D:93:LEU:O	2.16	0.46
1:A:95:PRO:HA	1:A:96:PRO:HA	1.51	0.45
2:B:51:ILE:C	2:B:52:SER:O	2.57	0.45
2:D:51:ILE:CG2	2:D:70:ILE:HG21	2.45	0.45
2:D:201:LEU:HD23	2:D:201:LEU:HA	1.67	0.45
1:A:34:THR:HG23	1:A:48:ILE:C	2.41	0.45
1:C:31:GLY:O	1:C:50:ASP:HA	2.16	0.45
2:D:28:LYS:C	2:D:30:GLY:N	2.74	0.45
2:D:62:ASP:OD1	2:D:65:LYS:HE3	2.17	0.45
2:D:156:ASP:HB3	2:D:187:LEU:HD13	1.99	0.45
1:A:130:THR:CG2	1:A:131:ALA:N	2.79	0.45
1:C:91:ARG:HD2	2:D:109:SER:O	2.17	0.45
2:D:54:ASP:O	2:D:56:THR:CG2	2.65	0.45
1:C:168:ASP:OD1	1:C:169:SER:N	2.50	0.45
1:A:213:GLY:CA	3:A:319:HOH:O	2.65	0.44
2:B:35:HIS:NE2	2:B:99:ASP:OD1	2.49	0.44
2:B:83:MET:HB2	2:B:86:LEU:HD11	1.98	0.44
1:C:101:GLN:CD	1:C:101:GLN:H	2.22	0.44
1:A:94:TRP:CZ2	2:B:108:TYR:HB3	2.52	0.44
1:A:148:GLN:OE1	1:A:155:LEU:HG	2.17	0.44
2:B:33:GLY:N	2:B:53:SER:HB3	2.32	0.44
1:A:28:SER:HA	1:A:68:GLY:C	2.41	0.44
1:C:36:TYR:HE2	1:C:89:GLN:OE1	2.01	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:2:ILE:H	1:A:2:ILE:HG12	1.34	0.44
2:B:70:ILE:HG12	2:B:71:SER:N	2.32	0.44
1:C:141:TYR:CD1	1:C:142:PRO:HA	2.52	0.44
2:D:43:GLU:HB2	2:D:44:GLY:H	1.46	0.44
2:D:52:SER:CB	2:D:54:ASP:OD1	2.65	0.44
1:C:50:ASP:HB2	1:C:53:ILE:HD12	1.98	0.44
2:D:31:ASP:OD2	2:D:31:ASP:N	2.50	0.44
2:D:54:ASP:C	2:D:56:THR:HG23	2.43	0.44
2:D:84:ASN:HB3	3:D:272:HOH:O	2.18	0.44
2:B:17:SER:HB3	2:B:84:ASN:HA	2.00	0.44
2:D:33:GLY:C	2:D:34:ILE:HG12	2.42	0.44
2:B:17:SER:HB3	2:B:84:ASN:OD1	2.18	0.44
1:A:207:THR:HG21	3:A:303:HOH:O	2.18	0.43
1:C:1:GLU:HG2	1:C:2:ILE:N	2.31	0.43
1:A:139:ASN:HB3	1:A:173:THR:HG21	2.00	0.43
1:A:144:GLU:CD	1:A:144:GLU:H	2.26	0.43
1:C:126:LEU:C	1:C:128:SER:H	2.26	0.43
2:B:8:GLY:O	2:B:9:GLY:O	2.35	0.43
2:B:156:ASP:HB3	2:B:187:LEU:HD13	2.00	0.43
2:D:49:THR:OG1	2:D:50:VAL:N	2.51	0.43
1:A:61:ARG:CZ	1:A:82:ASP:OD2	2.66	0.43
1:A:106:GLU:HG3	1:A:174:TYR:OH	2.19	0.43
2:B:156:ASP:HA	2:B:187:LEU:HB3	2.01	0.43
1:C:168:ASP:OD1	1:C:168:ASP:C	2.62	0.43
2:D:47:TRP:HZ2	2:D:50:VAL:HG12	1.83	0.43
2:B:165:SER:HB2	2:B:209:ASN:HB2	1.99	0.43
1:C:59:PRO:HG2	1:C:62:PHE:CE2	2.54	0.43
2:D:53:SER:HA	2:D:72:ARG:NH1	2.34	0.43
1:A:53:ILE:O	1:A:53:ILE:HG13	2.09	0.43
2:B:89:GLU:HG3	2:B:90:ASP:N	2.33	0.43
1:A:89:GLN:HG2	1:A:98:THR:O	2.19	0.43
1:A:153:ASN:CB	3:A:298:HOH:O	2.66	0.43
1:C:61:ARG:O	1:C:75:ILE:HA	2.19	0.43
2:D:74:ASN:C	2:D:77:ASN:H	2.27	0.43
1:A:95:PRO:HD3	3:A:279:HOH:O	2.18	0.43
2:B:73:ASP:OD1	2:B:75:SER:OG	2.34	0.43
1:A:86:TYR:O	1:A:102:GLY:HA2	2.19	0.42
2:B:29:PHE:O	2:B:29:PHE:CD2	2.72	0.42
1:C:125:GLN:NE2	1:C:132:SER:H	2.17	0.42
1:A:30:GLY:HA2	1:A:68:GLY:N	2.34	0.42
1:C:144:GLU:CD	1:C:144:GLU:N	2.76	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:104:GLY:C	2:D:106:ARG:N	2.77	0.42
1:A:1:GLU:OE1	1:A:96:PRO:HG2	2.18	0.42
2:D:54:ASP:OD1	2:D:54:ASP:N	2.47	0.42
2:B:58:GLU:C	2:B:59:ARG:HD2	2.44	0.42
2:D:52:SER:HB2	2:D:54:ASP:OD1	2.19	0.42
2:D:91:MET:HE2	2:D:122:THR:CA	2.45	0.42
2:D:211:ASN:ND2	2:D:213:LYS:HE3	2.34	0.42
1:A:67:SER:HB2	3:A:324:HOH:O	2.19	0.42
2:B:38:ARG:HB3	2:B:48:LEU:HD11	2.02	0.42
2:D:29:PHE:O	2:D:29:PHE:CD2	2.72	0.42
2:D:51:ILE:HG22	2:D:70:ILE:HG21	2.01	0.42
1:C:38:HIS:CD2	1:C:44:PRO:HG3	2.55	0.42
1:A:213:GLY:HA2	3:A:319:HOH:O	2.20	0.42
2:D:51:ILE:HG22	2:D:70:ILE:CG2	2.50	0.42
1:A:71:PHE:CD2	1:A:71:PHE:N	2.87	0.42
2:D:84:ASN:O	2:D:85:ASN:CB	2.68	0.42
1:A:47:LEU:C	1:A:48:ILE:CG2	2.93	0.42
2:B:70:ILE:CG1	2:B:71:SER:N	2.83	0.42
2:B:178:PHE:O	2:B:190:LEU:HD22	2.20	0.42
1:C:90:GLN:OE1	1:C:98:THR:OG1	2.37	0.42
1:C:146:LYS:HA	1:C:146:LYS:HD2	1.80	0.42
1:C:191:LYS:HE3	1:C:191:LYS:HB3	1.76	0.42
2:D:54:ASP:C	2:D:56:THR:N	2.78	0.42
2:B:86:LEU:HD23	2:B:90:ASP:CB	2.49	0.41
1:C:77:ARG:HG3	1:C:79:GLU:OE2	2.20	0.41
1:C:110:THR:O	1:C:111:VAL:C	2.62	0.41
2:D:68:PHE:CE1	2:D:83:MET:HG2	2.54	0.41
2:D:101:LYS:HZ3	2:D:101:LYS:HG3	1.74	0.41
2:D:133:VAL:O	2:D:221:LYS:HE3	2.20	0.41
1:A:125:GLN:NE2	3:A:225:HOH:O	2.53	0.41
2:B:70:ILE:C	2:B:71:SER:OG	2.63	0.41
1:C:28:SER:OG	1:C:68:GLY:CA	2.47	0.41
1:C:123:ASP:CA	3:C:238:HOH:O	2.55	0.41
2:D:104:GLY:C	2:D:106:ARG:H	2.28	0.41
2:B:51:ILE:O	2:B:52:SER:C	2.63	0.41
2:B:77:ASN:HD22	2:B:77:ASN:HA	1.65	0.41
1:C:35:TRP:CE2	1:C:73:LEU:HB2	2.56	0.41
2:D:68:PHE:C	2:D:69:THR:CG2	2.94	0.41
2:B:64:VAL:CG1	3:B:318:HOH:O	2.68	0.41
2:D:211:ASN:OD1	2:D:218:LYS:CE	2.66	0.41
2:B:100:GLY:O	2:B:109:SER:OG	2.38	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:65:LYS:HG2	2:D:65:LYS:H	1.51	0.41
1:C:34:THR:HG23	1:C:46:LEU:HD11	2.03	0.41
1:A:38:HIS:CD2	1:A:44:PRO:HG3	2.56	0.41
2:B:8:GLY:O	2:B:9:GLY:C	2.63	0.41
2:B:40:ALA:HA	2:B:41:PRO:HD3	1.80	0.41
2:D:1:GLN:HB2	3:D:274:HOH:O	2.19	0.41
2:D:166:TRP:CH2	2:D:208:CYS:HB3	2.55	0.41
1:A:161:GLN:OE1	2:B:181:VAL:HG12	2.21	0.40
2:B:35:HIS:O	2:B:96:CYS:HA	2.21	0.40
1:C:36:TYR:CE2	1:C:89:GLN:OE1	2.74	0.40
1:A:194:ALA:HB2	1:A:209:SER:HB3	2.04	0.40
2:B:87:ARG:HA	2:B:88:PRO:HD3	1.79	0.40
2:D:51:ILE:CG2	2:D:70:ILE:CG1	2.98	0.40
1:A:207:THR:CG2	3:A:303:HOH:O	2.70	0.40
2:B:51:ILE:O	2:B:51:ILE:HG23	2.21	0.40
1:C:126:LEU:HA	1:C:126:LEU:HD23	1.84	0.40
1:A:30:GLY:HA2	1:A:68:GLY:H	1.86	0.40
1:A:33:LEU:HD22	1:A:71:PHE:CD1	2.57	0.40
1:C:13:LEU:HD23	1:C:78:LEU:CD1	2.49	0.40
1:C:76:THR:O	1:C:77:ARG:C	2.64	0.40
1:C:108:LYS:HZ3	1:C:108:LYS:HG3	1.67	0.40
1:C:141:TYR:CG	1:C:142:PRO:HA	2.56	0.40
1:A:67:SER:O	1:A:69:THR:N	2.53	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	211/216 (98%)	197 (93%)	9 (4%)	5 (2%)	4 8
1	C	212/216 (98%)	199 (94%)	12 (6%)	1 (0%)	24 43

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	B	206/242 (85%)	187 (91%)	17 (8%)	2 (1%)	12	24
2	D	214/242 (88%)	191 (89%)	17 (8%)	6 (3%)	4	6
All	All	843/916 (92%)	774 (92%)	55 (6%)	14 (2%)	7	13

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	32	TYR
2	B	52	SER
2	D	2	VAL
2	D	72	ARG
1	A	30	GLY
1	A	68	GLY
1	A	139	ASN
2	B	9	GLY
2	D	29	PHE
2	D	75	SER
2	D	201	LEU
1	A	28	SER
1	C	77	ARG
2	D	25	SER

5.3.2 Protein sidechains [i](#)

There are no protein residues with a non-rotameric sidechain to report in this entry.

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

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	213/216 (98%)	0.39	15 (7%)	22 20	14, 36, 60, 75	0
1	C	214/216 (99%)	0.58	8 (3%)	45 40	23, 42, 56, 69	0
2	B	212/242 (87%)	0.57	19 (8%)	15 13	14, 41, 72, 82	0
2	D	218/242 (90%)	0.52	22 (10%)	12 10	17, 39, 65, 73	0
All	All	857/916 (93%)	0.52	64 (7%)	20 18	14, 39, 66, 82	0

All (64) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	93	MET	5.0
1	A	31	GLY	4.3
2	B	108	TYR	4.3
2	B	114	TYR	3.9
1	A	213	GLY	3.7
2	D	107	CYS	3.7
2	B	30	GLY	3.6
2	D	139	SER	3.5
1	C	77	ARG	3.5
1	C	108	LYS	3.4
2	B	101	LYS	3.4
2	D	71	SER	3.4
1	A	77	ARG	3.2
2	B	29	PHE	3.1
1	A	33	LEU	3.1
2	D	105	GLY	3.1
1	A	32	TYR	3.0
1	C	76	THR	3.0
2	D	203	THR	2.9
2	B	139	SER	2.9
2	D	31	ASP	2.8

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Mol	Chain	Res	Type	RSRZ
1	C	213	GLY	2.6
1	A	92	SER	2.6
1	C	200	GLN	2.6
2	B	2	VAL	2.6
2	D	102	CYS	2.6
2	B	28	LYS	2.6
1	A	2	ILE	2.5
2	D	25	SER	2.5
2	B	56	THR	2.5
2	D	62	ASP	2.4
1	C	81	GLU	2.4
2	D	55	GLY	2.3
1	A	90	GLN	2.3
2	B	1	GLN	2.3
1	C	2	ILE	2.3
2	B	78	THR	2.3
2	B	27	PHE	2.3
2	D	29	PHE	2.3
2	B	51	ILE	2.2
2	D	114	TYR	2.2
2	B	26	GLY	2.2
2	B	111	LEU	2.2
2	D	198	SER	2.2
2	B	55	GLY	2.2
2	D	199	SER	2.2
2	B	58	GLU	2.2
1	A	94	TRP	2.2
1	A	61	ARG	2.2
1	A	25	ALA	2.1
1	C	93	MET	2.1
2	B	24	GLY	2.1
2	D	56	THR	2.1
2	D	201	LEU	2.1
2	D	54	ASP	2.1
2	D	106	ARG	2.1
2	D	75	SER	2.1
2	B	100	GLY	2.1
1	A	68	GLY	2.1
1	A	95	PRO	2.1
2	D	28	LYS	2.0
1	A	26	SER	2.0
2	D	74	ASN	2.0

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Mol	Chain	Res	Type	RSRZ
2	D	103	GLY	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.