



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 03:10 AM UTC

PDB ID : 1MIM / pdb_00001mim
Title : IGG FAB FRAGMENT (CD25-BINDING)
Authors : Mikol, V.
Deposited on : 1995-12-04
Resolution : 2.60 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

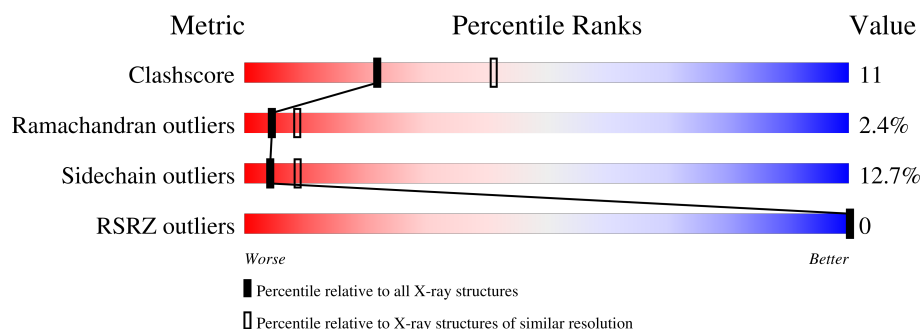
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	L	210	
2	H	215	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3303 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 CHIMERIC SDZ CHI621.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	210	Total	C	N	O	S	0	0	0
			1604	996	271	329	8			

There are 41 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
L	4	SER	LEU	conflict	UNP P01834
L	10	ILE	THR	conflict	UNP P01834
L	11	MET	LEU	conflict	UNP P01834
L	13	ALA	LEU	conflict	UNP P01834
L	18	LYS	ARG	conflict	UNP P01834
L	19	VAL	ALA	conflict	UNP P01834
L	21	MET	LEU	conflict	UNP P01834
L	22	THR	SER	conflict	UNP P01834
L	24	SER	ARG	conflict	UNP P01834
L	?	-	GLN	deletion	UNP P01834
L	28	SER	VAL	conflict	UNP P01834
L	29	ARG	GLY	conflict	UNP P01834
L	32	MET	LEU	conflict	UNP P01834
L	33	GLN	ALA	conflict	UNP P01834
L	41	THR	GLN	conflict	UNP P01834
L	42	SER	ALA	conflict	UNP P01834
L	44	LYS	ARG	conflict	UNP P01834
L	45	ARG	PRO	conflict	UNP P01834
L	46	TRP	LEU	conflict	UNP P01834
L	50	THR	ALA	conflict	UNP P01834
L	52	LYS	ASN	conflict	UNP P01834
L	53	LEU	ARG	conflict	UNP P01834
L	55	SER	THR	conflict	UNP P01834
L	57	VAL	ILE	conflict	UNP P01834
L	69	SER	ASP	conflict	UNP P01834
L	70	TYR	PHE	conflict	UNP P01834
L	71	SER	THR	conflict	UNP P01834

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Chain	Residue	Modelled	Actual	Comment	Reference
L	77	MET	LEU	conflict	UNP P01834
L	79	ALA	PRO	conflict	UNP P01834
L	82	ALA	PHE	conflict	UNP P01834
L	84	THR	VAL	conflict	UNP P01834
L	88	HIS	GLN	conflict	UNP P01834
L	89	GLN	HIS	conflict	UNP P01834
L	?	-	ASP	deletion	UNP P01834
L	?	-	ASN	deletion	UNP P01834
L	?	-	TRP	deletion	UNP P01834
L	?	-	PRO	deletion	UNP P01834
L	91	SER	PRO	conflict	UNP P01834
L	92	SER	GLY	conflict	UNP P01834
L	93	TYR	ALA	conflict	UNP P01834
L	101	LEU	VAL	conflict	UNP P01834

- Molecule 2 is a protein called CHIMERIC SDZ CHI621.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	215	Total	C	N	O	S	0	0	0
			1630	1032	270	321	7			

There are 33 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
H	3	GLN	-	insertion	GB 9857755
H	17	LYS	ARG	conflict	GB 9857755
H	29	ARG	SER	conflict	GB 9857755
H	32	MET	LEU	conflict	GB 9857755
H	48	ALA	GLY	conflict	GB 9857755
H	54	SER	ARG	conflict	GB 9857755
H	57	SER	ARG	conflict	GB 9857755
H	59	ASN	THR	conflict	GB 9857755
H	61	LYS	ARG	conflict	GB 9857755
H	63	GLU	LYS	conflict	GB 9857755
H	64	GLY	ASP	conflict	GB 9857755
H	75	SER	ASN	conflict	GB 9857755
H	86	HIS	ASN	conflict	GB 9857755
H	96	ARG	ILE	conflict	GB 9857755
H	97	ASP	ILE	conflict	GB 9857755
H	99	GLY	PHE	conflict	GB 9857755
H	100	TYR	ASP	conflict	GB 9857755
H	102	PHE	ALA	conflict	GB 9857755

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Chain	Residue	Modelled	Actual	Comment	Reference
H	?	-	ILE	deletion	GB 9857755
H	?	-	MET	deletion	GB 9857755
H	?	-	ASP	deletion	GB 9857755
H	?	-	TYR	deletion	GB 9857755
H	111	LEU	VAL	conflict	GB 9857755
H	115	SER	THR	conflict	GB 9857755
H	129	SER	CYS	conflict	GB 9857755
H	131	LYS	ARG	conflict	GB 9857755
H	135	GLY	GLU	conflict	GB 9857755
H	136	GLY	SER	conflict	GB 9857755
H	190	SER	ASN	conflict	GB 9857755
H	191	LEU	PHE	conflict	GB 9857755
H	197	ILE	THR	conflict	GB 9857755
H	201	ASN	ASP	conflict	GB 9857755
H	212	ARG	THR	conflict	GB 9857755

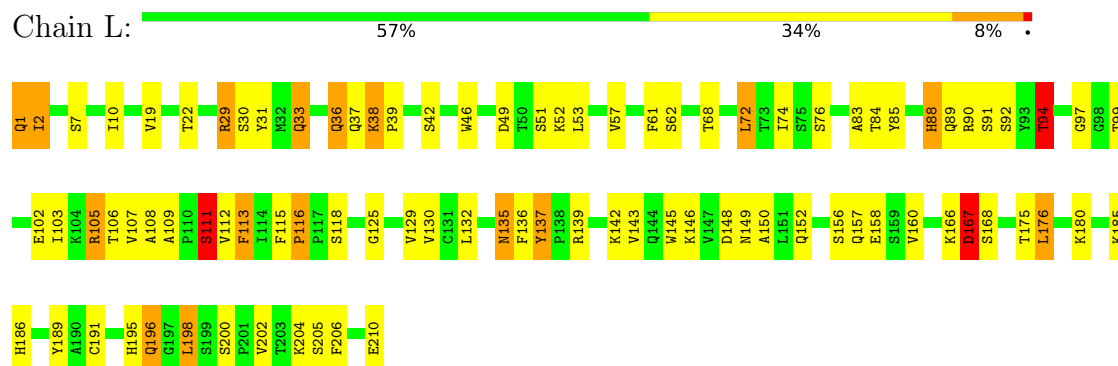
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	L	34	Total O 34 34	0	0
3	H	35	Total O 35 35	0	0

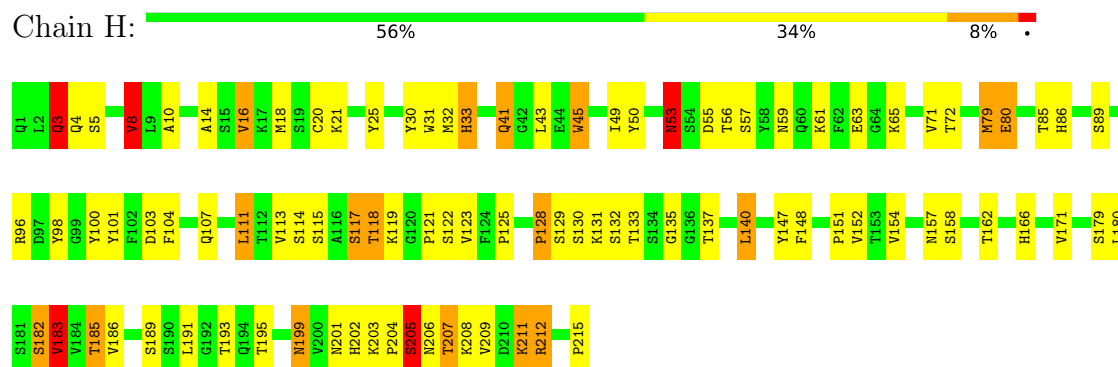
3 Residue-property plots

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: CHIMERIC SDZ CHI621



• Molecule 2: CHIMERIC SDZ CHI621



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	39.58Å 59.76Å 102.09Å 90.00° 99.98° 90.00°	Depositor
Resolution (Å)	8.00 – 2.60 8.00 – 2.60	Depositor EDS
% Data completeness (in resolution range)	(Not available) (8.00-2.60) 88.4 (8.00-2.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.15 (at 2.48Å)	Xtriage
Refinement program	X-PLOR	Depositor
R, R_{free}	0.196 , (Not available) 0.188 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	40.9	Xtriage
Anisotropy	0.443	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.19 , 59.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.028 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3303	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.84% 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	L	1.22	7/1639 (0.4%)	2.11	76/2220 (3.4%)
2	H	1.17	6/1675 (0.4%)	2.03	51/2281 (2.2%)
All	All	1.20	13/3314 (0.4%)	2.07	127/4501 (2.8%)

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	L	0	1

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	L	88	HIS	CD2-NE2	-7.06	1.30	1.37
2	H	202	HIS	CD2-NE2	-6.86	1.30	1.37
2	H	166	HIS	CD2-NE2	-6.32	1.30	1.37
1	L	195	HIS	CD2-NE2	-6.20	1.31	1.37
2	H	33	HIS	CD2-NE2	-6.19	1.31	1.37
2	H	85	THR	CA-CB	-6.09	1.44	1.53
1	L	202	VAL	CA-CB	-5.93	1.46	1.54
1	L	186	HIS	CD2-NE2	-5.78	1.31	1.37
1	L	107	VAL	CA-CB	-5.74	1.48	1.54
2	H	33	HIS	CG-ND1	-5.71	1.31	1.38
2	H	86	HIS	CD2-NE2	-5.50	1.31	1.37
1	L	108	ALA	CA-CB	-5.44	1.44	1.53
1	L	160	VAL	CA-CB	-5.01	1.47	1.55

All (127) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	103	ASP	CA-CB-CG	10.28	122.88	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	8	VAL	N-CA-CB	-9.46	99.34	111.64
1	L	36	GLN	OE1-CD-NE2	-9.07	113.53	122.60
1	L	30	SER	N-CA-C	8.69	120.36	111.07
1	L	167	ASP	CA-CB-CG	8.56	121.17	112.60
1	L	90	ARG	CA-CB-CG	-8.47	97.16	114.10
2	H	18	MET	O-C-N	-8.28	113.66	123.27
1	L	136	PHE	CA-CB-CG	8.27	122.07	113.80
1	L	2	ILE	N-CA-C	-8.19	96.78	108.58
1	L	74	ILE	N-CA-C	-8.14	96.76	108.48
2	H	53	ASN	CA-CB-CG	7.79	120.39	112.60
1	L	137	TYR	O-C-N	-7.69	115.84	121.84
1	L	90	ARG	N-CA-C	7.61	122.56	113.28
1	L	149	ASN	OD1-CG-ND2	-7.58	115.02	122.60
1	L	186	HIS	CB-CG-CD2	-7.45	121.52	131.20
1	L	22	THR	CA-CB-OG1	-7.39	98.51	109.60
1	L	72	LEU	N-CA-C	-7.38	96.29	108.76
2	H	53	ASN	CB-CG-ND2	7.31	127.37	116.40
2	H	104	PHE	CA-CB-CG	-7.25	106.55	113.80
1	L	130	VAL	O-C-N	-7.13	115.87	123.14
1	L	185	LYS	CB-CG-CD	-7.13	94.91	111.30
1	L	49	ASP	CA-CB-CG	7.11	119.71	112.60
2	H	5	SER	CA-CB-OG	7.03	125.16	111.10
2	H	182	SER	N-CA-C	-7.03	97.45	108.90
1	L	33	GLN	OE1-CD-NE2	-7.01	115.59	122.60
2	H	201	ASN	OD1-CG-ND2	-7.00	115.60	122.60
1	L	206	PHE	CA-CB-CG	6.99	120.79	113.80
1	L	118	SER	N-CA-C	-6.95	101.56	110.53
2	H	53	ASN	OD1-CG-ND2	-6.95	115.65	122.60
1	L	62	SER	N-CA-C	6.82	119.35	107.49
2	H	171	VAL	N-CA-C	-6.82	98.35	108.99
2	H	63	GLU	CA-CB-CG	-6.76	100.58	114.10
1	L	76	SER	CA-CB-OG	6.70	124.51	111.10
1	L	206	PHE	N-CA-C	-6.62	99.01	109.07
1	L	94	THR	N-CA-CB	-6.58	100.84	110.26
1	L	111	SER	O-C-N	-6.50	115.62	123.16
2	H	118	THR	CA-CB-OG1	-6.49	99.86	109.60
2	H	41	GLN	OE1-CD-NE2	-6.49	116.11	122.60
2	H	8	VAL	CB-CA-C	6.49	119.98	110.77
2	H	61	LYS	CA-CB-CG	-6.48	101.14	114.10
2	H	180	LEU	CD1-CG-CD2	-6.45	96.60	110.80
2	H	107	GLN	N-CA-C	6.44	120.59	112.87
1	L	176	LEU	N-CA-C	-6.42	98.74	109.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	71	VAL	CA-C-O	6.39	128.77	120.78
1	L	116	PRO	CA-C-N	6.37	126.32	119.76
1	L	116	PRO	C-N-CA	6.37	126.32	119.76
1	L	36	GLN	CG-CD-NE2	6.37	125.95	116.40
1	L	195	HIS	CB-CG-CD2	-6.32	122.98	131.20
1	L	1	GLN	OE1-CD-NE2	-6.32	116.28	122.60
2	H	166	HIS	CB-CG-CD2	-6.32	122.99	131.20
1	L	84	THR	N-CA-C	-6.26	99.50	109.59
2	H	55	ASP	CA-CB-CG	6.24	118.84	112.60
2	H	79	MET	O-C-N	-6.22	115.89	123.36
2	H	80	GLU	CA-CB-CG	6.22	126.55	114.10
1	L	97	GLY	N-CA-C	6.20	121.28	114.40
2	H	201	ASN	CB-CG-ND2	6.14	125.61	116.40
1	L	115	PHE	CA-C-N	6.10	126.66	120.38
1	L	115	PHE	C-N-CA	6.10	126.66	120.38
2	H	205	SER	N-CA-CB	6.08	120.76	110.49
2	H	212	ARG	CA-CB-CG	6.08	126.25	114.10
2	H	4	GLN	N-CA-C	6.05	117.70	110.19
2	H	212	ARG	N-CA-CB	6.00	118.78	109.97
1	L	196	GLN	CA-CB-CG	5.99	126.07	114.10
1	L	46	TRP	CG-CD1-NE1	-5.96	102.45	110.20
1	L	191	CYS	CA-C-O	-5.95	113.94	120.30
1	L	88	HIS	CB-CG-CD2	-5.88	123.56	131.20
1	L	196	GLN	N-CA-C	5.87	123.30	110.80
1	L	111	SER	N-CA-C	-5.86	100.11	109.96
1	L	186	HIS	CB-CG-ND1	5.85	131.48	122.70
1	L	76	SER	N-CA-C	-5.85	97.79	108.02
1	L	125	GLY	CA-C-O	-5.79	112.52	119.00
2	H	211	LYS	O-C-N	-5.79	116.50	123.27
2	H	31	TRP	CG-CD2-CE3	5.79	139.69	133.90
1	L	142	LYS	N-CA-C	-5.77	99.00	108.76
1	L	204	LYS	N-CA-C	-5.76	98.49	108.69
1	L	145	TRP	CG-CD1-NE1	-5.72	102.76	110.20
1	L	185	LYS	CA-CB-CG	5.71	125.51	114.10
1	L	196	GLN	CB-CA-C	-5.70	99.07	110.42
1	L	99	THR	N-CA-C	-5.68	97.97	107.80
1	L	62	SER	CA-CB-OG	-5.67	99.76	111.10
2	H	43	LEU	CD1-CG-CD2	-5.66	98.36	110.80
2	H	50	TYR	CA-C-O	-5.64	115.14	119.71
1	L	52	LYS	N-CA-C	-5.62	100.13	109.46
2	H	4	GLN	OE1-CD-NE2	-5.60	117.00	122.60
1	L	195	HIS	CB-CG-ND1	5.55	131.03	122.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L	115	PHE	CA-CB-CG	5.54	119.34	113.80
1	L	112	VAL	CA-C-O	-5.51	114.63	120.48
1	L	135	ASN	N-CA-C	5.50	122.52	110.80
2	H	205	SER	CB-CA-C	-5.47	99.54	110.42
2	H	205	SER	CA-CB-OG	5.46	122.02	111.10
2	H	212	ARG	CB-CA-C	-5.42	100.39	109.65
2	H	65	LYS	N-CA-C	5.41	119.83	112.04
1	L	102	GLU	CB-CG-CD	5.37	121.73	112.60
2	H	31	TRP	CB-CG-CD1	-5.35	118.87	126.90
1	L	22	THR	CA-CB-CG2	5.34	119.58	110.50
1	L	204	LYS	CG-CD-CE	-5.33	99.04	111.30
1	L	38	LYS	N-CA-C	-5.32	103.43	110.40
2	H	3	GLN	CB-CG-CD	5.32	121.65	112.60
1	L	176	LEU	CA-C-N	-5.32	115.49	122.99
1	L	176	LEU	C-N-CA	-5.32	115.49	122.99
2	H	208	LYS	N-CA-C	-5.31	99.45	108.26
2	H	183	VAL	N-CA-CB	-5.31	102.74	111.44
1	L	46	TRP	CG-CD2-CE3	5.27	139.17	133.90
2	H	45	TRP	CE2-CD2-CG	-5.27	100.87	107.20
2	H	199	ASN	OD1-CG-ND2	-5.27	117.33	122.60
1	L	46	TRP	CD1-CG-CD2	5.25	114.70	106.30
1	L	143	VAL	O-C-N	-5.20	117.77	123.18
2	H	31	TRP	CE2-CD2-CG	-5.20	100.96	107.20
1	L	46	TRP	CE2-CD2-CG	-5.19	100.97	107.20
1	L	200	SER	CA-C-N	5.18	125.34	119.90
1	L	200	SER	C-N-CA	5.18	125.34	119.90
2	H	89	SER	N-CA-C	-5.18	101.97	109.59
1	L	88	HIS	CB-CG-ND1	5.15	130.42	122.70
1	L	202	VAL	N-CA-CB	-5.09	103.28	110.26
2	H	59	ASN	N-CA-C	-5.09	103.32	110.50
1	L	33	GLN	CG-CD-NE2	5.06	124.00	116.40
2	H	157	ASN	OD1-CG-ND2	-5.06	117.54	122.60
2	H	119	LYS	CA-CB-CG	-5.06	103.98	114.10
1	L	148	ASP	CA-C-O	5.05	127.69	121.84
2	H	131	LYS	CB-CG-CD	-5.04	99.70	111.30
2	H	202	HIS	CB-CG-CD2	-5.04	124.65	131.20
1	L	113	PHE	N-CA-C	-5.03	102.75	110.14
1	L	168	SER	N-CA-C	5.03	118.72	112.58
2	H	57	SER	N-CA-C	-5.01	100.73	108.90
1	L	145	TRP	CD1-CG-CD2	5.01	114.31	106.30
1	L	145	TRP	CE2-CD2-CG	-5.01	101.19	107.20
1	L	157	GLN	O-C-N	-5.00	117.17	123.27

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	L	189	TYR	Sidechain

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	L	1604	0	1550	31	0
2	H	1630	0	1575	41	0
3	H	35	0	0	7	0
3	L	34	0	0	10	0
All	All	3303	0	3125	68	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 11.

All (68) 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:H:33:HIS:HD2	2:H:45:TRP:HE1	1.22	0.85
1:L:129:VAL:HA	3:L:237:HOH:O	1.80	0.81
1:L:38:LYS:HD3	1:L:83:ALA:HB2	1.66	0.77
2:H:20:CYS:SG	2:H:32:MET:HE3	2.28	0.73
2:H:158:SER:H	2:H:199:ASN:HD21	1.37	0.72
2:H:204:PRO:HB2	3:H:220:HOH:O	1.90	0.70
2:H:128:PRO:HG3	2:H:140:LEU:HG	1.74	0.69
2:H:128:PRO:HD2	2:H:215:PRO:HA	1.74	0.69
1:L:61:PHE:HD1	1:L:72:LEU:HD21	1.57	0.69
1:L:1:GLN:OE1	1:L:92:SER:HB3	1.96	0.66
2:H:33:HIS:CD2	2:H:45:TRP:HE1	2.10	0.65
2:H:158:SER:H	2:H:199:ASN:ND2	1.98	0.60
2:H:10:ALA:O	2:H:113:VAL:HA	2.04	0.58
2:H:25:TYR:CE2	2:H:96:ARG:HD2	2.39	0.58
1:L:33:GLN:NE2	2:H:101:TYR:HA	2.21	0.56
2:H:8:VAL:HG12	2:H:111:LEU:HD23	1.85	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:109:ALA:HB1	1:L:198:LEU:HD13	1.87	0.56
2:H:33:HIS:HD2	2:H:45:TRP:NE1	2.00	0.56
1:L:83:ALA:HB3	3:L:211:HOH:O	2.05	0.55
2:H:117:SER:HA	3:H:250:HOH:O	2.08	0.54
1:L:113:PHE:CE2	3:L:227:HOH:O	2.60	0.54
1:L:29:ARG:HD3	1:L:91:SER:OG	2.08	0.53
2:H:3:GLN:HG2	2:H:21:LYS:HB3	1.89	0.52
1:L:85:TYR:HE1	3:L:222:HOH:O	1.93	0.52
1:L:105:ARG:HD2	1:L:137:TYR:CB	2.40	0.52
1:L:132:LEU:HD22	2:H:183:VAL:HG21	1.92	0.51
1:L:36:GLN:HG3	3:L:211:HOH:O	2.10	0.50
2:H:79:MET:N	3:H:217:HOH:O	2.37	0.49
1:L:139:ARG:NH2	3:L:230:HOH:O	2.46	0.49
1:L:111:SER:C	3:L:227:HOH:O	2.55	0.49
2:H:140:LEU:HD11	2:H:186:VAL:HG22	1.93	0.49
2:H:114:SER:HG	2:H:148:PHE:HZ	1.58	0.49
1:L:156:SER:HA	1:L:175:THR:O	2.13	0.48
2:H:205:SER:HB2	2:H:207:THR:OG1	2.12	0.48
1:L:31:TYR:HB2	2:H:100:TYR:OH	2.14	0.48
1:L:88:HIS:HD2	1:L:94:THR:O	1.96	0.47
2:H:53:ASN:O	2:H:53:ASN:ND2	2.48	0.47
1:L:113:PHE:HD2	1:L:132:LEU:HD23	1.80	0.47
2:H:207:THR:HB	3:H:224:HOH:O	2.13	0.47
2:H:206:ASN:HB2	3:H:218:HOH:O	2.14	0.47
1:L:29:ARG:HA	1:L:29:ARG:HE	1.79	0.47
1:L:53:LEU:HB3	1:L:57:VAL:HB	1.96	0.46
1:L:61:PHE:CD1	1:L:72:LEU:HD21	2.43	0.46
2:H:96:ARG:HH21	2:H:98:TYR:HD1	1.63	0.46
1:L:37:GLN:O	1:L:83:ALA:HB1	2.16	0.46
2:H:203:LYS:O	2:H:205:SER:N	2.48	0.46
2:H:154:VAL:HG11	2:H:182:SER:HB2	1.98	0.45
2:H:49:ILE:HG13	2:H:56:THR:HG22	1.98	0.45
1:L:2:ILE:O	1:L:94:THR:HG21	2.17	0.45
1:L:146:LYS:HA	1:L:150:ALA:O	2.19	0.43
2:H:8:VAL:HG13	2:H:16:VAL:HG21	2.00	0.43
1:L:113:PHE:CD2	3:L:227:HOH:O	2.57	0.43
2:H:25:TYR:CD2	2:H:30:TYR:HD2	2.36	0.43
1:L:166:LYS:O	1:L:167:ASP:HB2	2.19	0.42
1:L:33:GLN:HE22	2:H:101:TYR:HA	1.83	0.42
2:H:128:PRO:HG3	2:H:140:LEU:CG	2.45	0.42
2:H:25:TYR:HD2	2:H:30:TYR:CD2	2.38	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:121:PRO:HB3	2:H:147:TYR:HB3	2.02	0.42
2:H:125:PRO:HD3	2:H:211:LYS:HE2	2.01	0.42
1:L:116:PRO:C	3:L:237:HOH:O	2.62	0.42
2:H:25:TYR:CD2	2:H:30:TYR:CD2	3.08	0.42
2:H:123:VAL:HG21	2:H:209:VAL:HG21	2.02	0.41
1:L:37:GLN:HA	3:L:240:HOH:O	2.21	0.41
2:H:118:THR:N	3:H:250:HOH:O	2.54	0.41
2:H:137:THR:HG23	2:H:185:THR:HG23	2.02	0.41
2:H:154:VAL:HG11	2:H:182:SER:CB	2.51	0.41
2:H:205:SER:O	2:H:206:ASN:HB3	2.20	0.41
1:L:33:GLN:HG3	3:H:234:HOH:O	2.22	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	L	208/210 (99%)	194 (93%)	10 (5%)	4 (2%)	6	13
2	H	213/215 (99%)	189 (89%)	18 (8%)	6 (3%)	4	6
All	All	421/425 (99%)	383 (91%)	28 (7%)	10 (2%)	4	9

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	L	167	ASP
2	H	130	SER
2	H	205	SER
1	L	135	ASN
1	L	196	GLN
1	L	51	SER
2	H	132	SER

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Mol	Chain	Res	Type
2	H	14	ALA
2	H	128	PRO
2	H	135	GLY

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	L	182/182 (100%)	162 (89%)	20 (11%)	6	13
2	H	181/181 (100%)	155 (86%)	26 (14%)	3	6
All	All	363/363 (100%)	317 (87%)	46 (13%)	4	9

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	7	SER
1	L	10	ILE
1	L	19	VAL
1	L	29	ARG
1	L	39	PRO
1	L	42	SER
1	L	68	THR
1	L	89	GLN
1	L	94	THR
1	L	103	ILE
1	L	105	ARG
1	L	106	THR
1	L	111	SER
1	L	152	GLN
1	L	158	GLU
1	L	176	LEU
1	L	180	LYS
1	L	198	LEU
1	L	205	SER
1	L	210	GLU

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Mol	Chain	Res	Type
2	H	3	GLN
2	H	8	VAL
2	H	16	VAL
2	H	41	GLN
2	H	53	ASN
2	H	72	THR
2	H	80	GLU
2	H	111	LEU
2	H	115	SER
2	H	117	SER
2	H	122	SER
2	H	129	SER
2	H	133	THR
2	H	140	LEU
2	H	151	PRO
2	H	152	VAL
2	H	162	THR
2	H	179	SER
2	H	183	VAL
2	H	185	THR
2	H	189	SER
2	H	191	LEU
2	H	193	THR
2	H	195	THR
2	H	207	THR
2	H	212	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	157	GLN
1	L	196	GLN
1	L	207	ASN
2	H	33	HIS
2	H	199	ASN
2	H	201	ASN

5.3.3 RNA ⓘ

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	L	210/210 (100%)	-0.83	0 100 100	4, 28, 58, 92	0
2	H	215/215 (100%)	-0.74	0 100 100	5, 29, 58, 104	0
All	All	425/425 (100%)	-0.78	0 100 100	4, 28, 58, 104	0

There are no RSRZ outliers to report.

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.