



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 8, 2026 – 04:43 PM UTC

PDB ID : 3NH7 / pdb_00003nh7
Title : Crystal structure of the neutralizing Fab fragment AbD1556 bound to the BMP type I receptor IA
Authors : Mueller, T.D.; Harth, S.; Sebald, W.
Deposited on : 2010-06-14
Resolution : 2.70 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

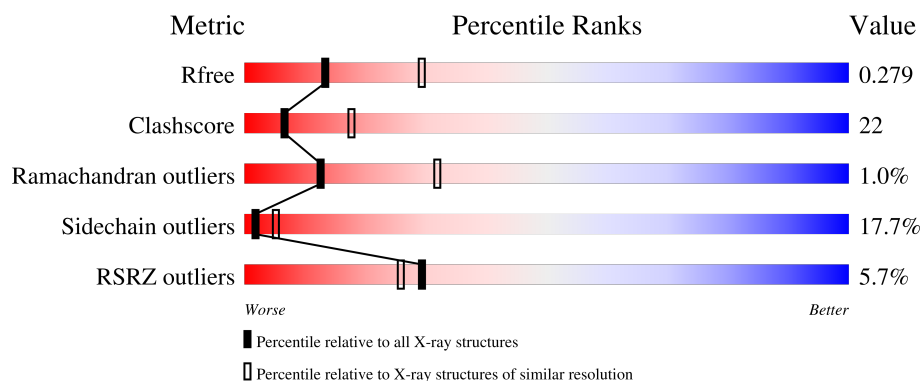
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	H	234	3% 44% 34% 9% 12%
1	I	234	% 56% 27% 8% 8%
1	J	234	4% 57% 29% 7% 6%
1	K	234	3% 57% 26% 8% 8%
2	L	213	8% 60% 31% 8%

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Mol	Chain	Length	Quality of chain			
2	M	213	3%	58%	32%	8%
2	N	213	6%	62%	29%	8%
2	O	213	2%	60%	29%	10%
3	A	129	8%	40%	19%	7%
3	B	129	9%	38%	21%	7%
3	C	129	10%	40%	18%	7%
3	D	129	10%	38%	20%	7%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 15350 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 Antibody fragment Fab AbD1556, heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	H	207	Total	C	N	O	S	0	0	0
			1569	992	269	303	5			
1	I	215	Total	C	N	O	S	0	0	0
			1620	1021	278	316	5			
1	J	221	Total	C	N	O	S	0	0	0
			1661	1045	286	325	5			
1	K	215	Total	C	N	O	S	0	0	0
			1625	1025	279	316	5			

- Molecule 2 is a protein called Antibody fragment Fab AbD1556, light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	L	210	Total	C	N	O	S	0	0	0
			1558	973	258	322	5			
2	M	210	Total	C	N	O	S	0	0	0
			1558	973	258	322	5			
2	N	210	Total	C	N	O	S	0	0	0
			1558	973	258	322	5			
2	O	210	Total	C	N	O	S	0	0	0
			1558	973	258	322	5			

- Molecule 3 is a protein called Bone morphogenetic protein receptor type-1A.

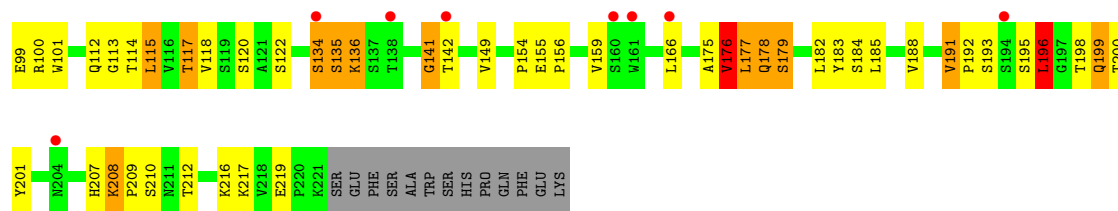
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	A	85	Total	C	N	O	S	0	0	0
			654	400	112	131	11			
3	B	85	Total	C	N	O	S	0	0	0
			654	400	112	131	11			
3	C	85	Total	C	N	O	S	0	0	0
			654	400	112	131	11			
3	D	85	Total	C	N	O	S	0	0	0
			654	400	112	131	11			

- Molecule 4 is water.

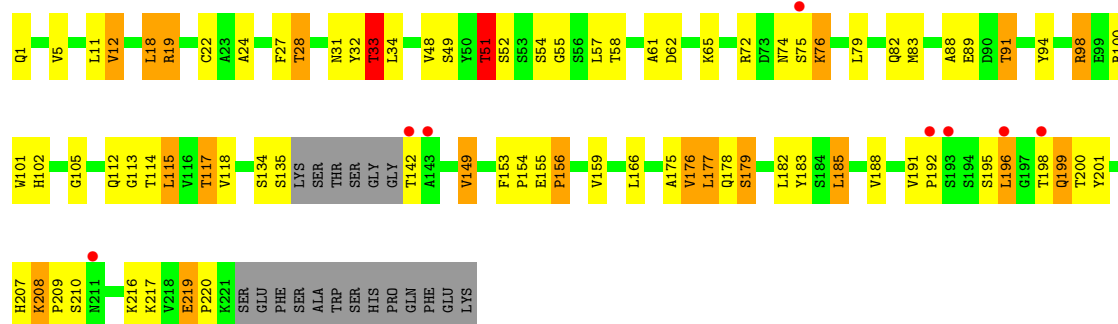
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	H	2	Total 2	O 2	0	0
4	L	3	Total 3	O 3	0	0
4	A	2	Total 2	O 2	0	0
4	I	5	Total 5	O 5	0	0
4	M	4	Total 4	O 4	0	0
4	B	1	Total 1	O 1	0	0
4	J	2	Total 2	O 2	0	0
4	N	1	Total 1	O 1	0	0
4	K	2	Total 2	O 2	0	0
4	O	4	Total 4	O 4	0	0
4	D	1	Total 1	O 1	0	0

- Molecule 1: Antibody fragment Fab AbD1556, heavy chain

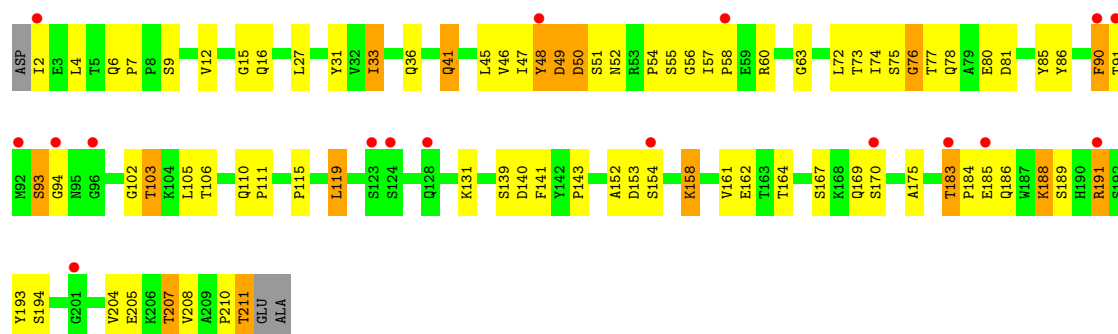




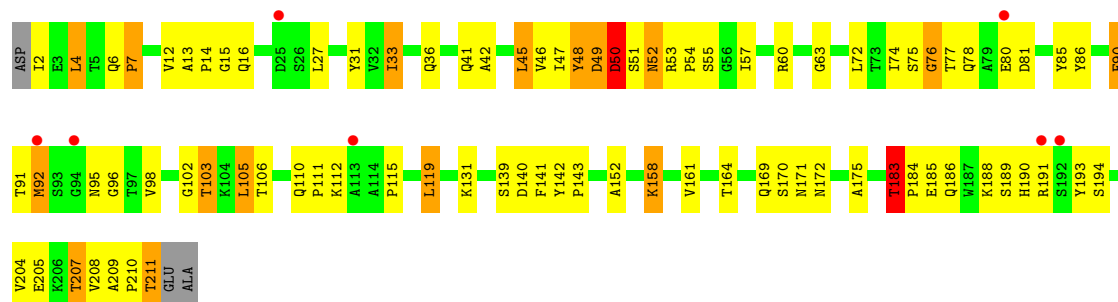
• Molecule 1: Antibody fragment Fab AbD1556, heavy chain



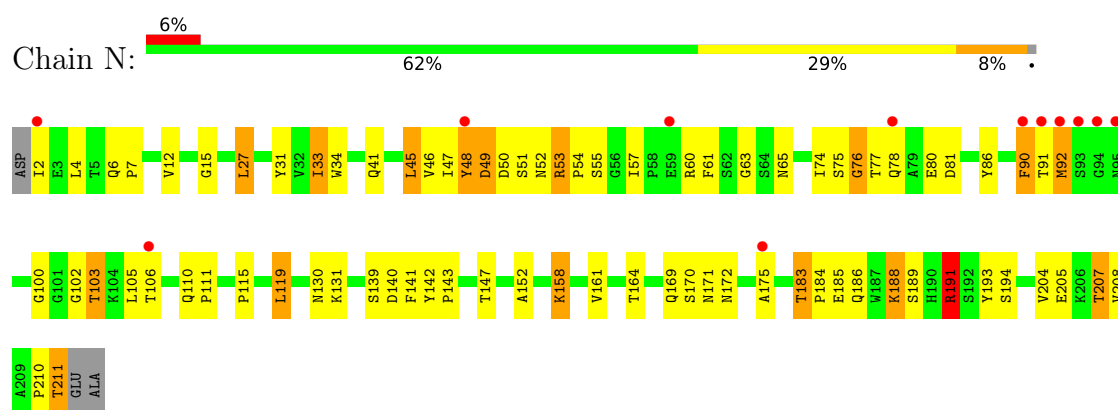
• Molecule 2: Antibody fragment Fab AbD1556, light chain



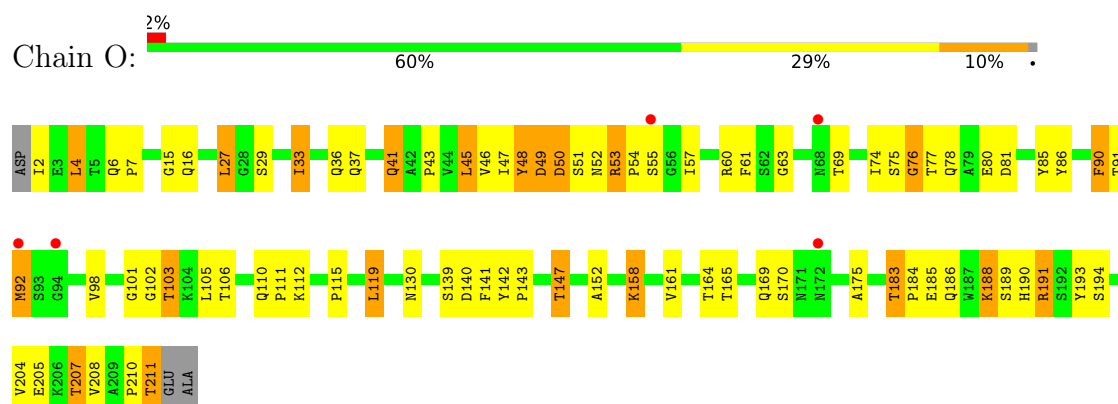
• Molecule 2: Antibody fragment Fab AbD1556, light chain



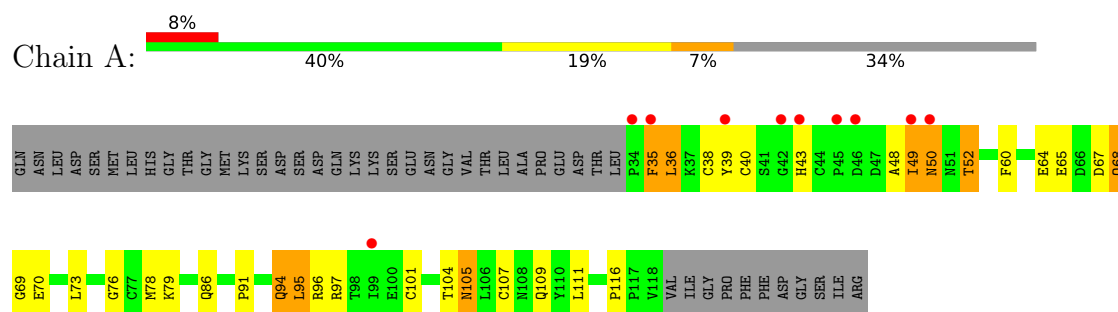
• Molecule 2: Antibody fragment Fab AbD1556, light chain



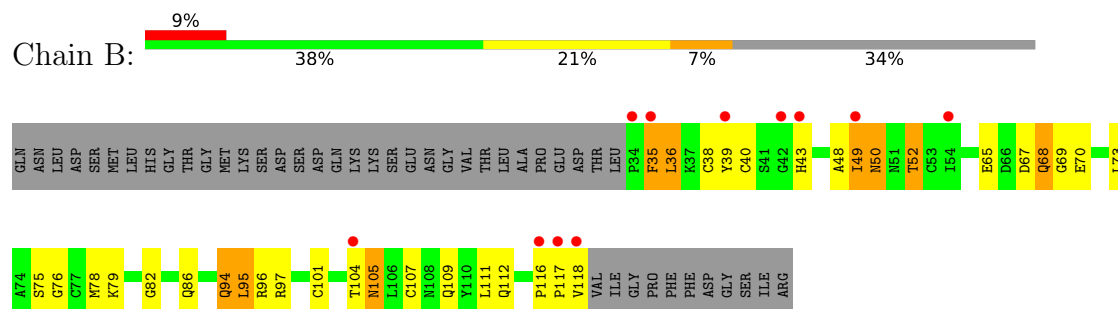
- Molecule 2: Antibody fragment Fab AbD1556, light chain



- Molecule 3: Bone morphogenetic protein receptor type-1A

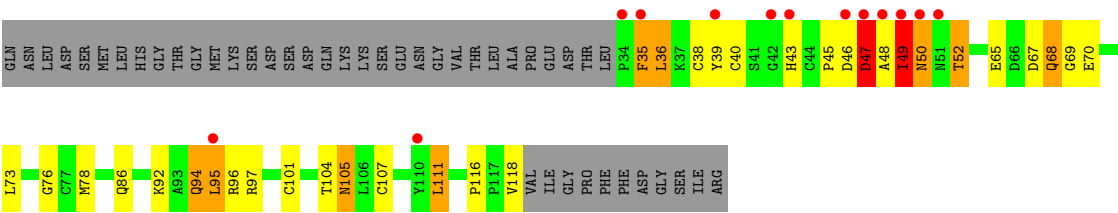


- Molecule 3: Bone morphogenetic protein receptor type-1A

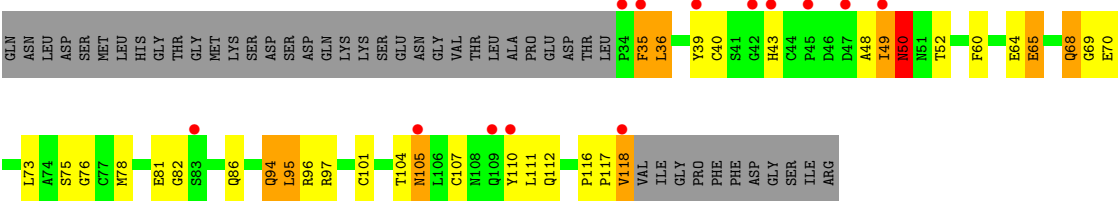


- Molecule 3: Bone morphogenetic protein receptor type-1A





● Molecule 3: Bone morphogenetic protein receptor type-1A



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	89.32Å 129.25Å 100.24Å 90.00° 92.27° 90.00°	Depositor
Resolution (Å)	30.40 – 2.70 30.40 – 2.70	Depositor EDS
% Data completeness (in resolution range)	95.3 (30.40-2.70) 95.3 (30.40-2.70)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.65 (at 2.68Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.234 , 0.280 0.235 , 0.279	Depositor DCC
R_{free} test set	3020 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	50.5	Xtriage
Anisotropy	0.040	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 64.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.025 for h,-k,-l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	15350	wwPDB-VP
Average B, all atoms (Å ²)	64.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.33% 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	H	0.82	0/1608	1.14	10/2190 (0.5%)
1	I	0.78	0/1660	1.08	5/2262 (0.2%)
1	J	0.73	0/1702	1.11	8/2318 (0.3%)
1	K	0.87	1/1665 (0.1%)	1.15	9/2268 (0.4%)
2	L	0.84	1/1595 (0.1%)	1.20	11/2180 (0.5%)
2	M	0.79	0/1595	1.18	20/2180 (0.9%)
2	N	0.84	0/1595	1.23	8/2180 (0.4%)
2	O	0.83	0/1595	1.16	14/2180 (0.6%)
3	A	0.68	0/667	0.97	1/903 (0.1%)
3	B	0.61	0/667	0.94	1/903 (0.1%)
3	C	0.69	0/667	1.06	4/903 (0.4%)
3	D	0.63	0/667	0.98	2/903 (0.2%)
All	All	0.79	2/15683 (0.0%)	1.13	93/21370 (0.4%)

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	L	0	1
3	A	0	1
3	B	0	1
3	C	0	1
3	D	0	1
All	All	0	5

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	L	191	ARG	CZ-NH2	5.88	1.41	1.33
1	K	33	THR	CB-CG2	-5.34	1.34	1.52

The worst 5 of 93 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	N	191	ARG	NE-CZ-NH1	-15.42	106.08	121.50
2	L	191	ARG	NE-CZ-NH2	14.29	132.06	119.20
2	O	49	ASP	N-CA-C	-12.41	90.45	109.94
2	N	49	ASP	N-CA-C	-11.72	91.53	109.94
2	N	191	ARG	NE-CZ-NH2	11.50	129.55	119.20

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	A	48	ALA	Peptide
3	B	48	ALA	Peptide
3	C	48	ALA	Peptide
3	D	48	ALA	Peptide
2	L	93	SER	Peptide

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	H	1569	0	1526	95	0
1	I	1620	0	1576	68	0
1	J	1661	0	1623	63	0
1	K	1625	0	1586	57	0
2	L	1558	0	1507	70	0
2	M	1558	0	1507	78	0
2	N	1558	0	1507	75	0
2	O	1558	0	1507	69	0
3	A	654	0	608	41	0
3	B	654	0	608	40	0
3	C	654	0	608	40	0
3	D	654	0	608	41	0
4	A	2	0	0	0	0
4	B	1	0	0	1	0
4	D	1	0	0	0	0
4	H	2	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	I	5	0	0	1	0
4	J	2	0	0	0	0
4	K	2	0	0	0	0
4	L	3	0	0	0	0
4	M	4	0	0	2	0
4	N	1	0	0	0	0
4	O	4	0	0	0	0
All	All	15350	0	14771	651	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 22.

The worst 5 of 651 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:O:52:ASN:ND2	3:D:43:HIS:CD2	1.99	1.31
2:L:52:ASN:ND2	3:A:43:HIS:CD2	2.07	1.22
1:I:33:THR:CG2	1:I:52:SER:HA	1.70	1.20
2:O:52:ASN:ND2	3:D:43:HIS:HD2	1.37	1.18
2:M:52:ASN:ND2	3:B:43:HIS:CD2	2.12	1.17

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	H	201/234 (86%)	171 (85%)	24 (12%)	6 (3%)	3 8
1	I	211/234 (90%)	202 (96%)	8 (4%)	1 (0%)	24 48
1	J	219/234 (94%)	205 (94%)	13 (6%)	1 (0%)	24 48
1	K	211/234 (90%)	201 (95%)	10 (5%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	L	208/213 (98%)	199 (96%)	7 (3%)	2 (1%)	12	32
2	M	208/213 (98%)	200 (96%)	7 (3%)	1 (0%)	24	48
2	N	208/213 (98%)	202 (97%)	5 (2%)	1 (0%)	24	48
2	O	208/213 (98%)	202 (97%)	5 (2%)	1 (0%)	24	48
3	A	83/129 (64%)	74 (89%)	8 (10%)	1 (1%)	10	27
3	B	83/129 (64%)	72 (87%)	9 (11%)	2 (2%)	4	12
3	C	83/129 (64%)	74 (89%)	7 (8%)	2 (2%)	4	12
3	D	83/129 (64%)	74 (89%)	7 (8%)	2 (2%)	4	12
All	All	2006/2304 (87%)	1876 (94%)	110 (6%)	20 (1%)	12	32

5 of 20 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	H	123	THR
1	H	192	PRO
2	L	76	GLY
2	L	94	GLY
2	M	76	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	H	175/198 (88%)	143 (82%)	32 (18%)	2	5
1	I	181/198 (91%)	146 (81%)	35 (19%)	1	4
1	J	186/198 (94%)	149 (80%)	37 (20%)	1	4
1	K	182/198 (92%)	145 (80%)	37 (20%)	1	3
2	L	176/178 (99%)	148 (84%)	28 (16%)	2	7
2	M	176/178 (99%)	149 (85%)	27 (15%)	3	7
2	N	176/178 (99%)	149 (85%)	27 (15%)	3	7
2	O	176/178 (99%)	148 (84%)	28 (16%)	2	7

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
3	A	76/114 (67%)	64 (84%)	12 (16%)	2	7
3	B	76/114 (67%)	61 (80%)	15 (20%)	1	4
3	C	76/114 (67%)	63 (83%)	13 (17%)	2	6
3	D	76/114 (67%)	61 (80%)	15 (20%)	1	4
All	All	1732/1960 (88%)	1426 (82%)	306 (18%)	2	5

5 of 306 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	K	22	CYS
2	O	191	ARG
1	K	65	LYS
1	K	208	LYS
3	D	75	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 79 such sidechains are listed below:

Mol	Chain	Res	Type
3	C	86	GLN
2	O	65	ASN
3	C	105	ASN
1	K	162	ASN
3	D	43	HIS

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 [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 ⓘ

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	H	207/234 (88%)	0.19	8 (3%) 43 39	37, 54, 91, 122	0
1	I	215/234 (91%)	0.22	2 (0%) 81 80	38, 56, 89, 114	0
1	J	221/234 (94%)	0.57	10 (4%) 38 34	44, 64, 102, 125	0
1	K	215/234 (91%)	0.35	8 (3%) 45 41	37, 54, 95, 123	0
2	L	210/213 (98%)	0.73	17 (8%) 18 15	44, 61, 86, 112	0
2	M	210/213 (98%)	0.42	7 (3%) 49 45	41, 59, 77, 92	0
2	N	210/213 (98%)	0.48	12 (5%) 29 26	48, 59, 78, 93	0
2	O	210/213 (98%)	0.38	5 (2%) 59 56	44, 57, 72, 86	0
3	A	85/129 (65%)	0.82	10 (11%) 9 7	51, 73, 107, 127	0
3	B	85/129 (65%)	0.96	11 (12%) 7 6	52, 79, 112, 135	0
3	C	85/129 (65%)	1.10	13 (15%) 5 4	58, 78, 108, 127	0
3	D	85/129 (65%)	1.10	13 (15%) 5 4	50, 80, 131, 157	0
All	All	2038/2304 (88%)	0.52	116 (5%) 29 26	37, 61, 99, 157	0

The worst 5 of 116 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	C	34	PRO	5.1
3	C	43	HIS	5.1
3	C	95	LEU	4.7
2	N	94	GLY	4.7
3	D	39	TYR	4.5

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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.