



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 04:19 PM UTC

PDB ID : 1ZVK / pdb_00001zvk
Title : Structure of Double mutant, D164N, E78H of Kumamolisin-As
Authors : Li, M.; Wlodawer, A.; Gustchina, A.; Nakayama, T.
Deposited on : 2005-06-02
Resolution : 2.04 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

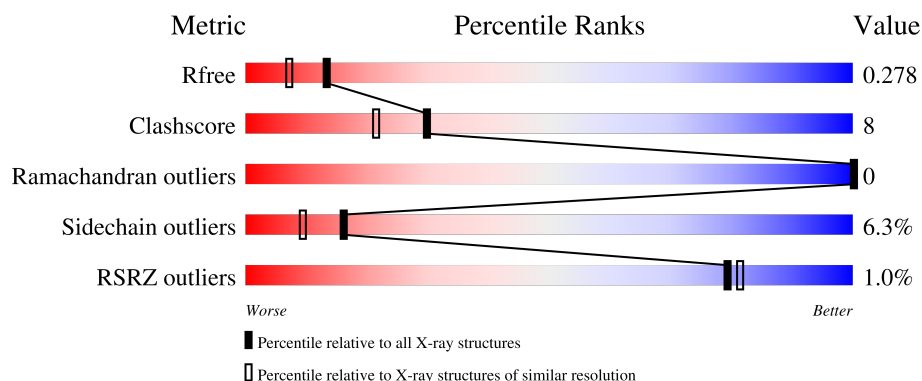
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.04 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	2260 (2.04-2.04)
Clashscore	190562	2333 (2.04-2.04)
Ramachandran outliers	187476	2318 (2.04-2.04)
Sidechain outliers	187428	2318 (2.04-2.04)
RSRZ outliers	180081	2260 (2.04-2.04)

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	358	% 72% 25% ..
1	B	358	% 77% 20% ..

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5502 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 kumamolisin-As.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	355	Total	C	N	O	S	0	0	1
			2528	1593	426	505	4			
1	B	355	Total	C	N	O	S	0	0	1
			2528	1593	426	505	4			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	78	HIS	GLU	engineered mutation	UNP Q8GB88
A	164	ASN	ASP	engineered mutation	UNP Q8GB88
B	78	HIS	GLU	engineered mutation	UNP Q8GB88
B	164	ASN	ASP	engineered mutation	UNP Q8GB88

- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Ca	0	0
			1	1		
2	B	1	Total	Ca	0	0
			1	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	207	Total	O	0	0
			207	207		
3	B	237	Total	O	0	0
			237	237		

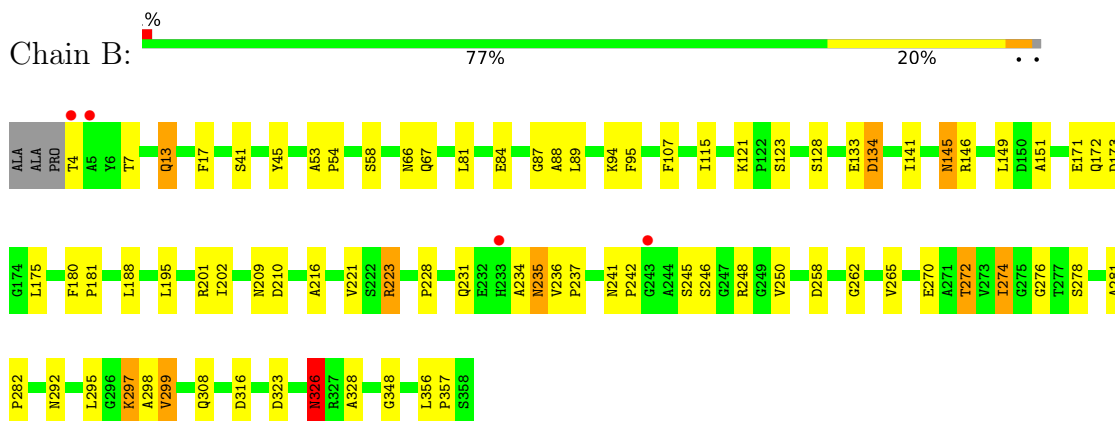
3 Residue-property plots

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: kumamolisin-As



• Molecule 1: kumamolisin-As



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	58.05Å 74.81Å 78.38Å 90.00° 103.08° 90.00°	Depositor
Resolution (Å)	10.00 – 2.04 10.00 – 2.04	Depositor EDS
% Data completeness (in resolution range)	93.9 (10.00-2.04) 95.9 (10.00-2.04)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.47 (at 2.05Å)	Xtriage
Refinement program	SHELXL-97	Depositor
R, R_{free}	0.188 , 0.292 0.189 , 0.278	Depositor DCC
R_{free} test set	2066 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	20.0	Xtriage
Anisotropy	0.433	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.50 , 92.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5502	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 20.73 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 8.1542e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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

Bond lengths and bond angles in the following residue types are not validated in this section:
CA

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	0.60	0/2592	1.79	49/3562 (1.4%)
1	B	0.60	0/2592	1.66	29/3562 (0.8%)
All	All	0.60	0/5184	1.73	78/7124 (1.1%)

There are no bond length outliers.

All (78) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	261	THR	CA-C-N	12.06	136.73	121.06
1	A	261	THR	C-N-CA	12.06	136.73	121.06
1	A	107	PHE	CA-CB-CG	9.68	123.48	113.80
1	A	121	LYS	O-C-N	-8.85	117.46	121.53
1	A	145	ASN	CA-CB-CG	-8.11	104.49	112.60
1	B	134	ASP	CA-CB-CG	8.08	120.68	112.60
1	B	221	VAL	CA-C-O	7.87	128.22	120.27
1	A	232	GLU	CA-C-N	-7.68	109.78	121.53
1	A	232	GLU	C-N-CA	-7.68	109.78	121.53
1	B	128	SER	O-C-N	7.28	131.01	121.99
1	A	330	ILE	CA-C-N	7.27	134.79	121.70
1	A	330	ILE	C-N-CA	7.27	134.79	121.70
1	B	121	LYS	CB-CA-C	-7.15	106.12	111.20
1	A	312	ASP	CA-CB-CG	7.10	119.70	112.60
1	A	243	GLY	N-CA-C	-6.98	105.57	114.37
1	B	66	ASN	CA-C-N	-6.96	114.60	123.16
1	B	66	ASN	C-N-CA	-6.96	114.60	123.16
1	A	243	GLY	O-C-N	6.86	129.96	122.56
1	A	277	THR	CA-C-N	6.84	130.14	120.28
1	A	277	THR	C-N-CA	6.84	130.14	120.28
1	A	232	GLU	O-C-N	-6.81	113.54	122.59
1	B	316	ASP	CA-CB-CG	6.79	119.39	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	173	ASP	CA-C-N	6.76	134.19	120.80
1	A	173	ASP	C-N-CA	6.76	134.19	120.80
1	A	105	ALA	CA-C-N	6.72	127.27	119.94
1	A	105	ALA	C-N-CA	6.72	127.27	119.94
1	B	123	SER	CA-C-N	6.71	132.48	122.75
1	B	123	SER	C-N-CA	6.71	132.48	122.75
1	A	66	ASN	CA-C-N	-6.50	114.42	122.42
1	A	66	ASN	C-N-CA	-6.50	114.42	122.42
1	B	107	PHE	CA-CB-CG	6.50	120.30	113.80
1	A	104	ASP	CA-CB-CG	6.20	118.80	112.60
1	A	12	ALA	CA-C-O	6.20	127.10	120.10
1	A	261	THR	O-C-N	6.18	129.65	121.99
1	B	250	VAL	CA-C-O	-6.10	111.77	119.95
1	B	88	ALA	CA-C-N	-6.03	112.67	122.26
1	B	88	ALA	C-N-CA	-6.03	112.67	122.26
1	B	221	VAL	O-C-N	-5.96	116.99	123.18
1	A	95	PHE	CA-CB-CG	-5.91	107.89	113.80
1	B	235	ASN	CA-CB-CG	5.88	118.48	112.60
1	B	146	ARG	CA-C-O	5.88	126.73	119.97
1	B	7	THR	N-CA-CB	5.85	116.52	109.74
1	A	262	GLY	N-CA-C	5.81	120.03	112.25
1	B	41	SER	CA-C-O	5.75	126.51	120.42
1	A	74	GLY	CA-C-O	-5.71	113.35	121.52
1	B	276	GLY	N-CA-C	-5.71	106.53	111.95
1	A	55	GLN	CB-CG-CD	5.70	122.29	112.60
1	A	354	ALA	CA-C-O	5.69	126.52	119.97
1	A	146	ARG	CA-C-O	5.67	126.56	120.55
1	B	94	LYS	CA-C-O	-5.65	114.61	121.05
1	B	262	GLY	N-CA-C	5.64	121.17	111.98
1	A	121	LYS	CA-C-O	5.62	124.75	120.26
1	A	67	GLN	O-C-N	5.57	125.73	121.27
1	A	21	LEU	CA-C-O	-5.57	114.14	120.32
1	B	13	GLN	CA-C-O	5.52	126.40	120.55
1	A	41	SER	CA-C-O	5.51	126.38	120.70
1	A	186	TYR	CA-C-O	5.50	125.99	119.56
1	B	173	ASP	N-CA-C	-5.46	104.89	113.02
1	A	64	ALA	O-C-N	-5.45	116.64	122.85
1	A	71	ASP	CA-C-O	5.42	122.81	119.29
1	A	190	CYS	CA-C-N	5.36	129.30	121.05
1	A	190	CYS	C-N-CA	5.36	129.30	121.05
1	B	348	GLY	N-CA-C	5.27	119.26	112.83
1	B	145	ASN	CA-CB-CG	-5.21	107.39	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	125	VAL	CA-C-O	-5.20	115.02	120.27
1	B	326	ASN	CA-C-N	-5.18	114.36	122.65
1	B	326	ASN	C-N-CA	-5.18	114.36	122.65
1	A	100	ALA	O-C-N	5.16	125.58	121.55
1	A	122	PRO	CA-C-N	-5.14	114.06	122.54
1	A	122	PRO	C-N-CA	-5.14	114.06	122.54
1	A	141	ILE	CA-C-N	5.09	127.52	120.29
1	A	141	ILE	C-N-CA	5.09	127.52	120.29
1	A	125	VAL	O-C-N	5.05	128.43	123.18
1	B	17	PHE	CA-CB-CG	5.05	118.85	113.80
1	A	166	GLY	O-C-N	5.01	128.21	123.05
1	A	113	THR	CA-C-N	5.01	127.26	120.65
1	A	113	THR	C-N-CA	5.01	127.26	120.65
1	B	236	VAL	CA-C-O	5.01	123.58	119.47

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2528	0	2418	41	0
1	B	2528	0	2418	44	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	207	0	0	8	0
3	B	237	0	0	12	0
All	All	5502	0	4836	84	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 8.

All (84) 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:B:295:LEU:HD12	1:B:299:VAL:HG12	1.55	0.85
1:B:326:ASN:HD22	1:B:328:ALA:H	1.22	0.85
1:B:292:ASN:OD1	1:B:299:VAL:HG13	1.82	0.79
1:B:265:VAL:HG21	1:B:274:ILE:HD13	1.67	0.77
1:B:81:LEU:HD13	1:B:274:ILE:HD11	1.70	0.74
1:B:242:PRO:HG2	3:B:1478:HOH:O	1.91	0.70
1:A:292:ASN:OD1	1:A:299:VAL:HG13	1.93	0.69
1:B:357:PRO:HD3	3:B:1166:HOH:O	1.93	0.67
1:B:228:PRO:HD2	1:B:231:GLN:HB2	1.79	0.64
1:B:145:ASN:HB2	3:B:1319:HOH:O	1.99	0.62
1:A:142:ALA:O	1:A:146:ARG:HG3	2.00	0.61
1:A:24:GLN:NE2	3:A:1362:HOH:O	2.32	0.61
1:A:152:ALA:HB1	3:A:1258:HOH:O	1.99	0.61
1:A:212:PRO:HG3	1:A:324:ILE:O	2.02	0.60
1:B:326:ASN:ND2	1:B:328:ALA:H	1.97	0.60
1:A:195:LEU:HD11	1:A:202:ILE:HG12	1.85	0.59
1:B:326:ASN:HD22	1:B:328:ALA:N	1.97	0.57
1:B:270:GLU:HG2	3:B:1341:HOH:O	2.04	0.57
1:B:175:LEU:HD13	1:B:223:ARG:HH21	1.69	0.57
1:B:223:ARG:HD3	3:B:2011:HOH:O	2.04	0.57
1:A:81:LEU:HD13	1:A:274:ILE:HD11	1.90	0.54
1:B:180:PHE:CG	1:B:181:PRO:HA	2.43	0.54
1:A:84:GLU:OE1	1:A:265:VAL:HB	2.08	0.54
1:A:294:LYS:HG2	1:A:352:LEU:HD21	1.90	0.54
1:A:5:ALA:HB1	1:A:264:GLU:HG3	1.90	0.53
1:B:87:GLY:HA2	1:B:95:PHE:CZ	2.44	0.52
1:A:81:LEU:HD13	1:A:274:ILE:CD1	2.40	0.52
1:A:87:GLY:HA2	1:A:95:PHE:CZ	2.44	0.52
1:B:228:PRO:HD3	1:B:248:ARG:HA	1.92	0.52
1:A:180:PHE:CG	1:A:181:PRO:HA	2.45	0.52
1:B:210:ASP:OD2	1:B:258:ASP:OD1	2.29	0.50
1:A:201:ARG:HA	3:A:1122:HOH:O	2.12	0.50
1:A:11:VAL:HG21	1:A:263:TYR:CE2	2.47	0.50
1:B:115:ILE:HD13	1:B:151:ALA:HB2	1.94	0.49
1:B:308:GLN:HB2	3:B:1294:HOH:O	2.12	0.49
1:B:134:ASP:OD1	1:B:171:GLU:OE1	2.30	0.49
1:B:270:GLU:O	1:B:272:THR:HG22	2.13	0.49
1:B:115:ILE:CD1	1:B:151:ALA:HB2	2.42	0.49
1:B:45:TYR:CE1	1:B:265:VAL:HA	2.48	0.49
1:B:81:LEU:HB2	1:B:274:ILE:HD11	1.96	0.48
1:B:81:LEU:HB2	1:B:274:ILE:CD1	2.44	0.48
1:B:195:LEU:HD11	1:B:202:ILE:HG12	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:298:ALA:HB3	3:B:1506:HOH:O	2.14	0.47
1:A:45:TYR:CE1	1:A:265:VAL:HA	2.50	0.47
1:A:337:TRP:HA	1:A:342:GLY:O	2.14	0.47
1:A:81:LEU:HD13	1:A:274:ILE:HG13	1.97	0.47
1:A:326:ASN:OD1	1:A:328:ALA:HB3	2.14	0.46
1:B:53:ALA:HA	1:B:54:PRO:HD3	1.86	0.46
1:A:265:VAL:HG21	1:A:274:ILE:CD1	2.46	0.46
1:B:265:VAL:CG2	1:B:274:ILE:HD13	2.44	0.46
1:A:23:GLY:O	1:A:26:GLN:HB2	2.16	0.46
1:A:45:TYR:CE2	1:A:84:GLU:HB3	2.50	0.46
1:B:13:GLN:NE2	3:B:1190:HOH:O	2.49	0.46
1:A:263:TYR:O	1:A:273:VAL:HA	2.16	0.45
1:A:303:ASN:HB2	1:A:304:PRO:HD3	1.99	0.45
1:B:223:ARG:NE	1:B:241:ASN:HD22	2.14	0.45
1:A:295:LEU:CD1	1:A:299:VAL:HG12	2.45	0.45
1:B:323:ASP:OD2	1:B:326:ASN:ND2	2.49	0.45
1:A:13:GLN:NE2	3:A:1330:HOH:O	2.50	0.45
1:A:327:ARG:HA	3:A:1017:HOH:O	2.17	0.45
1:B:45:TYR:CE2	1:B:84:GLU:HB3	2.52	0.44
1:A:43:ALA:HB2	3:A:1158:HOH:O	2.16	0.44
1:A:115:ILE:O	1:B:201:ARG:NE	2.48	0.44
1:A:352:LEU:O	1:A:356:LEU:HG	2.17	0.44
1:A:100:ALA:HB1	1:A:101:PRO:HD2	1.99	0.44
1:B:54:PRO:HD2	3:B:1183:HOH:O	2.17	0.44
1:B:67:GLN:NE2	3:B:1151:HOH:O	2.50	0.43
1:A:223:ARG:NE	3:A:1286:HOH:O	2.50	0.43
1:B:209:ASN:HA	1:B:216:ALA:CB	2.49	0.43
1:B:237:PRO:HD3	3:B:1241:HOH:O	2.16	0.43
1:A:309:LEU:HA	1:A:309:LEU:HD23	1.76	0.43
1:A:32:GLU:OE2	1:A:82:ASP:OD2	2.35	0.43
1:B:297:LYS:NZ	3:B:1280:HOH:O	2.50	0.43
1:A:281:ALA:HB3	1:A:282:PRO:HD3	2.00	0.42
1:B:281:ALA:N	1:B:282:PRO:HD2	2.34	0.42
1:A:270:GLU:HG2	1:A:271:ALA:O	2.19	0.42
1:A:332:GLN:HG2	3:A:1220:HOH:O	2.20	0.42
1:A:210:ASP:OD2	1:A:258:ASP:OD1	2.38	0.41
1:A:207:VAL:HG23	1:A:319:GLU:O	2.20	0.41
1:B:356:LEU:HD23	1:B:356:LEU:HA	1.88	0.41
1:A:81:LEU:HD13	1:A:274:ILE:CG1	2.50	0.41
1:B:234:ALA:O	1:B:235:ASN:HB3	2.20	0.41
1:A:76:ASP:OD2	1:A:268:ASP:OD2	2.39	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:209:ASN:HA	1:B:216:ALA:HB2	2.03	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	353/358 (99%)	341 (97%)	12 (3%)	0	100	100
1	B	353/358 (99%)	345 (98%)	8 (2%)	0	100	100
All	All	706/716 (99%)	686 (97%)	20 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	253/255 (99%)	238 (94%)	15 (6%)	18	10
1	B	253/255 (99%)	236 (93%)	17 (7%)	15	8
All	All	506/510 (99%)	474 (94%)	32 (6%)	16	9

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	GLU
1	A	24	GLN
1	A	67	GLN
1	A	89	LEU
1	A	133	GLU
1	A	175	LEU
1	A	188	LEU
1	A	204	GLN
1	A	239	SER
1	A	245	SER
1	A	297	LYS
1	A	299	VAL
1	A	317	ILE
1	A	327	ARG
1	A	332	GLN
1	B	4	THR
1	B	58	SER
1	B	89	LEU
1	B	133	GLU
1	B	141	ILE
1	B	149	LEU
1	B	172	GLN
1	B	188	LEU
1	B	223	ARG
1	B	245	SER
1	B	246	SER
1	B	272	THR
1	B	274	ILE
1	B	278	SER
1	B	297	LYS
1	B	299	VAL
1	B	326	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	172	GLN
1	A	308	GLN
1	A	332	GLN
1	B	13	GLN
1	B	24	GLN
1	B	55	GLN
1	B	67	GLN

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Mol	Chain	Res	Type
1	B	172	GLN
1	B	256	ASN
1	B	326	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 ⓘ

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	355/358 (99%)	-0.06	3 (0%) 82 84	12, 23, 42, 72	0
1	B	355/358 (99%)	-0.12	4 (1%) 78 80	14, 22, 39, 73	0
All	All	710/716 (99%)	-0.09	7 (0%) 79 81	12, 23, 41, 73	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	233	HIS	3.4
1	A	233	HIS	3.1
1	A	327	ARG	2.2
1	B	243	GLY	2.2
1	B	4	THR	2.0
1	A	244	ALA	2.0
1	B	5	ALA	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](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q < 0.9’ lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	CA	A	501	1/1	0.96	0.05	24,24,24,24	0
2	CA	B	502	1/1	0.98	0.02	22,22,22,22	0

6.5 Other polymers [i](#)

There are no such residues in this entry.