



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 01:52 PM UTC

PDB ID : 1VFL / pdb_00001vfl
Title : Adenosine deaminase
Authors : Kinoshita, T.
Deposited on : 2004-04-16
Resolution : 1.80 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

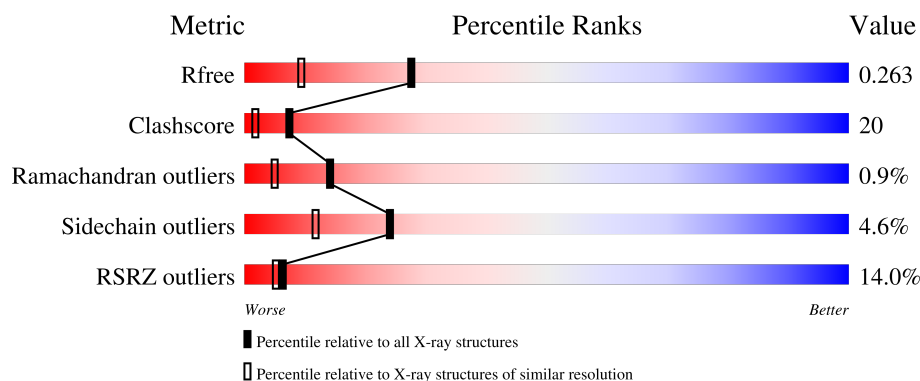
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (1.80-1.80)

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	356	14% 64% 30% ..

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3036 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 Adenosine deaminase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	349	Total	C	N	O	S	0	0	0
			2789	1772	471	534	12			

There are 15 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	5	ASP	ASN	SEE REMARK 999	UNP P56658
A	29	LYS	ARG	SEE REMARK 999	UNP P56658
A	30	ARG	LYS	SEE REMARK 999	UNP P56658
A	54	THR	SER	SEE REMARK 999	UNP P56658
A	57	ASP	GLU	SEE REMARK 999	UNP P56658
A	74	ASP	GLU	SEE REMARK 999	UNP P56658
A	76	ILE	VAL	SEE REMARK 999	UNP P56658
A	196	GLN	LYS	SEE REMARK 999	UNP P56658
A	243	THR	ALA	SEE REMARK 999	UNP P56658
A	258	ILE	VAL	SEE REMARK 999	UNP P56658
A	276	ALA	PRO	SEE REMARK 999	UNP P56658
A	278	ILE	VAL	SEE REMARK 999	UNP P56658
A	310	LYS	ASN	SEE REMARK 999	UNP P56658
A	311	ASP	GLU	SEE REMARK 999	UNP P56658
A	349	ARG	GLY	SEE REMARK 999	UNP P56658

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		

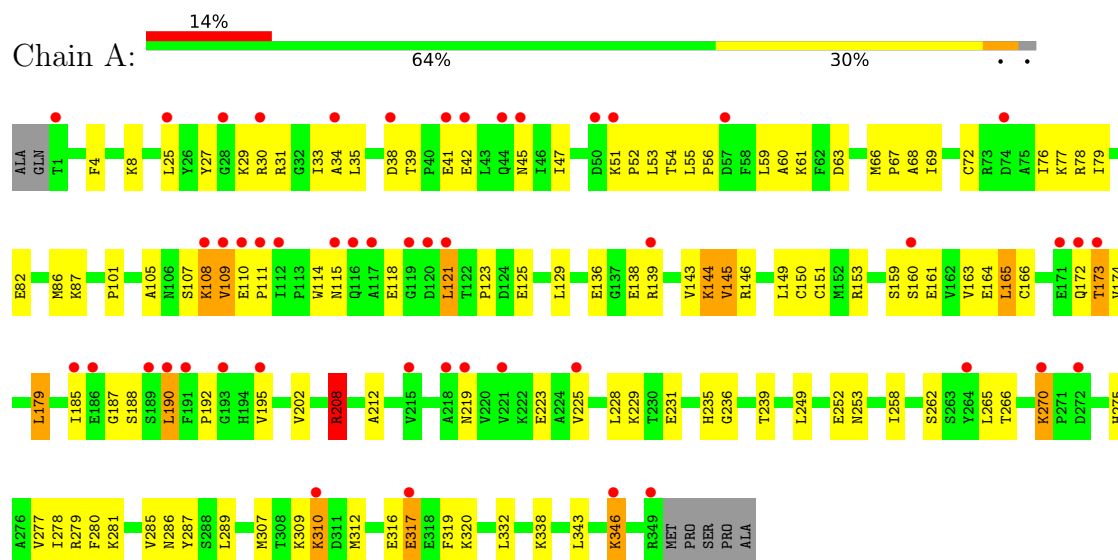
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	246	Total 246	O 246	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: Adenosine deaminase



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	73.73Å 73.73Å 134.01Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	28.43 – 1.80 28.43 – 1.80	Depositor EDS
% Data completeness (in resolution range)	96.7 (28.43-1.80) 96.8 (28.43-1.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.29 (at 1.80Å)	Xtriage
Refinement program	CNX 2002	Depositor
R, R_{free}	0.230 , 0.267 0.228 , 0.263	Depositor DCC
R_{free} test set	1212 reflections (3.55%)	wwPDB-VP
Wilson B-factor (Å ²)	25.9	Xtriage
Anisotropy	0.313	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 52.3	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	0.95	EDS
Total number of atoms	3036	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.89% 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](#)

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

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	0.36	0/2853	0.90	8/3867 (0.2%)

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	151	CYS	N-CA-C	-7.10	99.03	110.32
1	A	159	SER	N-CA-C	6.36	119.03	111.33
1	A	150	CYS	N-CA-C	5.84	118.75	109.24
1	A	145	VAL	N-CA-C	5.55	116.11	108.12
1	A	270	LYS	CA-C-N	5.36	126.53	119.84
1	A	270	LYS	C-N-CA	5.36	126.53	119.84
1	A	208	ARG	N-CA-C	5.20	117.92	109.24
1	A	225	VAL	N-CA-C	5.15	115.82	110.82

There are no chirality outliers.

There are no planarity outliers.

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	2789	0	2746	113	1
2	A	1	0	0	0	0
3	A	246	0	0	20	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	3036	0	2746	113	1

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 20.

All (113) 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:208:ARG:HH12	1:A:231:GLU:HG3	1.00	1.12
1:A:208:ARG:HH11	1:A:208:ARG:HB2	1.14	1.06
1:A:258:ILE:HD11	1:A:280:PHE:HD2	1.27	1.00
1:A:208:ARG:NH1	1:A:231:GLU:HG3	1.80	0.97
1:A:144:LYS:HE3	1:A:332:LEU:HD11	1.55	0.89
1:A:110:GLU:HB3	1:A:111:PRO:HD3	1.53	0.89
1:A:208:ARG:HB2	1:A:208:ARG:NH1	1.91	0.86
1:A:195:VAL:HA	3:A:629:HOH:O	1.79	0.80
1:A:258:ILE:HD11	1:A:280:PHE:CD2	2.14	0.79
1:A:4:PHE:HB3	1:A:8:LYS:NZ	1.97	0.79
1:A:67:PRO:HD3	3:A:640:HOH:O	1.80	0.79
1:A:228:LEU:HG	3:A:629:HOH:O	1.81	0.79
1:A:190:LEU:HD21	1:A:223:GLU:OE2	1.84	0.77
1:A:219:ASN:O	1:A:223:GLU:HG3	1.87	0.74
1:A:66:MET:HB2	3:A:640:HOH:O	1.89	0.72
1:A:258:ILE:HD12	1:A:287:TYR:CD2	2.26	0.71
1:A:310:LYS:HE3	1:A:310:LYS:HA	1.73	0.70
1:A:139:ARG:NH1	1:A:139:ARG:HB3	2.08	0.69
1:A:138:GLU:HG3	1:A:143:VAL:O	1.94	0.67
1:A:208:ARG:NE	3:A:637:HOH:O	2.27	0.67
1:A:343:LEU:O	1:A:346:LYS:HG2	1.95	0.66
1:A:30:ARG:HG2	3:A:626:HOH:O	1.96	0.66
1:A:38:ASP:HB3	3:A:660:HOH:O	1.96	0.66
1:A:190:LEU:HD21	1:A:223:GLU:CD	2.21	0.66
1:A:187:GLY:HA2	1:A:190:LEU:HD23	1.78	0.65
1:A:107:SER:O	1:A:108:LYS:HB3	1.97	0.65
1:A:108:LYS:HG2	1:A:109:VAL:H	1.63	0.63
1:A:51:LYS:O	1:A:53:LEU:HD13	1.99	0.63
1:A:4:PHE:HB3	1:A:8:LYS:HZ1	1.63	0.63
1:A:190:LEU:HD22	1:A:190:LEU:N	2.13	0.63
1:A:144:LYS:HZ1	1:A:146:ARG:HG2	1.64	0.62
1:A:192:PRO:HA	1:A:195:VAL:HG12	1.81	0.62
1:A:139:ARG:HB3	1:A:139:ARG:HH11	1.62	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:87:LYS:HA	1:A:87:LYS:HE2	1.83	0.60
1:A:108:LYS:HD3	1:A:123:PRO:HD2	1.84	0.59
1:A:125:GLU:O	1:A:129:LEU:HD23	2.01	0.59
1:A:316:GLU:O	1:A:320:LYS:HG3	2.03	0.58
1:A:82:GLU:O	1:A:86:MET:HG3	2.02	0.58
1:A:55:LEU:HB3	1:A:56:PRO:HD3	1.86	0.58
1:A:63:ASP:HA	3:A:640:HOH:O	2.03	0.57
1:A:212:ALA:O	1:A:236:GLY:HA2	2.05	0.56
1:A:110:GLU:CB	1:A:111:PRO:HD3	2.31	0.56
1:A:208:ARG:CZ	3:A:637:HOH:O	2.53	0.55
1:A:166:CYS:HA	1:A:174:VAL:HG11	1.88	0.55
1:A:338:LYS:HG3	3:A:597:HOH:O	2.06	0.55
1:A:239:THR:HB	3:A:674:HOH:O	2.08	0.54
1:A:144:LYS:HZ3	1:A:145:VAL:N	2.06	0.53
1:A:60:ALA:HA	1:A:63:ASP:OD2	2.07	0.53
1:A:72:CYS:O	1:A:76:ILE:HG12	2.08	0.53
1:A:208:ARG:HH11	1:A:208:ARG:CB	2.05	0.53
1:A:34:ALA:C	1:A:35:LEU:HD12	2.34	0.53
1:A:146:ARG:HD3	1:A:173:THR:HA	1.92	0.52
1:A:107:SER:O	1:A:108:LYS:CB	2.58	0.52
1:A:164:GLU:HG2	3:A:622:HOH:O	2.11	0.51
1:A:190:LEU:HD22	1:A:190:LEU:H	1.75	0.51
1:A:172:GLN:C	1:A:173:THR:HG23	2.35	0.50
1:A:192:PRO:HA	1:A:195:VAL:CG1	2.42	0.49
1:A:278:ILE:HD11	1:A:312:MET:HG2	1.94	0.49
1:A:41:GLU:HG3	3:A:625:HOH:O	2.13	0.49
1:A:172:GLN:O	1:A:172:GLN:HG2	2.14	0.48
1:A:108:LYS:O	1:A:109:VAL:C	2.57	0.48
1:A:31:ARG:HH21	1:A:68:ALA:C	2.21	0.48
1:A:153:ARG:HG3	1:A:179:LEU:HD22	1.95	0.48
1:A:54:THR:HG23	3:A:733:HOH:O	2.14	0.47
1:A:31:ARG:HG3	1:A:33:ILE:HG12	1.96	0.47
1:A:270:LYS:HG2	3:A:727:HOH:O	2.15	0.47
1:A:67:PRO:HG3	1:A:114:TRP:HB3	1.96	0.47
1:A:27:TYR:HE2	1:A:78:ARG:HD3	1.80	0.46
1:A:139:ARG:NH1	1:A:139:ARG:CB	2.77	0.46
1:A:125:GLU:HG2	1:A:129:LEU:HD23	1.98	0.46
1:A:228:LEU:HD12	1:A:228:LEU:N	2.30	0.46
1:A:4:PHE:HB3	1:A:8:LYS:HZ3	1.74	0.46
1:A:30:ARG:HH11	1:A:30:ARG:HG3	1.80	0.46
1:A:108:LYS:HE2	1:A:161:GLU:CD	2.41	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:144:LYS:HE3	1:A:332:LEU:CD1	2.37	0.46
1:A:66:MET:N	1:A:67:PRO:CD	2.79	0.46
1:A:115:ASN:ND2	3:A:603:HOH:O	2.49	0.46
1:A:47:ILE:O	1:A:61:LYS:HD2	2.16	0.46
1:A:39:THR:HG23	1:A:42:GLU:HB3	1.99	0.45
1:A:105:ALA:HB1	1:A:121:LEU:O	2.16	0.45
1:A:101:PRO:HG3	1:A:149:LEU:HD22	1.98	0.45
1:A:149:LEU:CD1	1:A:165:LEU:HB3	2.47	0.44
1:A:39:THR:CG2	1:A:42:GLU:HB3	2.47	0.44
1:A:202:VAL:HG22	1:A:208:ARG:HE	1.81	0.44
1:A:275:HIS:ND1	1:A:277:VAL:HG12	2.32	0.44
1:A:121:LEU:HD23	1:A:121:LEU:HA	1.76	0.44
1:A:69:ILE:HG22	1:A:76:ILE:HD13	2.00	0.44
1:A:253:ASN:CG	1:A:253:ASN:O	2.60	0.44
1:A:346:LYS:NZ	3:A:641:HOH:O	2.50	0.44
1:A:25:LEU:O	1:A:29:LYS:HG2	2.19	0.43
1:A:27:TYR:CE2	1:A:78:ARG:HD3	2.54	0.43
1:A:4:PHE:HD2	1:A:8:LYS:NZ	2.17	0.43
1:A:52:PRO:HA	1:A:265:LEU:O	2.19	0.43
1:A:108:LYS:HB2	3:A:610:HOH:O	2.19	0.42
1:A:208:ARG:NH1	1:A:229:LYS:O	2.52	0.42
1:A:317:GLU:O	1:A:317:GLU:OE1	2.37	0.42
1:A:160:SER:O	1:A:163:VAL:HG12	2.19	0.42
1:A:139:ARG:CB	1:A:139:ARG:CZ	2.97	0.42
1:A:309:LYS:HE2	1:A:319:PHE:CE1	2.54	0.42
1:A:146:ARG:CD	1:A:173:THR:HA	2.49	0.42
1:A:262:SER:O	1:A:266:THR:HG23	2.19	0.42
1:A:77:LYS:NZ	1:A:136:GLU:OE2	2.53	0.41
1:A:185:ILE:HB	1:A:188:SER:HB3	2.03	0.41
1:A:307:MET:HE2	1:A:312:MET:CE	2.50	0.41
1:A:29:LYS:HG3	3:A:626:HOH:O	2.20	0.41
1:A:59:LEU:HD11	3:A:667:HOH:O	2.20	0.41
1:A:69:ILE:HG21	1:A:79:ILE:HD13	2.03	0.41
1:A:281:LYS:HE2	1:A:281:LYS:HB3	1.96	0.40
1:A:285:VAL:CG1	1:A:286:ASN:N	2.84	0.40
1:A:144:LYS:NZ	1:A:145:VAL:N	2.68	0.40
1:A:192:PRO:O	1:A:195:VAL:HG12	2.21	0.40
1:A:252:GLU:O	1:A:253:ASN:HB3	2.21	0.40
1:A:346:LYS:NZ	1:A:346:LYS:CB	2.84	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:139:ARG:O	1:A:279:ARG:NH2[4_565]	2.19	0.01

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	347/356 (98%)	331 (95%)	13 (4%)	3 (1%)	14 5

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	109	VAL
1	A	108	LYS
1	A	235	HIS

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	304/309 (98%)	290 (95%)	14 (5%)	24 12

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	45	ASN
1	A	118	GLU
1	A	121	LEU

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Mol	Chain	Res	Type
1	A	144	LYS
1	A	165	LEU
1	A	173	THR
1	A	179	LEU
1	A	190	LEU
1	A	208	ARG
1	A	249	LEU
1	A	289	LEU
1	A	310	LYS
1	A	317	GLU
1	A	346	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	45	ASN
1	A	131	ASN
1	A	154	HIS
1	A	172	GLN
1	A	207	HIS
1	A	253	ASN
1	A	306	GLN

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 [i](#)

Of 1 ligands modelled in this entry, 1 is 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 [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	A	349/356 (98%)	0.92	49 (14%) 6 5	18, 31, 49, 56	0

All (49) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	190	LEU	5.4
1	A	120	ASP	4.7
1	A	171	GLU	4.4
1	A	189	SER	4.1
1	A	191	PHE	4.0
1	A	264	TYR	3.9
1	A	272	ASP	3.9
1	A	185	ILE	3.9
1	A	30	ARG	3.8
1	A	108	LYS	3.7
1	A	215	VAL	3.7
1	A	38	ASP	3.4
1	A	115	ASN	3.4
1	A	109	VAL	3.3
1	A	111	PRO	3.3
1	A	41	GLU	3.2
1	A	25	LEU	3.2
1	A	172	GLN	3.2
1	A	186	GLU	3.2
1	A	119	GLY	3.1
1	A	44	GLN	3.1
1	A	173	THR	3.1
1	A	349	ARG	3.1
1	A	112	ILE	3.0
1	A	110	GLU	3.0
1	A	28	GLY	2.9
1	A	42	GLU	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	346	LYS	2.6
1	A	50	ASP	2.6
1	A	195	VAL	2.6
1	A	193	GLY	2.5
1	A	270	LYS	2.5
1	A	117	ALA	2.5
1	A	218	ALA	2.5
1	A	74	ASP	2.5
1	A	139	ARG	2.4
1	A	219	ASN	2.4
1	A	1	THR	2.3
1	A	34	ALA	2.3
1	A	51	LYS	2.2
1	A	317	GLU	2.2
1	A	310	LYS	2.2
1	A	116	GLN	2.2
1	A	57	ASP	2.1
1	A	121	LEU	2.1
1	A	45	ASN	2.1
1	A	221	VAL	2.1
1	A	160	SER	2.0
1	A	225	VAL	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(Å ²)	Q<0.9
2	ZN	A	501	1/1	0.99	0.03	24,24,24,24	0

6.5 Other polymers [i](#)

There are no such residues in this entry.