



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 03:28 AM UTC

PDB ID : 1ADU / pdb_00001adu
Title : EARLY E2A DNA-BINDING PROTEIN
Authors : Tucker, P.A.; Kanellopoulos, P.N.; Tsernoglou, D.; Van Der Vliet, P.C.
Deposited on : 1995-05-11
Resolution : 3.00 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

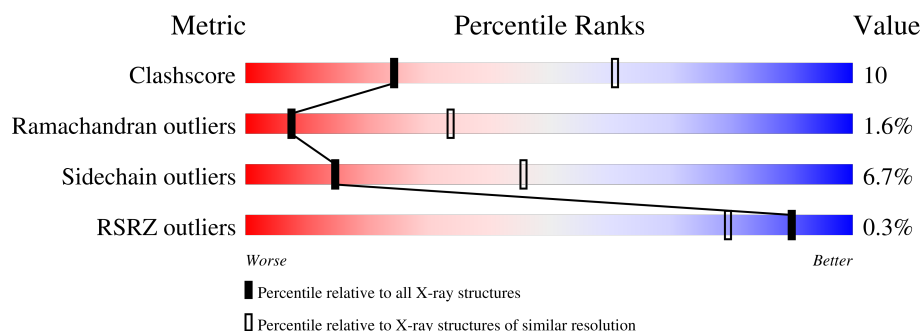
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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	
1	B	356	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4534 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 ADENOVIRUS SINGLE-STRANDED DNA-BINDING PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	287	Total	C	N	O	S	0	0	0
			2268	1444	397	404	23			
1	B	286	Total	C	N	O	S	0	0	0
			2262	1441	393	406	22			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	237	ASN	LYS	conflict	UNP P03265
B	237	ASN	LYS	conflict	UNP P03265

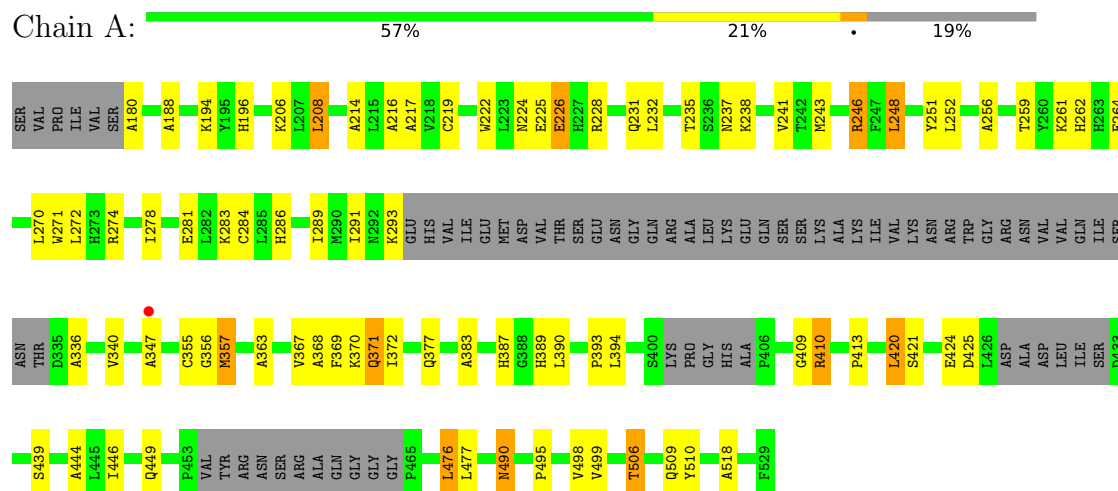
- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

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

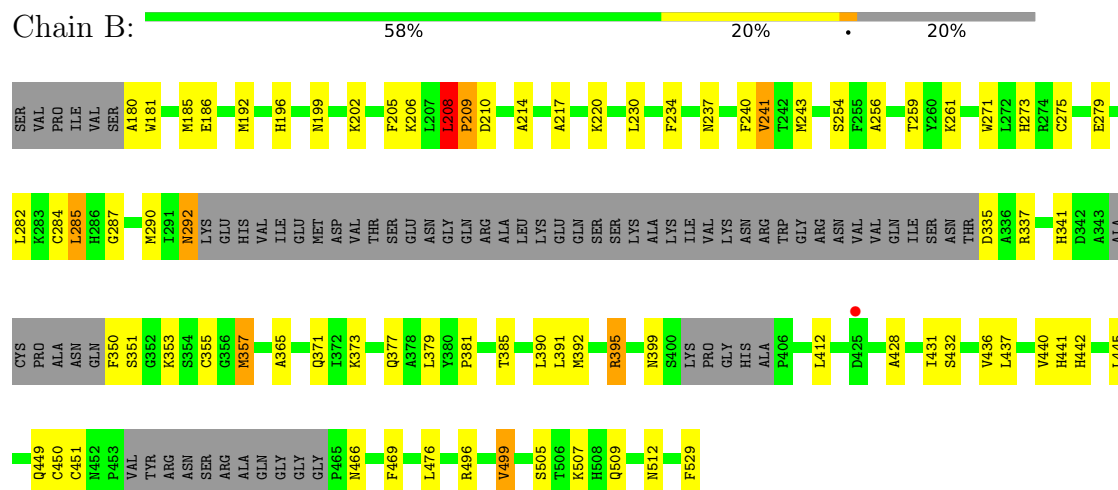
3 Residue-property plots [i](#)

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: ADENOVIRUS SINGLE-STRANDED DNA-BINDING PROTEIN



• Molecule 1: ADENOVIRUS SINGLE-STRANDED DNA-BINDING PROTEIN



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	61.00Å 91.20Å 149.40Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	6.00 – 3.00 6.00 – 3.00	Depositor EDS
% Data completeness (in resolution range)	86.8 (6.00-3.00) 82.3 (6.00-3.00)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.47 (at 2.98Å)	Xtriage
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.201 , 0.296 0.198 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	55.9	Xtriage
Anisotropy	0.467	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 85.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4534	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.52% 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	0.39	0/2325	0.90	6/3139 (0.2%)
1	B	0.39	0/2318	0.92	6/3130 (0.2%)
All	All	0.39	0/4643	0.91	12/6269 (0.2%)

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	499	VAL	CA-C-N	6.48	126.50	119.89
1	B	499	VAL	C-N-CA	6.48	126.50	119.89
1	A	499	VAL	N-CA-C	-6.41	103.11	108.63
1	B	208	LEU	CA-C-N	6.18	127.57	119.84
1	B	208	LEU	C-N-CA	6.18	127.57	119.84
1	A	355	CYS	N-CA-C	-6.10	105.87	113.38
1	A	264	GLU	CA-C-N	5.76	125.23	119.24
1	A	264	GLU	C-N-CA	5.76	125.23	119.24
1	B	357	MET	N-CA-C	5.42	117.56	109.59
1	A	425	ASP	N-CA-C	5.41	117.42	108.49
1	A	498	VAL	N-CA-C	5.26	115.35	107.51
1	B	287	GLY	N-CA-C	-5.08	108.11	115.27

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	2268	0	2214	42	0
1	B	2262	0	2205	48	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
All	All	4534	0	4419	89	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 10.

All (89) 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:391:LEU:HD12	1:B:476:LEU:HG	1.69	0.75
1:B:192:MET:HE3	1:B:214:ALA:HB1	1.72	0.70
1:B:290:MET:HB2	1:B:337:ARG:HH11	1.56	0.70
1:A:188:ALA:HB2	1:A:252:LEU:HD21	1.79	0.64
1:A:383:ALA:HB3	1:A:387:HIS:NE2	2.13	0.64
1:B:220:LYS:HD3	1:B:241:VAL:HG11	1.80	0.62
1:B:395:ARG:HB3	1:B:399:ASN:HB2	1.82	0.61
1:A:421:SER:O	1:A:424:GLU:HG2	2.02	0.60
1:A:271:TRP:HH2	1:A:357:MET:HE1	1.66	0.59
1:B:373:LYS:O	1:B:377:GLN:HG3	2.03	0.59
1:B:290:MET:HB2	1:B:337:ARG:NH1	2.18	0.59
1:A:377:GLN:HA	1:A:387:HIS:CE1	2.37	0.58
1:B:181:TRP:HZ2	1:B:206:LYS:HG3	1.68	0.58
1:A:518:ALA:HB1	1:B:379:LEU:HG	1.86	0.57
1:A:222:TRP:O	1:A:226:GLU:HB2	2.06	0.55
1:B:353:LYS:HA	1:B:509:GLN:NE2	2.21	0.55
1:B:275:CYS:HB3	1:B:282:LEU:HD23	1.88	0.54
1:B:192:MET:CE	1:B:202:LYS:HA	2.38	0.54
1:B:292:ASN:OD1	1:B:335:ASP:N	2.43	0.52
1:A:368:ALA:HB1	1:A:446:ILE:HG13	1.90	0.52
1:B:505:SER:OG	1:B:507:LYS:HG2	2.10	0.52
1:A:232:LEU:HB3	1:A:235:THR:HB	1.92	0.52
1:B:371:GLN:NE2	1:B:445:LEU:HD23	2.27	0.50
1:B:254:SER:HB3	1:B:496:ARG:O	2.11	0.50
1:A:284:CYS:HA	1:A:357:MET:HE2	1.94	0.49
1:B:192:MET:HE2	1:B:202:LYS:HA	1.94	0.49
1:A:272:LEU:HD12	1:A:389:HIS:NE2	2.28	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:371:GLN:HG3	1:A:372:ILE:N	2.25	0.49
1:A:420:LEU:HD23	1:A:439:SER:HB3	1.94	0.49
1:A:224:ASN:HA	1:A:228:ARG:HG3	1.93	0.49
1:A:274:ARG:HB2	1:A:283:LYS:HD2	1.95	0.49
1:A:371:GLN:HG2	1:A:444:ALA:O	2.12	0.49
1:A:393:PRO:HD3	1:A:476:LEU:HD11	1.94	0.49
1:B:450:CYS:SG	1:B:466:ASN:HA	2.53	0.48
1:B:285:LEU:HD22	1:B:357:MET:SD	2.53	0.48
1:B:284:CYS:HB3	1:B:290:MET:SD	2.53	0.48
1:B:180:ALA:O	1:B:256:ALA:HB1	2.13	0.47
1:A:219:CYS:SG	1:A:248:LEU:HD12	2.54	0.47
1:B:234:PHE:HD2	1:B:445:LEU:HD12	1.80	0.47
1:B:282:LEU:H	1:B:337:ARG:NH1	2.13	0.47
1:A:369:PHE:O	1:A:372:ILE:HG22	2.15	0.47
1:B:351:SER:HB3	1:B:512:ASN:OD1	2.15	0.46
1:A:356:GLY:O	1:A:410:ARG:NH1	2.48	0.46
1:B:371:GLN:HB2	1:B:442:HIS:HB2	1.97	0.46
1:A:284:CYS:SG	1:A:286:HIS:HB2	2.56	0.46
1:A:271:TRP:CH2	1:A:357:MET:HE1	2.48	0.46
1:A:278:ILE:HB	1:A:281:GLU:HB3	1.98	0.46
1:B:271:TRP:HB2	1:B:392:MET:SD	2.57	0.45
1:A:180:ALA:N	1:A:256:ALA:O	2.50	0.44
1:A:291:ILE:HD11	1:A:340:VAL:HG23	1.99	0.44
1:A:194:LYS:HD3	1:A:490:ASN:HD21	1.83	0.44
1:B:185:MET:HE1	1:B:205:PHE:O	2.17	0.44
1:A:293:LYS:HB2	1:A:336:ALA:HB3	2.00	0.44
1:A:251:TYR:HD1	1:A:495:PRO:HG2	1.84	0.43
1:B:428:ALA:O	1:B:437:LEU:HD11	2.18	0.43
1:A:363:ALA:O	1:A:367:VAL:HG23	2.18	0.43
1:B:282:LEU:HD11	1:B:365:ALA:HB3	2.00	0.43
1:B:273:HIS:O	1:B:273:HIS:CG	2.71	0.43
1:A:225:GLU:O	1:A:226:GLU:HG2	2.17	0.43
1:A:214:ALA:O	1:A:217:ALA:HB3	2.19	0.42
1:A:420:LEU:CD2	1:A:439:SER:HB3	2.50	0.42
1:A:243:MET:HA	1:A:246:ARG:HD3	2.01	0.42
1:A:413:PRO:HB3	1:A:449:GLN:O	2.20	0.42
1:B:271:TRP:CE3	1:B:390:LEU:HD23	2.55	0.42
1:B:436:VAL:O	1:B:440:VAL:HG23	2.19	0.42
1:B:243:MET:HE2	1:B:476:LEU:HD11	2.02	0.42
1:B:437:LEU:HD12	1:B:437:LEU:HA	1.83	0.42
1:A:261:LYS:O	1:A:262:HIS:HB2	2.20	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:220:LYS:NZ	1:B:237:ASN:HD21	2.18	0.41
1:A:409:GLY:HA2	1:A:510:TYR:CZ	2.55	0.41
1:A:506:THR:O	1:A:509:GLN:HG2	2.19	0.41
1:A:216:ALA:HB1	1:A:241:VAL:HG12	2.03	0.41
1:B:206:LYS:HB3	1:B:208:LEU:HD13	2.03	0.41
1:B:449:GLN:HA	1:B:469:PHE:O	2.21	0.41
1:A:206:LYS:HB2	1:A:208:LEU:HD22	2.03	0.41
1:B:192:MET:HE1	1:B:202:LYS:HA	2.02	0.41
1:B:341:HIS:CD2	1:B:341:HIS:N	2.89	0.41
1:B:284:CYS:HB2	1:B:355:CYS:SG	2.60	0.41
1:B:181:TRP:CZ2	1:B:206:LYS:HG3	2.52	0.40
1:A:237:ASN:O	1:A:241:VAL:HG23	2.21	0.40
1:B:199:ASN:HA	1:B:202:LYS:HB3	2.03	0.40
1:B:230:LEU:HD21	1:B:240:PHE:CE2	2.56	0.40
1:B:282:LEU:C	1:B:290:MET:HG2	2.47	0.40
1:B:428:ALA:HB1	1:B:441:HIS:CE1	2.56	0.40
1:B:214:ALA:O	1:B:217:ALA:HB3	2.21	0.40
1:A:383:ALA:HB3	1:A:387:HIS:HE2	1.86	0.40
1:A:291:ILE:HG12	1:A:340:VAL:HA	2.03	0.40
1:B:185:MET:HE1	1:B:206:LYS:HA	2.04	0.40
1:B:412:LEU:O	1:B:451:CYS:HA	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	277/356 (78%)	251 (91%)	23 (8%)	3 (1%)	11	43
1	B	276/356 (78%)	248 (90%)	22 (8%)	6 (2%)	5	26
All	All	553/712 (78%)	499 (90%)	45 (8%)	9 (2%)	7	34

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	226	GLU
1	A	347	ALA
1	B	261	LYS
1	B	196	HIS
1	A	196	HIS
1	B	209	PRO
1	B	432	SER
1	B	431	ILE
1	B	381	PRO

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	A	245/303 (81%)	226 (92%)	19 (8%)	11	39
1	B	245/303 (81%)	231 (94%)	14 (6%)	18	52
All	All	490/606 (81%)	457 (93%)	33 (7%)	15	46

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	208	LEU
1	A	231	GLN
1	A	238	LYS
1	A	246	ARG
1	A	248	LEU
1	A	259	THR
1	A	270	LEU
1	A	289	ILE
1	A	357	MET
1	A	370	LYS
1	A	371	GLN
1	A	390	LEU
1	A	394	LEU
1	A	410	ARG

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	420	LEU
1	A	476	LEU
1	A	477	LEU
1	A	490	ASN
1	A	506	THR
1	B	186	GLU
1	B	208	LEU
1	B	209	PRO
1	B	210	ASP
1	B	241	VAL
1	B	259	THR
1	B	279	GLU
1	B	285	LEU
1	B	292	ASN
1	B	350	PHE
1	B	385	THR
1	B	395	ARG
1	B	499	VAL
1	B	529	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	227	HIS
1	A	231	GLN
1	A	253	GLN
1	A	349	GLN
1	A	490	ASN
1	A	508	HIS
1	A	512	ASN
1	B	237	ASN
1	B	253	GLN
1	B	262	HIS
1	B	292	ASN
1	B	371	GLN
1	B	377	GLN
1	B	449	GLN
1	B	490	ASN

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 4 ligands modelled in this entry, 4 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 [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	A	287/356 (80%)	-0.78	1 (0%) 90 79	7, 28, 61, 73	0
1	B	286/356 (80%)	-0.66	1 (0%) 90 79	11, 39, 66, 78	0
All	All	573/712 (80%)	-0.72	2 (0%) 90 79	7, 34, 64, 78	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	347	ALA	2.4
1	B	425	ASP	2.3

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	B	531	1/1	0.96	0.06	52,52,52,52	0
2	ZN	A	531	1/1	0.99	0.01	41,41,41,41	0
2	ZN	B	530	1/1	0.99	0.01	44,44,44,44	0

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	ZN	A	530	1/1	0.99	0.03	18,18,18,18	0

6.5 Other polymers [i](#)

There are no such residues in this entry.