



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 10:21 AM UTC

PDB ID : 2NYT / pdb_00002nyt
Title : The APOBEC2 Crystal Structure and Functional Implications for AID
Authors : Prochnow, C.; Bransteitter, R.; Klein, M.; Goodman, M.; Chen, X.
Deposited on : 2006-11-21
Resolution : 2.50 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

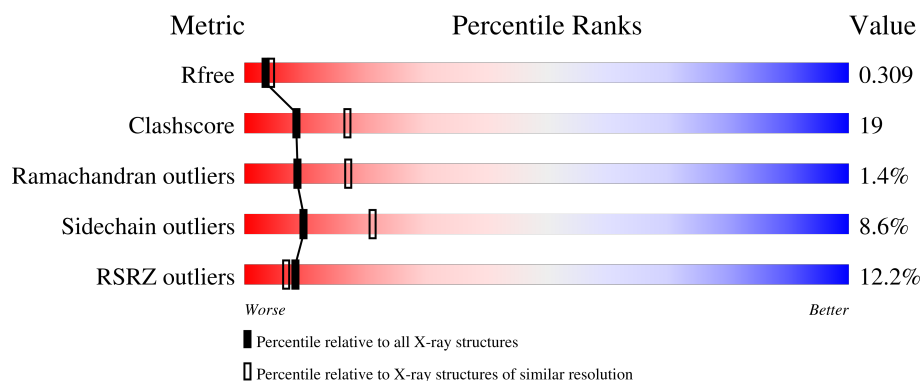
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	190	11% 54% 38% ..
1	B	190	16% 67% 30% ..
1	C	190	12% 54% 40% ...
1	D	190	9% 58% 31% 5% • 6%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6088 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 Probable C->U-editing enzyme APOBEC-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	185	Total	C	N	O	S	0	0	0
			1500	968	251	274	7			
1	B	190	Total	C	N	O	S	0	0	0
			1523	981	256	279	7			
1	C	185	Total	C	N	O	S	0	0	0
			1501	970	251	273	7			
1	D	179	Total	C	N	O	S	0	0	0
			1472	953	245	267	7			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	35	GLY	-	cloning artifact	UNP Q9Y235
A	36	SER	-	cloning artifact	UNP Q9Y235
A	37	GLY	-	cloning artifact	UNP Q9Y235
A	38	GLY	-	cloning artifact	UNP Q9Y235
A	39	GLY	-	cloning artifact	UNP Q9Y235
A	40	MET	-	cloning artifact	UNP Q9Y235
B	35	GLY	-	cloning artifact	UNP Q9Y235
B	36	SER	-	cloning artifact	UNP Q9Y235
B	37	GLY	-	cloning artifact	UNP Q9Y235
B	38	GLY	-	cloning artifact	UNP Q9Y235
B	39	GLY	-	cloning artifact	UNP Q9Y235
B	40	MET	-	cloning artifact	UNP Q9Y235
C	35	GLY	-	cloning artifact	UNP Q9Y235
C	36	SER	-	cloning artifact	UNP Q9Y235
C	37	GLY	-	cloning artifact	UNP Q9Y235
C	38	GLY	-	cloning artifact	UNP Q9Y235
C	39	GLY	-	cloning artifact	UNP Q9Y235
C	40	MET	-	cloning artifact	UNP Q9Y235
D	35	GLY	-	cloning artifact	UNP Q9Y235
D	36	SER	-	cloning artifact	UNP Q9Y235
D	37	GLY	-	cloning artifact	UNP Q9Y235

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Chain	Residue	Modelled	Actual	Comment	Reference
D	38	GLY	-	cloning artifact	UNP Q9Y235
D	39	GLY	-	cloning artifact	UNP Q9Y235
D	40	MET	-	cloning artifact	UNP Q9Y235

- 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		
2	B	1	Total	Zn	0	0
			1	1		
2	C	1	Total	Zn	0	0
			1	1		
2	D	1	Total	Zn	0	0
			1	1		

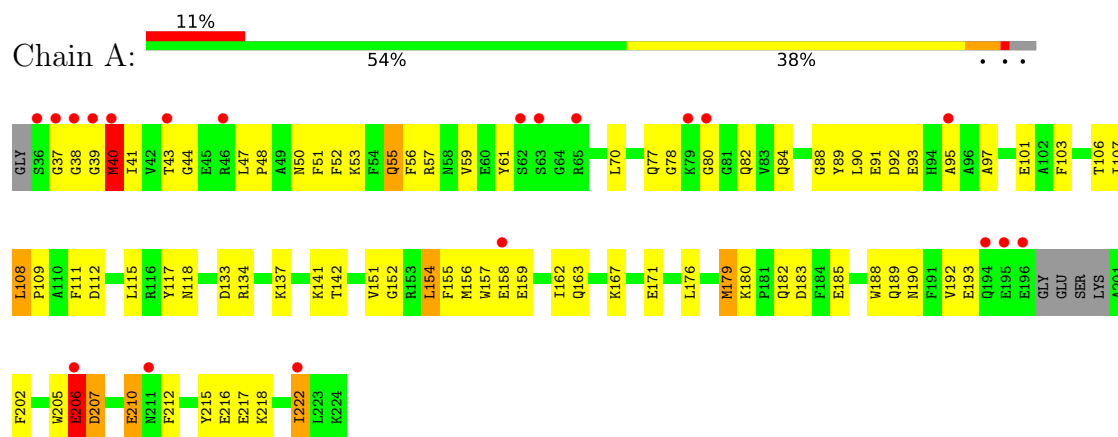
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	32	Total	O	0	0
			32	32		
3	B	23	Total	O	0	0
			23	23		
3	C	19	Total	O	0	0
			19	19		
3	D	14	Total	O	0	0
			14	14		

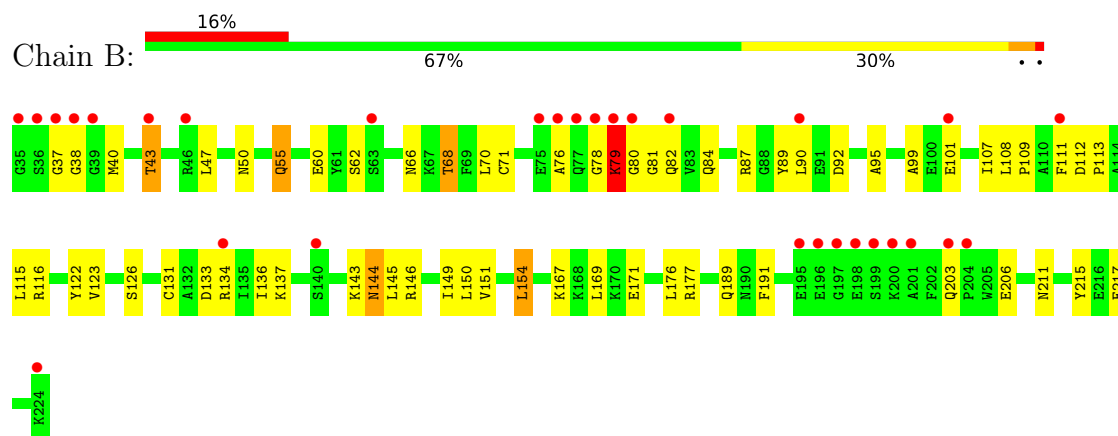
3 Residue-property plots [i](#)

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.

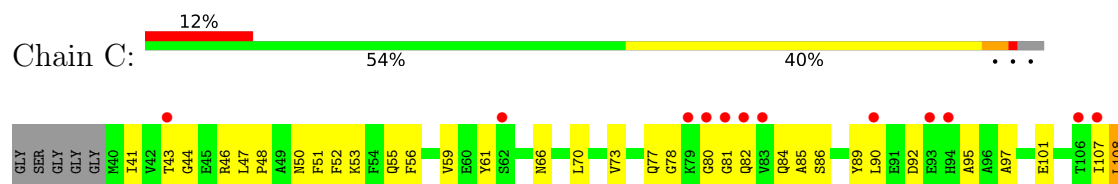
- Molecule 1: Probable C->U-editing enzyme APOBEC-2



- Molecule 1: Probable C->U-editing enzyme APOBEC-2

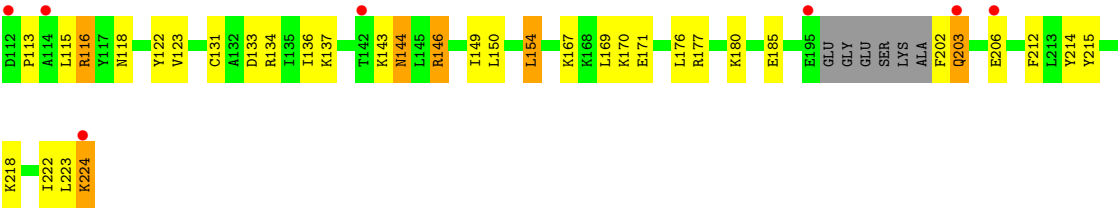
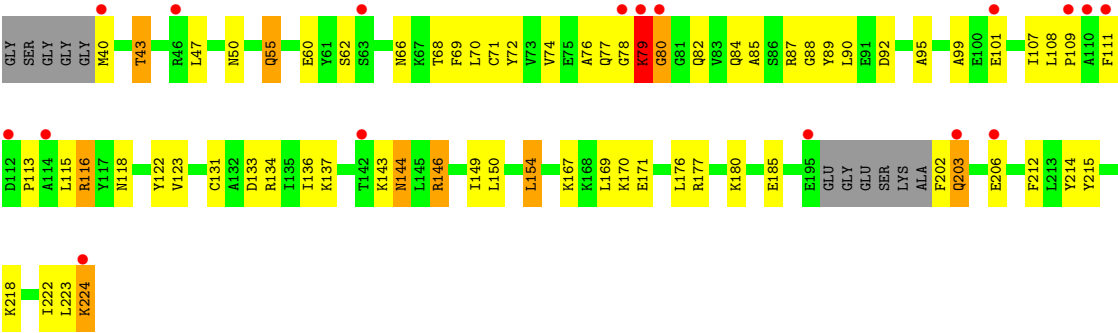


- Molecule 1: Probable C->U-editing enzyme APOBEC-2





● Molecule 1: Probable C->U-editing enzyme APOBEC-2



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	37.84Å 89.41Å 245.77Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.50 50.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	(Not available) (50.00-2.50) 92.5 (50.00-2.50)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.80 (at 2.24Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.246 , 0.295 0.263 , 0.309	Depositor DCC
R_{free} test set	2095 reflections (5.83%)	wwPDB-VP
Wilson B-factor (Å ²)	23.6	Xtriage
Anisotropy	0.772	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 43.3	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	0.89	EDS
Total number of atoms	6088	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.39% 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	0.61	2/1536 (0.1%)	0.81	3/2071 (0.1%)
1	B	0.54	0/1560	0.81	1/2105 (0.0%)
1	C	0.54	0/1538	0.80	2/2077 (0.1%)
1	D	0.49	0/1508	0.78	0/2034
All	All	0.55	2/6142 (0.0%)	0.80	6/8287 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	179	MET	SD-CE	-7.04	1.61	1.79
1	A	40	MET	SD-CE	6.32	1.95	1.79

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	108	LEU	CA-C-N	5.94	125.15	118.97
1	A	108	LEU	C-N-CA	5.94	125.15	118.97
1	C	108	LEU	CA-C-N	5.87	125.07	118.97
1	C	108	LEU	C-N-CA	5.87	125.07	118.97
1	A	216	GLU	N-CA-C	-5.51	105.35	111.36
1	B	206	GLU	N-CA-C	5.46	117.67	111.11

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	1500	0	1464	75	0
1	B	1523	0	1477	55	0
1	C	1501	0	1460	65	0
1	D	1472	0	1443	53	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	32	0	0	0	0
3	B	23	0	0	2	0
3	C	19	0	0	0	0
3	D	14	0	0	2	0
All	All	6088	0	5844	224	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 19.

All (224) 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:D:131:CYS:HA	1:D:134:ARG:HD2	1.49	0.91
1:C:46:ARG:HE	1:C:195:GLU:HG3	1.37	0.89
1:B:131:CYS:HA	1:B:134:ARG:HD2	1.53	0.89
1:C:47:LEU:HD12	1:C:48:PRO:HD2	1.66	0.78
1:D:78:GLY:HA3	1:D:115:LEU:HD23	1.66	0.77
1:B:78:GLY:HA3	1:B:115:LEU:HD23	1.65	0.76
1:B:101:GLU:HG2	1:B:134:ARG:HD3	1.68	0.76
1:A:47:LEU:HD12	1:A:48:PRO:HD2	1.66	0.75
1:D:60:GLU:HA	1:D:66:ASN:HB3	1.68	0.75
1:D:101:GLU:HG2	1:D:134:ARG:HD3	1.70	0.74
1:A:39:GLY:HA2	1:A:193:GLU:HA	1.73	0.71
1:B:60:GLU:HA	1:B:66:ASN:HB3	1.71	0.71
1:A:61:TYR:OH	1:C:159:GLU:HB3	1.91	0.70
1:D:40:MET:HE1	1:D:87:ARG:HG3	1.74	0.69
1:A:40:MET:H	1:A:193:GLU:HB2	1.58	0.69
1:B:144:ASN:HD22	1:B:144:ASN:H	1.40	0.69
1:C:41:ILE:HG12	1:C:89:TYR:CD2	2.29	0.67
1:A:93:GLU:HG3	1:B:81:GLY:HA3	1.75	0.67
1:D:144:ASN:H	1:D:144:ASN:HD22	1.43	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:41:ILE:HG12	1:A:89:TYR:CD2	2.31	0.66
1:C:84:GLN:HG2	1:D:90:LEU:HD22	1.78	0.66
1:B:154:LEU:HB2	1:B:215:TYR:CD2	2.31	0.65
1:A:218:LYS:O	1:A:222:ILE:HG22	1.98	0.63
1:B:47:LEU:HD11	1:B:203:GLN:HG2	1.81	0.63
1:B:111:PHE:O	1:B:113:PRO:HD3	1.99	0.62
1:B:144:ASN:HD22	1:B:144:ASN:N	1.95	0.62
1:A:38:GLY:HA3	1:B:87:ARG:HH22	1.65	0.62
1:B:107:ILE:C	1:B:109:PRO:HD3	2.24	0.62
1:A:92:ASP:CG	1:B:82:GLN:HB3	2.25	0.61
1:D:170:LYS:HD2	1:D:223:LEU:HD23	1.82	0.61
1:C:152:GLY:N	1:C:179:MET:HE3	2.15	0.61
1:C:218:LYS:O	1:C:222:ILE:HG22	2.01	0.61
1:D:144:ASN:HD22	1:D:144:ASN:N	1.96	0.61
1:D:107:ILE:C	1:D:109:PRO:HD3	2.26	0.61
1:B:167:LYS:O	1:B:171:GLU:HG2	2.00	0.60
1:B:71:CYS:HB2	1:B:122:TYR:HB2	1.82	0.60
1:C:81:GLY:O	1:D:92:ASP:HA	2.02	0.59
1:D:111:PHE:O	1:D:113:PRO:HD3	2.02	0.59
1:B:40:MET:HE3	1:B:191:PHE:HA	1.83	0.59
1:D:167:LYS:O	1:D:171:GLU:HG2	2.02	0.59
1:B:37:GLY:HA3	1:B:189:GLN:O	2.02	0.59
1:C:51:PHE:HA	1:C:55:GLN:NE2	2.17	0.59
1:D:60:GLU:HA	1:D:66:ASN:CB	2.33	0.58
1:C:210:GLU:H	1:C:210:GLU:CD	2.10	0.58
1:A:51:PHE:HA	1:A:55:GLN:NE2	2.18	0.58
1:B:90:LEU:HD21	1:B:107:ILE:HD11	1.86	0.58
1:A:154:LEU:HB3	1:A:157:TRP:HB3	1.85	0.57
1:C:51:PHE:HA	1:C:55:GLN:HE21	1.69	0.57
1:C:167:LYS:O	1:C:171:GLU:HG3	2.05	0.57
1:D:90:LEU:HD21	1:D:107:ILE:HD11	1.87	0.57
1:A:167:LYS:O	1:A:171:GLU:HG3	2.05	0.57
1:A:51:PHE:HA	1:A:55:GLN:HE21	1.68	0.57
1:A:154:LEU:HB2	1:A:215:TYR:CD2	2.38	0.57
1:C:154:LEU:HB3	1:C:157:TRP:HB3	1.87	0.57
1:C:46:ARG:HE	1:C:195:GLU:CG	2.16	0.56
1:D:92:ASP:HB3	1:D:95:ALA:HB2	1.87	0.56
1:B:60:GLU:HA	1:B:66:ASN:CB	2.36	0.56
1:A:40:MET:SD	1:A:88:GLY:HA2	2.46	0.56
1:A:157:TRP:CZ3	1:A:218:LYS:HD2	2.41	0.56
1:A:37:GLY:HA2	1:A:190:ASN:OD1	2.05	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:71:CYS:HB2	1:D:122:TYR:HB2	1.87	0.55
1:B:144:ASN:H	1:B:144:ASN:ND2	2.04	0.55
1:B:47:LEU:HB3	1:B:50:ASN:HD22	1.71	0.55
1:A:152:GLY:N	1:A:179:MET:HE3	2.21	0.55
1:D:70:LEU:HB2	1:D:99:ALA:HB1	1.89	0.55
1:A:155:PHE:CE2	1:A:156:MET:HE2	2.42	0.55
1:D:47:LEU:HB3	1:D:50:ASN:HD22	1.71	0.55
1:B:154:LEU:HB2	1:B:215:TYR:HD2	1.72	0.54
1:C:137:LYS:O	1:C:141:LYS:HG2	2.07	0.54
1:A:90:LEU:HD13	1:B:84:GLN:OE1	2.08	0.54
1:A:91:GLU:O	1:B:82:GLN:HA	2.07	0.54
1:B:92:ASP:HB3	1:B:95:ALA:HB2	1.89	0.54
1:D:144:ASN:H	1:D:144:ASN:ND2	2.05	0.54
1:C:155:PHE:CE2	1:C:156:MET:HE2	2.43	0.54
1:A:137:LYS:O	1:A:141:LYS:HG2	2.09	0.53
1:C:188:TRP:O	1:C:192:VAL:HB	2.09	0.53
1:C:70:LEU:C	1:C:70:LEU:HD23	2.34	0.53
1:A:92:ASP:HA	1:B:82:GLN:HA	1.89	0.53
1:A:207:ASP:HB2	1:A:210:GLU:HG2	1.90	0.53
1:B:145:LEU:HD12	3:B:1032:HOH:O	2.08	0.53
1:D:78:GLY:O	1:D:79:LYS:C	2.51	0.53
1:C:46:ARG:NE	1:C:195:GLU:HG3	2.17	0.52
1:C:154:LEU:HB2	1:C:215:TYR:CD2	2.44	0.52
1:D:90:LEU:HD21	1:D:107:ILE:CD1	2.39	0.52
1:B:90:LEU:HD21	1:B:107:ILE:CD1	2.39	0.52
1:B:203:GLN:HG3	1:B:203:GLN:O	2.09	0.52
1:A:57:ARG:NH2	1:C:160:PRO:HD2	2.24	0.51
1:B:78:GLY:O	1:B:79:LYS:C	2.54	0.51
1:A:185:GLU:O	1:A:189:GLN:HG3	2.11	0.51
1:A:207:ASP:CB	1:A:210:GLU:HG2	2.40	0.51
1:B:70:LEU:HB2	1:B:99:ALA:HB1	1.91	0.51
1:A:92:ASP:HB3	1:A:95:ALA:HB2	1.92	0.50
1:D:108:LEU:N	1:D:109:PRO:HD3	2.26	0.50
1:B:68:THR:HG21	3:B:1028:HOH:O	2.11	0.50
1:C:47:LEU:HD23	1:C:50:ASN:ND2	2.27	0.50
1:C:82:GLN:HG2	1:D:92:ASP:OD2	2.11	0.50
1:A:70:LEU:C	1:A:70:LEU:HD23	2.37	0.50
1:A:47:LEU:HD23	1:A:50:ASN:ND2	2.27	0.49
1:A:163:GLN:HG2	1:A:222:ILE:HD12	1.92	0.49
1:B:169:LEU:HD23	1:B:176:LEU:HD21	1.93	0.49
1:C:219:LEU:O	1:C:222:ILE:HG23	2.11	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:86:SER:HB3	1:D:72:TYR:OH	2.13	0.49
1:C:41:ILE:HG12	1:C:89:TYR:CE2	2.48	0.49
1:C:86:SER:HA	1:D:88:GLY:HA3	1.94	0.49
1:B:108:LEU:N	1:B:109:PRO:HD3	2.28	0.48
1:C:92:ASP:HB3	1:C:95:ALA:HB2	1.94	0.48
1:A:39:GLY:O	1:A:40:MET:HB2	2.14	0.48
1:A:159:GLU:OE2	1:C:153:ARG:NH2	2.45	0.48
1:C:46:ARG:HD2	1:C:195:GLU:CD	2.39	0.48
1:C:90:LEU:HD22	1:C:107:ILE:HD11	1.96	0.48
1:A:188:TRP:O	1:A:192:VAL:HB	2.14	0.48
1:C:107:ILE:C	1:C:109:PRO:HD3	2.39	0.47
1:D:47:LEU:HD11	1:D:203:GLN:NE2	2.29	0.47
1:D:180:LYS:HG2	1:D:212:PHE:CE1	2.49	0.47
1:C:219:LEU:HG	1:C:223:LEU:HD22	1.96	0.47
1:D:136:ILE:HG13	1:D:169:LEU:HD12	1.96	0.47
1:A:111:PHE:CD1	1:A:142:THR:HG21	2.50	0.47
1:C:85:ALA:HB3	1:D:89:TYR:HE2	1.78	0.47
1:A:43:THR:HG22	1:A:44:GLY:N	2.29	0.47
1:C:163:GLN:O	1:C:167:LYS:HG3	2.14	0.47
1:B:136:ILE:HG13	1:B:169:LEU:HD12	1.96	0.47
1:C:86:SER:HA	1:D:88:GLY:CA	2.45	0.47
1:C:185:GLU:O	1:C:189:GLN:HG3	2.15	0.47
1:D:43:THR:HG21	1:D:55:GLN:HG2	1.96	0.47
1:C:47:LEU:HD12	1:C:48:PRO:CD	2.41	0.47
1:D:214:TYR:CZ	1:D:218:LYS:HE3	2.50	0.47
1:D:47:LEU:HB3	1:D:50:ASN:ND2	2.30	0.46
1:A:182:GLN:H	1:A:182:GLN:NE2	2.13	0.46
1:C:111:PHE:HB3	1:C:117:TYR:CE1	2.50	0.46
1:C:108:LEU:N	1:C:109:PRO:HD3	2.31	0.46
1:B:43:THR:HG21	1:B:55:GLN:HG2	1.97	0.46
1:A:163:GLN:O	1:A:167:LYS:HG3	2.16	0.46
1:D:77:GLN:HB3	3:D:1065:HOH:O	2.16	0.46
1:A:37:GLY:HA3	1:A:189:GLN:O	2.16	0.46
1:D:150:LEU:HD23	1:D:177:ARG:HB2	1.97	0.46
1:C:219:LEU:O	1:C:223:LEU:HB2	2.16	0.46
1:A:111:PHE:HB3	1:A:117:TYR:CE1	2.51	0.45
1:B:55:GLN:H	1:B:55:GLN:NE2	2.14	0.45
1:A:180:LYS:HG2	1:A:183:ASP:OD2	2.16	0.45
1:A:205:TRP:O	1:A:206:GLU:C	2.59	0.45
1:B:150:LEU:HD23	1:B:177:ARG:HB2	1.96	0.45
1:D:76:ALA:HB3	1:D:84:GLN:HB3	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:41:ILE:HG12	1:A:89:TYR:CE2	2.51	0.45
1:A:107:ILE:C	1:A:109:PRO:HD3	2.41	0.45
1:B:47:LEU:HB3	1:B:50:ASN:ND2	2.32	0.45
1:B:76:ALA:HB3	1:B:84:GLN:HB3	1.98	0.45
1:C:111:PHE:N	1:C:111:PHE:CD1	2.85	0.45
1:A:78:GLY:HA3	1:A:115:LEU:HD23	1.99	0.44
1:B:101:GLU:HG2	1:B:134:ARG:CD	2.45	0.44
1:A:84:GLN:HG2	1:B:90:LEU:HD22	2.00	0.44
1:D:80:GLY:HA3	3:D:1063:HOH:O	2.16	0.44
1:A:47:LEU:HD12	1:A:48:PRO:CD	2.44	0.44
1:A:90:LEU:HD22	1:A:107:ILE:HD11	1.99	0.44
1:A:108:LEU:N	1:A:109:PRO:HD3	2.32	0.44
1:C:163:GLN:HG2	1:C:222:ILE:HD12	1.98	0.44
1:B:89:TYR:O	1:B:90:LEU:HD23	2.18	0.43
1:D:149:ILE:HB	1:D:176:LEU:HD23	1.99	0.43
1:A:103:PHE:HA	1:A:107:ILE:HD12	2.00	0.43
1:C:43:THR:HG22	1:C:44:GLY:N	2.33	0.43
1:A:37:GLY:N	1:B:38:GLY:HA2	2.33	0.43
1:A:92:ASP:CB	1:A:95:ALA:HB2	2.47	0.43
1:A:93:GLU:CG	1:B:81:GLY:HA3	2.43	0.43
1:C:111:PHE:CD1	1:C:142:THR:HG21	2.54	0.43
1:D:82:GLN:H	1:D:82:GLN:HG2	1.66	0.43
1:D:167:LYS:HE2	1:D:222:ILE:O	2.18	0.43
1:A:38:GLY:HA2	1:B:37:GLY:HA2	1.98	0.43
1:A:111:PHE:CD1	1:A:111:PHE:N	2.86	0.43
1:D:55:GLN:H	1:D:55:GLN:NE2	2.16	0.43
1:B:149:ILE:HB	1:B:176:LEU:HD23	1.99	0.43
1:A:106:THR:HG21	1:B:112:ASP:HB2	2.00	0.43
1:D:185:GLU:HA	1:D:202:PHE:CE1	2.54	0.43
1:B:144:ASN:N	1:B:144:ASN:ND2	2.63	0.43
1:C:92:ASP:CB	1:C:95:ALA:HB2	2.49	0.43
1:D:154:LEU:HB2	1:D:215:TYR:HD2	1.84	0.43
1:A:52:PHE:HA	1:A:56:PHE:HB3	2.00	0.42
1:C:170:LYS:NZ	1:C:224:LYS:HG2	2.33	0.42
1:A:82:GLN:HG2	1:B:92:ASP:OD2	2.19	0.42
1:B:133:ASP:O	1:B:137:LYS:HG3	2.18	0.42
1:C:180:LYS:HG2	1:C:183:ASP:OD2	2.18	0.42
1:A:55:GLN:NE2	1:A:55:GLN:H	2.18	0.42
1:D:74:VAL:O	1:D:85:ALA:HA	2.20	0.42
1:A:40:MET:HB2	1:A:40:MET:HE3	1.98	0.42
1:C:166:LEU:HB2	1:C:222:ILE:HD11	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:154:LEU:HD12	1:B:154:LEU:HA	1.73	0.42
1:D:133:ASP:O	1:D:137:LYS:HG3	2.19	0.42
1:C:52:PHE:HA	1:C:56:PHE:HB3	2.02	0.42
1:D:169:LEU:HD23	1:D:176:LEU:HD21	2.00	0.42
1:A:97:ALA:HB1	1:A:101:GLU:HB2	2.01	0.42
1:B:70:LEU:HD23	1:B:70:LEU:C	2.44	0.42
1:A:59:VAL:HG12	1:A:61:TYR:H	1.85	0.42
1:A:179:MET:HG2	1:A:183:ASP:HB3	2.02	0.41
1:A:52:PHE:HA	1:A:56:PHE:CB	2.49	0.41
1:C:55:GLN:NE2	1:C:55:GLN:H	2.17	0.41
1:C:97:ALA:HB1	1:C:101:GLU:HB2	2.02	0.41
1:C:182:GLN:H	1:C:182:GLN:NE2	2.18	0.41
1:A:156:MET:HG2	1:C:156:MET:HG2	2.02	0.41
1:A:157:TRP:CH2	1:A:218:LYS:HD2	2.56	0.41
1:D:154:LEU:HD12	1:D:154:LEU:HA	1.75	0.41
1:A:163:GLN:HA	1:A:222:ILE:HD11	2.02	0.41
1:C:78:GLY:HA3	1:C:115:LEU:HD23	2.01	0.41
1:D:170:LYS:NZ	1:D:224:LYS:HD2	2.36	0.41
1:A:101:GLU:HG2	1:A:134:ARG:NE	2.36	0.41
1:C:59:VAL:HB	1:C:66:ASN:HB2	2.02	0.41
1:C:59:VAL:HG12	1:C:61:TYR:H	1.85	0.41
1:C:183:ASP:O	1:C:187:VAL:HG23	2.21	0.41
1:A:112:ASP:OD1	1:A:115:LEU:HD12	2.20	0.41
1:C:73:VAL:HG23	1:C:191:PHE:HZ	1.86	0.41
1:D:69:PHE:HA	1:D:90:LEU:O	2.20	0.41
1:D:90:LEU:HD11	1:D:107:ILE:HG13	2.03	0.41
1:A:156:MET:HB2	1:A:162:ILE:HG13	2.02	0.41
1:A:37:GLY:H	1:B:38:GLY:HA2	1.85	0.41
1:C:112:ASP:OD1	1:C:115:LEU:HD12	2.21	0.41
1:C:156:MET:HE3	1:C:162:ILE:HD11	2.03	0.41
1:C:194:GLN:O	1:C:195:GLU:C	2.63	0.41
1:C:169:LEU:HD23	1:C:176:LEU:HD12	2.03	0.41
1:B:126:SER:HA	1:B:151:VAL:HG11	2.03	0.40
1:C:151:VAL:C	1:C:179:MET:HE3	2.45	0.40
1:D:116:ARG:HH21	1:D:146:ARG:NH1	2.19	0.40
1:D:214:TYR:O	1:D:218:LYS:HG2	2.21	0.40
1:A:163:GLN:HA	1:A:222:ILE:CD1	2.51	0.40
1:A:185:GLU:HB2	1:A:202:PHE:CD2	2.56	0.40
1:A:151:VAL:C	1:A:179:MET:HE3	2.46	0.40
1:A:180:LYS:HB3	1:A:212:PHE:CE1	2.56	0.40
1:C:207:ASP:HA	1:C:210:GLU:HG2	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	181/190 (95%)	166 (92%)	12 (7%)	3 (2%)	7	13
1	B	188/190 (99%)	175 (93%)	11 (6%)	2 (1%)	11	22
1	C	183/190 (96%)	166 (91%)	14 (8%)	3 (2%)	7	14
1	D	175/190 (92%)	163 (93%)	10 (6%)	2 (1%)	11	22
All	All	727/760 (96%)	670 (92%)	47 (6%)	10 (1%)	9	17

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	206	GLU
1	C	223	LEU
1	A	40	MET
1	B	79	LYS
1	C	206	GLU
1	D	79	LYS
1	A	80	GLY
1	B	80	GLY
1	C	80	GLY
1	D	80	GLY

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	155/159 (98%)	142 (92%)	13 (8%)	10	22
1	B	155/159 (98%)	142 (92%)	13 (8%)	10	22
1	C	154/159 (97%)	142 (92%)	12 (8%)	11	24
1	D	154/159 (97%)	139 (90%)	15 (10%)	8	17
All	All	618/636 (97%)	565 (91%)	53 (9%)	10	21

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	53	LYS
1	A	55	GLN
1	A	77	GLN
1	A	118	ASN
1	A	133	ASP
1	A	154	LEU
1	A	158	GLU
1	A	176	LEU
1	A	206	GLU
1	A	207	ASP
1	A	210	GLU
1	A	217	GLU
1	A	222	ILE
1	B	43	THR
1	B	55	GLN
1	B	62	SER
1	B	68	THR
1	B	79	LYS
1	B	116	ARG
1	B	123	VAL
1	B	143	LYS
1	B	144	ASN
1	B	146	ARG
1	B	154	LEU
1	B	211	ASN
1	B	217	GLU
1	C	53	LYS
1	C	77	GLN
1	C	118	ASN
1	C	133	ASP
1	C	154	LEU
1	C	158	GLU
1	C	176	LEU

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Mol	Chain	Res	Type
1	C	194	GLN
1	C	209	GLN
1	C	210	GLU
1	C	222	ILE
1	C	223	LEU
1	D	43	THR
1	D	55	GLN
1	D	62	SER
1	D	68	THR
1	D	79	LYS
1	D	116	ARG
1	D	118	ASN
1	D	123	VAL
1	D	143	LYS
1	D	144	ASN
1	D	146	ARG
1	D	154	LEU
1	D	203	GLN
1	D	206	GLU
1	D	224	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	50	ASN
1	A	55	GLN
1	A	77	GLN
1	A	182	GLN
1	A	211	ASN
1	B	50	ASN
1	B	55	GLN
1	B	82	GLN
1	B	94	HIS
1	B	144	ASN
1	B	163	GLN
1	B	211	ASN
1	C	50	ASN
1	C	55	GLN
1	C	77	GLN
1	C	98	HIS
1	C	182	GLN
1	C	194	GLN

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Mol	Chain	Res	Type
1	D	50	ASN
1	D	55	GLN
1	D	82	GLN
1	D	94	HIS
1	D	144	ASN
1	D	163	GLN
1	D	189	GLN
1	D	194	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 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 ⓘ

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	185/190 (97%)	0.74	20 (10%) 11 9	9, 27, 49, 62	0
1	B	190/190 (100%)	0.99	30 (15%) 5 4	14, 30, 58, 79	0
1	C	185/190 (97%)	0.77	23 (12%) 8 6	9, 27, 51, 73	0
1	D	179/190 (94%)	0.81	17 (9%) 14 12	14, 31, 48, 61	0
All	All	739/760 (97%)	0.83	90 (12%) 8 7	9, 29, 51, 79	0

All (90) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	199	SER	5.3
1	C	197	GLY	4.7
1	A	36	SER	4.6
1	C	201	ALA	4.5
1	B	201	ALA	4.5
1	A	196	GLU	4.4
1	B	46	ARG	4.4
1	A	39	GLY	4.3
1	B	224	LYS	4.3
1	B	35	GLY	4.3
1	B	199	SER	4.3
1	C	198	GLU	4.2
1	B	198	GLU	4.1
1	B	78	GLY	4.1
1	A	37	GLY	3.8
1	B	200	LYS	3.8
1	B	196	GLU	3.8
1	C	81	GLY	3.8
1	B	197	GLY	3.7
1	A	38	GLY	3.4
1	B	80	GLY	3.4

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Mol	Chain	Res	Type	RSRZ
1	B	36	SER	3.4
1	B	38	GLY	3.3
1	A	79	LYS	3.3
1	C	195	GLU	3.2
1	B	37	GLY	3.2
1	D	79	LYS	3.1
1	C	94	HIS	3.1
1	A	195	GLU	3.0
1	A	62	SER	3.0
1	B	134	ARG	2.9
1	C	79	LYS	2.8
1	A	80	GLY	2.7
1	C	62	SER	2.7
1	B	203	GLN	2.7
1	A	222	ILE	2.7
1	C	90	LEU	2.7
1	C	196	GLU	2.7
1	B	204	PRO	2.6
1	A	63	SER	2.6
1	C	106	THR	2.6
1	D	78	GLY	2.5
1	C	107	ILE	2.5
1	B	82	GLN	2.5
1	B	63	SER	2.5
1	A	95	ALA	2.5
1	D	109	PRO	2.5
1	A	206	GLU	2.5
1	A	194	GLN	2.4
1	A	46	ARG	2.4
1	D	46	ARG	2.4
1	D	114	ALA	2.4
1	C	207	ASP	2.4
1	B	43	THR	2.4
1	D	101	GLU	2.4
1	C	93	GLU	2.4
1	D	203	GLN	2.3
1	B	79	LYS	2.3
1	A	158	GLU	2.3
1	C	80	GLY	2.3
1	B	101	GLU	2.3
1	D	195	GLU	2.3
1	D	110	ALA	2.2

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Mol	Chain	Res	Type	RSRZ
1	D	142	THR	2.2
1	B	39	GLY	2.2
1	C	224	LYS	2.2
1	B	195	GLU	2.2
1	D	206	GLU	2.2
1	C	43	THR	2.2
1	C	222	ILE	2.2
1	C	83	VAL	2.2
1	B	75	GLU	2.2
1	D	40	MET	2.1
1	A	43	THR	2.1
1	D	112	ASP	2.1
1	C	141	LYS	2.1
1	B	77	GLN	2.1
1	B	90	LEU	2.1
1	D	80	GLY	2.1
1	D	63	SER	2.1
1	C	82	GLN	2.1
1	A	65	ARG	2.1
1	B	111	PHE	2.1
1	B	76	ALA	2.1
1	A	211	ASN	2.1
1	D	224	LYS	2.1
1	B	140	SER	2.1
1	A	40	MET	2.1
1	D	111	PHE	2.0
1	C	223	LEU	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	ZN	D	2003	1/1	0.93	0.07	52,52,52,52	0
2	ZN	B	2001	1/1	0.98	0.03	31,31,31,31	0
2	ZN	A	2000	1/1	0.99	0.03	25,25,25,25	0
2	ZN	C	2002	1/1	1.00	0.02	28,28,28,28	0

6.5 Other polymers [i](#)

There are no such residues in this entry.