



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 08:38 AM UTC

PDB ID : 3Q05 / pdb_00003q05
Title : An induced fit mechanism regulates p53 DNA binding kinetics to confer sequence specificity
Authors : Petty, T.J.; Halazonetis, T.D.
Deposited on : 2010-12-15
Resolution : 2.40 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

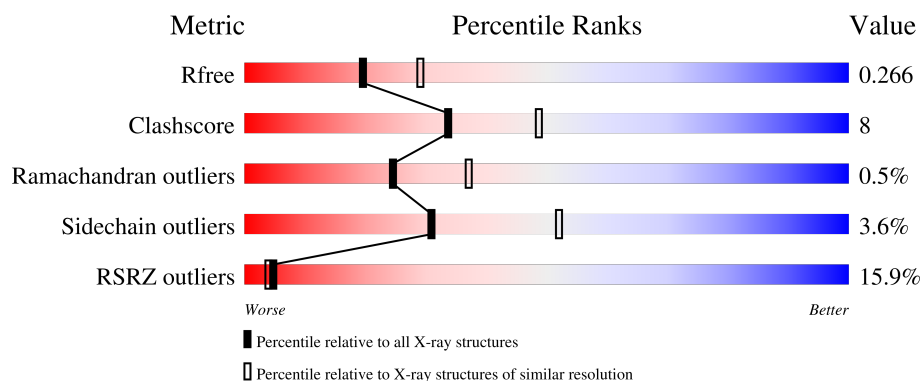
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	234	12% 84% 13% .
1	B	234	27% 70% 28% .
1	C	234	7% 84% 14% .
1	D	234	15% 85% 12% .
2	K	26	27% 50% 50% .

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Mol	Chain	Length	Quality of chain
3	L	26	 A horizontal bar chart showing the quality of chain L. The bar is divided into three segments: red (31%), green (62%), and yellow (38%). The segments are stacked horizontally, with red on the left, green in the middle, and yellow on the right. The percentages are labeled above each segment.

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8677 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 Cellular tumor antigen p53.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	234	Total	C	N	O	S	0	0	0
			1863	1157	339	354	13			
1	C	234	Total	C	N	O	S	0	0	0
			1863	1157	339	354	13			
1	D	234	Total	C	N	O	S	0	0	0
			1863	1157	339	354	13			
1	B	234	Total	C	N	O	S	0	0	0
			1863	1157	339	354	13			

There are 72 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	135	VAL	CYS	engineered mutation	UNP P04637
A	141	VAL	CYS	engineered mutation	UNP P04637
A	146	TYR	TRP	engineered mutation	UNP P04637
A	182	SER	CYS	engineered mutation	UNP P04637
A	203	ALA	VAL	engineered mutation	UNP P04637
A	209	PRO	ARG	engineered mutation	UNP P04637
A	229	TYR	CYS	engineered mutation	UNP P04637
A	233	TYR	HIS	engineered mutation	UNP P04637
A	234	PHE	TYR	engineered mutation	UNP P04637
A	235	LYS	ASN	engineered mutation	UNP P04637
A	236	PHE	TYR	engineered mutation	UNP P04637
A	253	VAL	THR	engineered mutation	UNP P04637
A	268	ASP	ASN	engineered mutation	UNP P04637
A	322	THR	PRO	engineered mutation	UNP P04637
A	323	MET	LEU	engineered mutation	UNP P04637
A	340	GLN	MET	engineered mutation	UNP P04637
A	344	ARG	LEU	engineered mutation	UNP P04637
A	356	THR	GLY	engineered mutation	UNP P04637
C	135	VAL	CYS	engineered mutation	UNP P04637
C	141	VAL	CYS	engineered mutation	UNP P04637
C	146	TYR	TRP	engineered mutation	UNP P04637

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Chain	Residue	Modelled	Actual	Comment	Reference
C	182	SER	CYS	engineered mutation	UNP P04637
C	203	ALA	VAL	engineered mutation	UNP P04637
C	209	PRO	ARG	engineered mutation	UNP P04637
C	229	TYR	CYS	engineered mutation	UNP P04637
C	233	TYR	HIS	engineered mutation	UNP P04637
C	234	PHE	TYR	engineered mutation	UNP P04637
C	235	LYS	ASN	engineered mutation	UNP P04637
C	236	PHE	TYR	engineered mutation	UNP P04637
C	253	VAL	THR	engineered mutation	UNP P04637
C	268	ASP	ASN	engineered mutation	UNP P04637
C	322	THR	PRO	engineered mutation	UNP P04637
C	323	MET	LEU	engineered mutation	UNP P04637
C	340	GLN	MET	engineered mutation	UNP P04637
C	344	ARG	LEU	engineered mutation	UNP P04637
C	356	THR	GLY	engineered mutation	UNP P04637
D	135	VAL	CYS	engineered mutation	UNP P04637
D	141	VAL	CYS	engineered mutation	UNP P04637
D	146	TYR	TRP	engineered mutation	UNP P04637
D	182	SER	CYS	engineered mutation	UNP P04637
D	203	ALA	VAL	engineered mutation	UNP P04637
D	209	PRO	ARG	engineered mutation	UNP P04637
D	229	TYR	CYS	engineered mutation	UNP P04637
D	233	TYR	HIS	engineered mutation	UNP P04637
D	234	PHE	TYR	engineered mutation	UNP P04637
D	235	LYS	ASN	engineered mutation	UNP P04637
D	236	PHE	TYR	engineered mutation	UNP P04637
D	253	VAL	THR	engineered mutation	UNP P04637
D	268	ASP	ASN	engineered mutation	UNP P04637
D	322	THR	PRO	engineered mutation	UNP P04637
D	323	MET	LEU	engineered mutation	UNP P04637
D	340	GLN	MET	engineered mutation	UNP P04637
D	344	ARG	LEU	engineered mutation	UNP P04637
D	356	THR	GLY	engineered mutation	UNP P04637
B	135	VAL	CYS	engineered mutation	UNP P04637
B	141	VAL	CYS	engineered mutation	UNP P04637
B	146	TYR	TRP	engineered mutation	UNP P04637
B	182	SER	CYS	engineered mutation	UNP P04637
B	203	ALA	VAL	engineered mutation	UNP P04637
B	209	PRO	ARG	engineered mutation	UNP P04637
B	229	TYR	CYS	engineered mutation	UNP P04637
B	233	TYR	HIS	engineered mutation	UNP P04637
B	234	PHE	TYR	engineered mutation	UNP P04637

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Chain	Residue	Modelled	Actual	Comment	Reference
B	235	LYS	ASN	engineered mutation	UNP P04637
B	236	PHE	TYR	engineered mutation	UNP P04637
B	253	VAL	THR	engineered mutation	UNP P04637
B	268	ASP	ASN	engineered mutation	UNP P04637
B	322	THR	PRO	engineered mutation	UNP P04637
B	323	MET	LEU	engineered mutation	UNP P04637
B	340	GLN	MET	engineered mutation	UNP P04637
B	344	ARG	LEU	engineered mutation	UNP P04637
B	356	THR	GLY	engineered mutation	UNP P04637

- Molecule 2 is a DNA chain called DNA (26-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	K	26	Total	C	N	O	P	0	0	0
			527	252	96	154	25			

- Molecule 3 is a DNA chain called DNA (26-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	L	26	Total	C	N	O	P	0	0	0
			533	254	100	154	25			

- Molecule 4 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Zn	0	0
			1	1		
4	C	1	Total	Zn	0	0
			1	1		
4	D	1	Total	Zn	0	0
			1	1		
4	B	1	Total	Zn	0	0
			1	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	34	Total	O	0	0
			34	34		
5	C	66	Total	O	0	0
			66	66		

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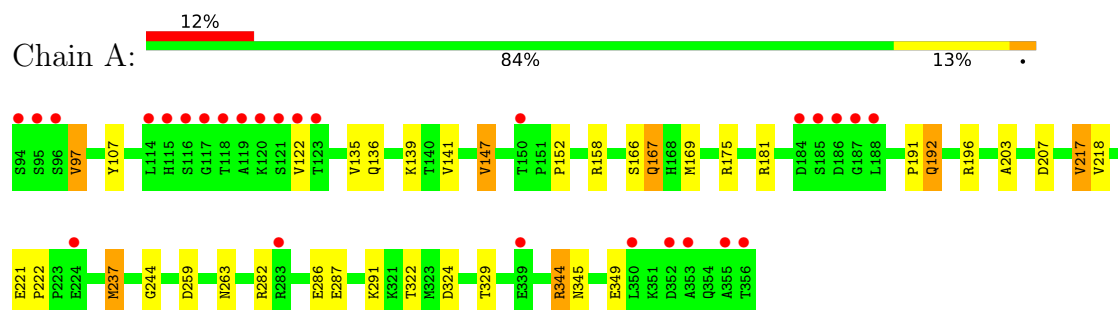
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	D	47	Total 47	O 47	0	0
5	B	7	Total 7	O 7	0	0
5	K	4	Total 4	O 4	0	0
5	L	3	Total 3	O 3	0	0

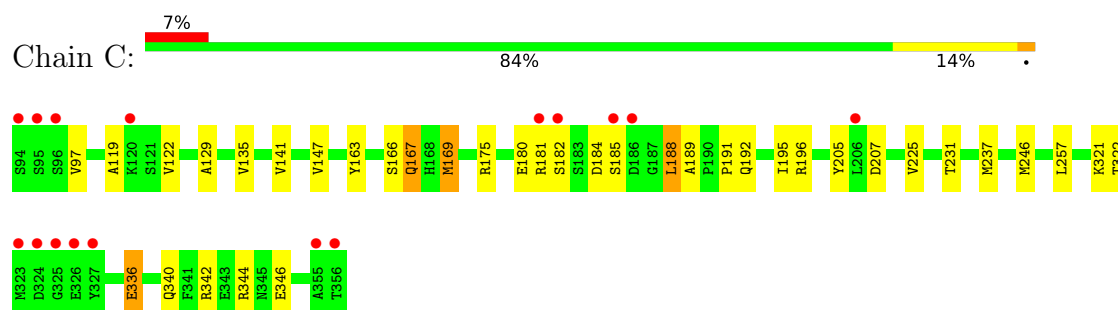
3 Residue-property plots [i](#)

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.

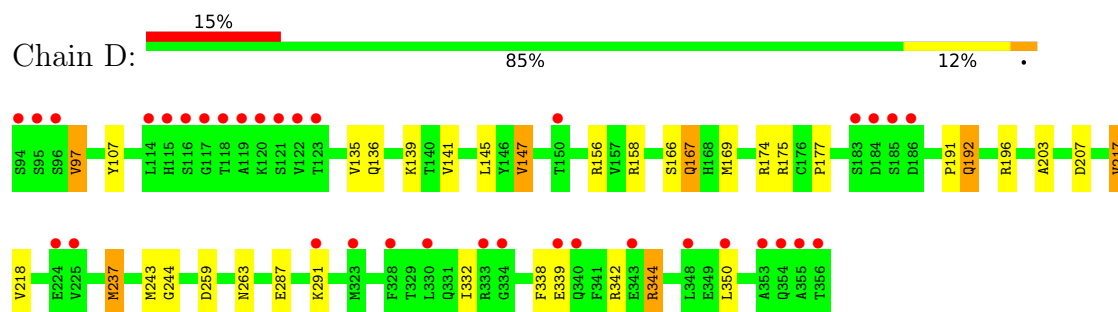
• Molecule 1: Cellular tumor antigen p53



• Molecule 1: Cellular tumor antigen p53

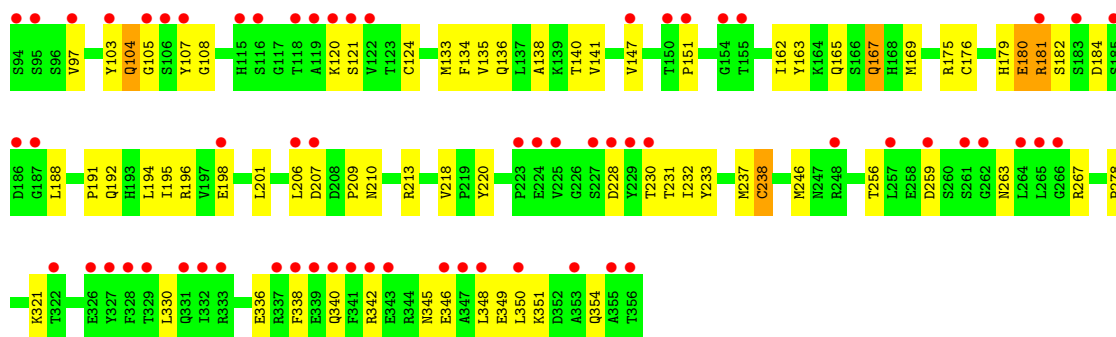


• Molecule 1: Cellular tumor antigen p53



• Molecule 1: Cellular tumor antigen p53





• Molecule 2: DNA (26-MER)



• Molecule 3: DNA (26-MER)



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	163.30Å 169.76Å 55.66Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	41.07 – 2.40 41.07 – 2.40	Depositor EDS
% Data completeness (in resolution range)	97.3 (41.07-2.40) 97.2 (41.07-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.24 (at 2.39Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.238 , 0.267 0.238 , 0.266	Depositor DCC
R_{free} test set	3005 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	45.7	Xtriage
Anisotropy	0.323	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 45.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.009 for k,h,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	8677	wwPDB-VP
Average B, all atoms (Å ²)	54.0	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 31.72 % 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 1.0569e-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: 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	0.42	0/1903	0.90	3/2571 (0.1%)
1	B	0.37	0/1903	0.86	4/2571 (0.2%)
1	C	0.50	0/1903	0.94	5/2571 (0.2%)
1	D	0.41	0/1903	0.89	3/2571 (0.1%)
2	K	0.30	0/590	0.72	0/908
3	L	0.29	0/598	0.68	0/922
All	All	0.41	0/8800	0.87	15/12114 (0.1%)

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	K	0	1
3	L	0	1
All	All	0	2

There are no bond length outliers.

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	237	MET	N-CA-C	8.71	122.69	112.93
1	D	237	MET	N-CA-C	8.49	122.44	112.93
1	A	237	MET	N-CA-C	8.12	122.54	113.21
1	B	237	MET	N-CA-C	7.69	121.54	112.93
1	A	244	GLY	N-CA-C	-6.73	107.07	115.08
1	D	244	GLY	N-CA-C	-6.23	107.11	115.21
1	D	97	VAL	N-CA-C	5.97	113.79	107.76
1	C	188	LEU	N-CA-C	-5.89	107.91	114.62

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	97	VAL	N-CA-C	5.86	113.67	107.76
1	C	195	ILE	N-CA-C	5.56	116.59	108.58
1	B	180	GLU	N-CA-C	-5.26	105.62	111.36
1	C	169	MET	N-CA-C	5.22	117.71	111.71
1	C	257	LEU	N-CA-C	-5.03	100.98	109.07
1	B	138	ALA	N-CA-C	5.02	118.36	111.54
1	B	195	ILE	N-CA-C	5.00	115.78	108.58

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	K	13	DC	Sidechain
3	L	37	DC	Sidechain

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	1863	0	1822	27	0
1	B	1863	0	1822	39	0
1	C	1863	0	1822	18	0
1	D	1863	0	1822	30	0
2	K	527	0	294	10	0
3	L	533	0	294	9	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
5	A	34	0	0	0	0
5	B	7	0	0	0	0
5	C	66	0	0	2	0
5	D	47	0	0	2	0
5	K	4	0	0	0	0
5	L	3	0	0	0	0
All	All	8677	0	7876	126	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 (126) 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:167:GLN:HE21	1:A:167:GLN:H	1.17	0.92
1:D:167:GLN:H	1:D:167:GLN:HE21	1.14	0.90
1:A:192:GLN:HE21	1:A:192:GLN:H	1.31	0.77
1:D:192:GLN:NE2	1:D:192:GLN:H	1.83	0.77
1:A:192:GLN:H	1:A:192:GLN:NE2	1.83	0.76
1:D:192:GLN:H	1:D:192:GLN:HE21	1.32	0.76
3:L:30:DG:H2''	3:L:31:DG:H5''	1.68	0.74
1:C:167:GLN:HE21	1:C:167:GLN:H	1.35	0.72
3:L:33:DA:H2''	3:L:34:DT:H5'	1.72	0.71
1:D:167:GLN:H	1:D:167:GLN:NE2	1.88	0.68
1:A:167:GLN:H	1:A:167:GLN:NE2	1.90	0.68
1:B:167:GLN:NE2	1:B:167:GLN:H	1.92	0.67
1:C:97:VAL:HG11	1:C:169:MET:HE2	1.76	0.66
1:B:184:ASP:HB2	1:B:196:ARG:HH22	1.61	0.66
1:B:97:VAL:HG11	1:B:169:MET:HE2	1.77	0.65
1:B:167:GLN:H	1:B:167:GLN:HE21	1.46	0.63
1:D:342:ARG:NH2	1:D:342:ARG:HB3	2.14	0.63
1:D:167:GLN:HE21	1:D:167:GLN:N	1.92	0.62
1:D:97:VAL:HG11	1:D:169:MET:HE2	1.81	0.62
1:C:129:ALA:HB2	1:C:346:GLU:HB3	1.80	0.61
1:A:97:VAL:HG11	1:A:169:MET:HE2	1.81	0.61
1:D:344:ARG:CZ	1:D:344:ARG:HA	2.31	0.61
1:A:181:ARG:NH2	1:B:181:ARG:HH22	1.98	0.61
1:A:192:GLN:HE21	1:A:192:GLN:N	1.99	0.60
1:D:342:ARG:HB3	1:D:342:ARG:HH21	1.67	0.59
2:K:23:DC:H2''	2:K:24:DC:H5'	1.84	0.58
1:C:119:ALA:O	1:C:122:VAL:HG22	2.04	0.58
1:A:167:GLN:HE21	1:A:167:GLN:N	1.95	0.57
1:C:184:ASP:HB2	1:C:196:ARG:HH22	1.69	0.57
1:D:192:GLN:HE21	1:D:192:GLN:N	2.00	0.57
1:D:339:GLU:HA	1:D:342:ARG:HG3	1.86	0.57
1:B:256:THR:HG22	1:B:267:ARG:HD3	1.87	0.57
1:D:107:TYR:HB3	1:D:147:VAL:HG22	1.87	0.56
1:B:135:VAL:HG21	1:B:141:VAL:HG22	1.87	0.56
1:C:167:GLN:H	1:C:167:GLN:NE2	2.04	0.56
1:B:259:ASP:OD2	1:B:263:ASN:HB2	2.07	0.55
1:A:107:TYR:HB3	1:A:147:VAL:HG22	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:166:SER:HA	1:A:169:MET:HG3	1.89	0.54
1:D:166:SER:HA	1:D:169:MET:HG3	1.90	0.53
1:A:329:THR:HG23	5:C:1145:HOH:O	2.08	0.53
1:A:158:ARG:HB2	1:A:217:VAL:HG13	1.92	0.52
1:D:158:ARG:HB2	1:D:217:VAL:HG13	1.91	0.52
1:B:151:PRO:HD2	1:B:220:TYR:CE2	2.45	0.51
1:B:147:VAL:HG21	1:B:151:PRO:HD3	1.92	0.51
1:D:338:PHE:HE1	1:B:330:LEU:HB2	1.74	0.51
1:B:218:VAL:HG21	1:B:232:ILE:HD12	1.92	0.50
2:K:27:DT:H2''	2:K:28:DT:H71	1.93	0.50
2:K:11:DG:H1'	2:K:12:DC:H5''	1.93	0.50
1:B:135:VAL:HG22	1:B:136:GLN:N	2.27	0.50
1:D:332:ILE:HG13	1:D:338:PHE:HD1	1.76	0.50
3:L:30:DG:C2'	3:L:31:DG:H5''	2.40	0.50
2:K:7:DA:H2''	2:K:8:DC:O5'	2.11	0.49
1:C:167:GLN:HE21	1:C:167:GLN:N	2.06	0.49
1:D:259:ASP:OD2	1:D:263:ASN:HB2	2.12	0.49
1:B:103:TYR:CE2	1:B:105:GLY:HA2	2.47	0.49
1:B:209:PRO:HG2	1:B:210:ASN:OD1	2.12	0.49
2:K:21:DG:H2''	2:K:22:DC:H5'	1.94	0.49
1:B:338:PHE:CZ	1:B:342:ARG:HD2	2.48	0.48
1:C:175:ARG:HD3	1:C:191:PRO:O	2.13	0.48
1:C:342:ARG:O	1:C:346:GLU:HG3	2.14	0.48
1:B:124:CYS:HB2	1:B:135:VAL:HG23	1.96	0.48
1:A:259:ASP:OD2	1:A:263:ASN:HB2	2.13	0.47
1:D:350:LEU:HD23	1:D:350:LEU:O	2.14	0.47
1:A:122:VAL:HB	3:L:28:DC:OP2	2.14	0.47
1:B:176:CYS:O	1:B:180:GLU:HB2	2.13	0.47
3:L:34:DT:H1'	3:L:35:DG:C8	2.50	0.47
1:A:135:VAL:HG21	1:A:141:VAL:HG22	1.97	0.46
5:C:1028:HOH:O	1:D:243:MET:HE1	2.14	0.46
1:D:135:VAL:HG21	1:D:141:VAL:HG22	1.97	0.46
1:B:162:ILE:HG22	1:B:213:ARG:NH1	2.30	0.46
1:A:136:GLN:HB2	1:A:139:LYS:HG3	1.97	0.46
1:C:321:LYS:HE3	1:C:322:THR:HG23	1.96	0.46
1:B:135:VAL:HG21	1:B:141:VAL:CG2	2.45	0.46
1:D:136:GLN:HB2	1:D:139:LYS:HG3	1.98	0.46
1:B:165:GLN:O	1:B:169:MET:HG3	2.15	0.46
1:D:344:ARG:HA	1:D:344:ARG:NE	2.30	0.46
1:B:201:LEU:HD22	1:B:201:LEU:N	2.31	0.46
1:A:135:VAL:HG22	1:A:136:GLN:N	2.31	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:134:PHE:HB3	1:B:278:PRO:HB3	1.98	0.46
1:A:175:ARG:HD3	1:A:191:PRO:O	2.15	0.45
1:B:175:ARG:NH2	1:B:179:HIS:HB3	2.31	0.45
1:D:203:ALA:HA	1:D:218:VAL:HG12	1.98	0.45
1:B:350:LEU:O	1:B:354:GLN:HG2	2.16	0.45
1:D:174:ARG:NE	5:D:1022:HOH:O	2.49	0.45
1:A:345:ASN:O	1:A:349:GLU:HG2	2.17	0.45
1:B:133:MET:HE3	1:B:141:VAL:CG1	2.47	0.45
1:B:140:THR:HG23	1:B:233:TYR:HB3	1.97	0.45
1:C:180:GLU:C	1:C:182:SER:H	2.24	0.45
3:L:48:DT:H2''	3:L:49:DC:C6	2.50	0.45
1:B:107:TYR:CE1	1:B:151:PRO:HA	2.52	0.45
2:K:27:DT:H2''	2:K:28:DT:C5	2.52	0.45
1:C:166:SER:HA	1:C:169:MET:HG3	1.98	0.45
1:B:206:LEU:HD12	1:B:207:ASP:H	1.82	0.45
1:D:156:ARG:HD3	5:D:1055:HOH:O	2.16	0.44
1:A:287:GLU:HG3	1:A:291:LYS:HE3	2.00	0.44
1:C:181:ARG:HH11	1:D:177:PRO:HG3	1.81	0.44
1:D:135:VAL:HG22	1:D:136:GLN:N	2.31	0.44
1:B:104:GLN:NE2	1:B:108:GLY:HA2	2.33	0.44
3:L:49:DC:H2''	3:L:50:DA:C8	2.52	0.44
1:B:342:ARG:O	1:B:346:GLU:HG3	2.18	0.44
2:K:27:DT:H2''	2:K:28:DT:C7	2.48	0.44
1:D:175:ARG:HD3	1:D:191:PRO:O	2.17	0.43
1:A:203:ALA:HA	1:A:218:VAL:HG12	2.00	0.43
1:B:163:TYR:OH	1:B:246:MET:HA	2.18	0.43
1:B:175:ARG:HD3	1:B:191:PRO:O	2.18	0.43
1:D:287:GLU:HG3	1:D:291:LYS:HE3	2.00	0.42
1:B:194:LEU:CD1	1:B:238:CYS:HB2	2.49	0.42
1:B:345:ASN:O	1:B:349:GLU:HG3	2.20	0.42
1:A:344:ARG:HG2	1:C:344:ARG:HH12	1.85	0.42
1:A:196:ARG:NH1	1:A:237:MET:HE2	2.34	0.42
1:C:336:GLU:O	1:C:340:GLN:HG3	2.19	0.42
1:B:348:LEU:HD23	1:B:351:LYS:HD2	2.01	0.42
1:A:221:GLU:HA	1:A:222:PRO:HD3	1.94	0.42
1:B:188:LEU:HD23	1:B:188:LEU:O	2.19	0.42
1:C:135:VAL:HG21	1:C:141:VAL:HG22	2.01	0.41
2:K:4:DG:H1	3:L:49:DC:H42	1.67	0.41
1:C:163:TYR:OH	1:C:246:MET:HA	2.21	0.41
1:A:107:TYR:OH	1:A:152:PRO:HD3	2.20	0.41
1:B:140:THR:CG2	1:B:233:TYR:HB3	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:282:ARG:O	1:A:286:GLU:HG3	2.21	0.41
1:C:189:ALA:HB2	1:C:205:TYR:CZ	2.56	0.41
1:B:336:GLU:O	1:B:340:GLN:HG2	2.21	0.41
2:K:19:DA:H2	3:L:34:DT:O2	2.04	0.41
1:D:196:ARG:NH1	1:D:237:MET:HE2	2.36	0.40
1:A:322:THR:C	1:A:324:ASP:H	2.29	0.40
2:K:23:DC:C2'	2:K:24:DC:H5'	2.50	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	232/234 (99%)	222 (96%)	10 (4%)	0	100	100
1	B	232/234 (99%)	217 (94%)	11 (5%)	4 (2%)	7	10
1	C	232/234 (99%)	219 (94%)	12 (5%)	1 (0%)	30	43
1	D	232/234 (99%)	222 (96%)	10 (4%)	0	100	100
All	All	928/936 (99%)	880 (95%)	43 (5%)	5 (0%)	24	37

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	121	SER
1	B	120	LYS
1	B	182	SER
1	C	185	SER
1	B	228	ASP

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	208/208 (100%)	202 (97%)	6 (3%)	37	60
1	B	208/208 (100%)	199 (96%)	9 (4%)	26	44
1	C	208/208 (100%)	200 (96%)	8 (4%)	29	49
1	D	208/208 (100%)	201 (97%)	7 (3%)	32	54
All	All	832/832 (100%)	802 (96%)	30 (4%)	31	52

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	147	VAL
1	A	167	GLN
1	A	192	GLN
1	A	207	ASP
1	A	217	VAL
1	A	344	ARG
1	C	147	VAL
1	C	167	GLN
1	C	188	LEU
1	C	192	GLN
1	C	207	ASP
1	C	225	VAL
1	C	231	THR
1	C	336	GLU
1	D	145	LEU
1	D	147	VAL
1	D	167	GLN
1	D	192	GLN
1	D	207	ASP
1	D	217	VAL
1	D	344	ARG
1	B	104	GLN
1	B	167	GLN
1	B	181	ARG

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Mol	Chain	Res	Type
1	B	192	GLN
1	B	198	GLU
1	B	230	THR
1	B	231	THR
1	B	238	CYS
1	B	321	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	104	GLN
1	A	167	GLN
1	A	192	GLN
1	A	239	ASN
1	C	104	GLN
1	C	136	GLN
1	C	144	GLN
1	C	165	GLN
1	C	167	GLN
1	C	168	HIS
1	C	192	GLN
1	C	247	ASN
1	C	340	GLN
1	C	345	ASN
1	D	104	GLN
1	D	167	GLN
1	D	192	GLN
1	D	239	ASN
1	B	100	GLN
1	B	104	GLN
1	B	144	GLN
1	B	165	GLN
1	B	167	GLN
1	B	168	HIS
1	B	192	GLN
1	B	247	ASN
1	B	340	GLN

5.3.3 RNA ⓘ

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

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	234/234 (100%)	0.64	27 (11%) 9 7	28, 44, 94, 95	0
1	B	234/234 (100%)	1.32	64 (27%) 1 1	41, 72, 95, 95	0
1	C	234/234 (100%)	0.06	16 (6%) 23 20	18, 34, 82, 95	0
1	D	234/234 (100%)	0.78	35 (14%) 5 4	26, 43, 95, 95	0
2	K	26/26 (100%)	1.00	7 (26%) 1 1	33, 52, 86, 91	0
3	L	26/26 (100%)	1.14	8 (30%) 1 1	38, 55, 85, 91	0
All	All	988/988 (100%)	0.72	157 (15%) 5 4	18, 47, 95, 95	0

All (157) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	119	ALA	9.4
1	A	122	VAL	7.0
1	B	120	LYS	6.9
1	D	355	ALA	5.5
1	D	186	ASP	5.2
1	A	119	ALA	5.2
1	D	118	THR	5.0
3	L	28	DC	4.6
1	B	95	SER	4.5
1	A	95	SER	4.5
1	A	114	LEU	4.5
1	A	115	HIS	4.3
1	D	116	SER	4.2
1	D	122	VAL	4.2
1	B	356	THR	4.1
1	D	356	THR	3.9
1	A	186	ASP	3.9
1	D	120	LYS	3.9
1	C	94	SER	3.9

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Mol	Chain	Res	Type	RSRZ
1	A	94	SER	3.9
1	A	118	THR	3.9
1	D	95	SER	3.8
1	A	121	SER	3.8
2	K	4	DG	3.7
3	L	27	DA	3.6
1	C	356	THR	3.6
1	C	95	SER	3.6
1	C	185	SER	3.5
1	B	228	ASP	3.5
1	D	121	SER	3.5
1	B	329	THR	3.4
1	B	187	GLY	3.4
1	D	184	ASP	3.3
1	D	115	HIS	3.3
1	D	225	VAL	3.3
1	A	353	ALA	3.3
1	D	323	MET	3.3
1	A	120	LYS	3.3
1	B	94	SER	3.3
3	L	51	DA	3.3
1	C	355	ALA	3.3
1	A	96	SER	3.2
1	D	183	SER	3.2
2	K	5	DA	3.2
1	D	114	LEU	3.2
1	B	229	TYR	3.2
1	A	117	GLY	3.2
1	B	262	GLY	3.2
1	B	350	LEU	3.1
1	B	264	LEU	3.1
3	L	29	DG	3.1
1	B	353	ALA	3.1
1	D	353	ALA	3.1
1	D	94	SER	3.1
1	A	150	THR	3.0
1	C	186	ASP	3.0
1	C	325	GLY	3.0
1	C	323	MET	3.0
1	D	350	LEU	3.0
2	K	25	DG	2.9
1	C	96	SER	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	206	LEU	2.9
1	B	348	LEU	2.9
1	A	116	SER	2.9
1	B	339	GLU	2.9
1	C	324	ASP	2.9
1	A	184	ASP	2.8
1	D	291	LYS	2.8
1	B	181	ARG	2.8
1	B	183	SER	2.8
1	A	355	ALA	2.8
1	D	340	GLN	2.8
1	B	327	TYR	2.7
1	D	185	SER	2.7
1	A	350	LEU	2.7
1	C	327	TYR	2.7
1	D	330	LEU	2.7
1	B	225	VAL	2.7
1	A	339	GLU	2.7
1	B	119	ALA	2.7
1	B	355	ALA	2.7
1	A	123	THR	2.6
1	B	118	THR	2.6
1	B	105	GLY	2.6
1	B	186	ASP	2.6
1	B	322	THR	2.6
2	K	24	DC	2.6
1	B	341	PHE	2.5
1	A	356	THR	2.5
1	D	96	SER	2.5
1	B	333	ARG	2.5
2	K	27	DT	2.5
1	B	151	PRO	2.5
1	B	332	ILE	2.5
1	B	265	LEU	2.5
1	C	120	LYS	2.5
1	B	122	VAL	2.5
1	A	185	SER	2.5
1	D	334	GLY	2.4
1	B	224	GLU	2.4
1	B	346	GLU	2.4
3	L	52	DA	2.4
1	B	115	HIS	2.4

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Mol	Chain	Res	Type	RSRZ
1	D	333	ARG	2.4
3	L	50	DA	2.4
1	D	117	GLY	2.4
1	B	103	TYR	2.4
1	B	338	PHE	2.3
1	B	227	SER	2.3
3	L	49	DC	2.3
1	D	328	PHE	2.3
1	A	224	GLU	2.3
1	D	123	THR	2.3
1	B	230	THR	2.3
1	B	248	ARG	2.3
1	A	188	LEU	2.3
1	B	207	ASP	2.3
1	D	348	LEU	2.3
1	B	97	VAL	2.3
1	B	223	PRO	2.3
1	A	187	GLY	2.2
1	B	266	GLY	2.2
1	B	116	SER	2.2
1	C	326	GLU	2.2
1	C	181	ARG	2.2
1	B	337	ARG	2.2
1	C	182	SER	2.2
1	B	328	PHE	2.2
1	B	147	VAL	2.2
1	B	326	GLU	2.2
1	B	342	ARG	2.2
1	D	150	THR	2.2
1	B	107	TYR	2.2
1	B	150	THR	2.2
1	B	121	SER	2.2
1	B	185	SER	2.2
3	L	48	DT	2.1
1	B	257	LEU	2.1
1	D	224	GLU	2.1
1	D	343	GLU	2.1
2	K	6	DG	2.1
1	B	331	GLN	2.1
1	B	347	ALA	2.1
1	B	106	SER	2.1
1	C	206	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	339	GLU	2.1
1	B	259	ASP	2.1
1	B	340	GLN	2.1
1	A	283	ARG	2.1
1	B	154	GLY	2.1
1	B	261	SER	2.1
2	K	3	DT	2.1
1	B	155	THR	2.0
1	D	354	GLN	2.0
1	B	198	GLU	2.0
1	B	343	GLU	2.0
1	A	352	ASP	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
4	ZN	C	1	1/1	0.99	0.05	30,30,30,30	0
4	ZN	D	1	1/1	0.99	0.06	31,31,31,31	0
4	ZN	B	1	1/1	0.99	0.06	49,49,49,49	0
4	ZN	A	1	1/1	1.00	0.06	43,43,43,43	0

6.5 Other polymers [i](#)

There are no such residues in this entry.