



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 1, 2016 – 02:29 AM GMT

PDB ID : 2H1L
Title : The Structure of the Oncoprotein SV40 Large T Antigen and p53 Tumor
Suppressor Complex
Authors : Lilyestrom, W.
Deposited on : 2006-05-16
Resolution : 3.16 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

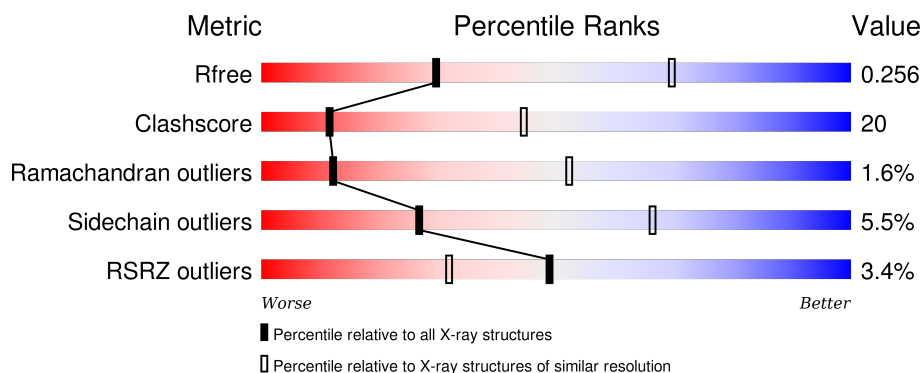
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.16 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}	91344	1112 (3.20-3.12)
Clashscore	102246	1249 (3.20-3.12)
Ramachandran outliers	100387	1222 (3.20-3.12)
Sidechain outliers	100360	1221 (3.20-3.12)
RSRZ outliers	91569	1117 (3.20-3.12)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	370	<div> <div>2%</div> <div>56%</div> <div>38%</div> <div>• •</div> </div>
1	B	370	<div> <div>50%</div> <div>43%</div> <div>5%</div> <div>•</div> </div>
1	C	370	<div> <div>2%</div> <div>52%</div> <div>41%</div> <div>5%</div> <div>•</div> </div>
1	D	370	<div> <div>%</div> <div>55%</div> <div>41%</div> <div>• •</div> </div>
1	E	370	<div> <div>%</div> <div>54%</div> <div>39%</div> <div>• •</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	370	
1	G	370	
1	H	370	
1	I	370	
1	J	370	
1	K	370	
1	L	370	
2	M	203	
2	N	203	
2	O	203	
2	P	203	
2	Q	203	
2	R	203	
2	S	203	
2	T	203	
2	U	203	
2	V	203	
2	W	203	
2	X	203	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 53920 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 Large T antigen.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	362	Total	C	N	O	S	0	0	0
			2933	1888	493	531	21			
1	B	362	Total	C	N	O	S	0	0	0
			2933	1888	493	531	21			
1	C	363	Total	C	N	O	S	0	0	0
			2940	1892	494	533	21			
1	D	362	Total	C	N	O	S	0	0	0
			2933	1888	493	531	21			
1	E	362	Total	C	N	O	S	0	0	0
			2933	1888	493	531	21			
1	F	363	Total	C	N	O	S	0	0	0
			2940	1892	494	533	21			
1	G	362	Total	C	N	O	S	0	0	0
			2933	1888	493	531	21			
1	H	362	Total	C	N	O	S	0	0	0
			2933	1888	493	531	21			
1	I	363	Total	C	N	O	S	0	0	0
			2940	1892	494	533	21			
1	J	362	Total	C	N	O	S	0	0	0
			2933	1888	493	531	21			
1	K	362	Total	C	N	O	S	0	0	0
			2933	1888	493	531	21			
1	L	363	Total	C	N	O	S	0	0	0
			2940	1892	494	533	21			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	258	GLY	-	CLONING ARTIFACT	UNP Q9NP68
A	259	SER	-	CLONING ARTIFACT	UNP Q9NP68
B	258	GLY	-	CLONING ARTIFACT	UNP Q9NP68
B	259	SER	-	CLONING ARTIFACT	UNP Q9NP68
C	258	GLY	-	CLONING ARTIFACT	UNP Q9NP68

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Chain	Residue	Modelled	Actual	Comment	Reference
C	259	SER	-	CLONING ARTIFACT	UNP Q9NP68
D	258	GLY	-	CLONING ARTIFACT	UNP Q9NP68
D	259	SER	-	CLONING ARTIFACT	UNP Q9NP68
E	258	GLY	-	CLONING ARTIFACT	UNP Q9NP68
E	259	SER	-	CLONING ARTIFACT	UNP Q9NP68
F	258	GLY	-	CLONING ARTIFACT	UNP Q9NP68
F	259	SER	-	CLONING ARTIFACT	UNP Q9NP68
G	258	GLY	-	CLONING ARTIFACT	UNP Q9NP68
G	259	SER	-	CLONING ARTIFACT	UNP Q9NP68
H	258	GLY	-	CLONING ARTIFACT	UNP Q9NP68
H	259	SER	-	CLONING ARTIFACT	UNP Q9NP68
I	258	GLY	-	CLONING ARTIFACT	UNP Q9NP68
I	259	SER	-	CLONING ARTIFACT	UNP Q9NP68
J	258	GLY	-	CLONING ARTIFACT	UNP Q9NP68
J	259	SER	-	CLONING ARTIFACT	UNP Q9NP68
K	258	GLY	-	CLONING ARTIFACT	UNP Q9NP68
K	259	SER	-	CLONING ARTIFACT	UNP Q9NP68
L	258	GLY	-	CLONING ARTIFACT	UNP Q9NP68
L	259	SER	-	CLONING ARTIFACT	UNP Q9NP68

- Molecule 2 is a protein called Cellular tumor antigen p53.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	M	199	Total	C	N	O	S	0	0	0
			1556	959	286	295	16			
2	N	199	Total	C	N	O	S	0	0	0
			1556	959	286	295	16			
2	O	199	Total	C	N	O	S	0	0	0
			1556	959	286	295	16			
2	P	199	Total	C	N	O	S	0	0	0
			1556	959	286	295	16			
2	Q	199	Total	C	N	O	S	0	0	0
			1556	959	286	295	16			
2	R	199	Total	C	N	O	S	0	0	0
			1556	959	286	295	16			
2	S	199	Total	C	N	O	S	0	0	0
			1556	959	286	295	16			
2	T	199	Total	C	N	O	S	0	0	0
			1556	959	286	295	16			
2	U	199	Total	C	N	O	S	0	0	0
			1556	959	286	295	16			
2	V	199	Total	C	N	O	S	0	0	0
			1556	959	286	295	16			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	W	199	Total	C	N	O	S	0	0	0
			1556	959	286	295	16			
2	X	199	Total	C	N	O	S	0	0	0
			1556	959	286	295	16			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
M	90	GLY	-	CLONING ARTIFACT	UNP Q9NP68
M	91	SER	-	CLONING ARTIFACT	UNP Q9NP68
N	90	GLY	-	CLONING ARTIFACT	UNP Q9NP68
N	91	SER	-	CLONING ARTIFACT	UNP Q9NP68
O	90	GLY	-	CLONING ARTIFACT	UNP Q9NP68
O	91	SER	-	CLONING ARTIFACT	UNP Q9NP68
P	90	GLY	-	CLONING ARTIFACT	UNP Q9NP68
P	91	SER	-	CLONING ARTIFACT	UNP Q9NP68
Q	90	GLY	-	CLONING ARTIFACT	UNP Q9NP68
Q	91	SER	-	CLONING ARTIFACT	UNP Q9NP68
R	90	GLY	-	CLONING ARTIFACT	UNP Q9NP68
R	91	SER	-	CLONING ARTIFACT	UNP Q9NP68
S	90	GLY	-	CLONING ARTIFACT	UNP Q9NP68
S	91	SER	-	CLONING ARTIFACT	UNP Q9NP68
T	90	GLY	-	CLONING ARTIFACT	UNP Q9NP68
T	91	SER	-	CLONING ARTIFACT	UNP Q9NP68
U	90	GLY	-	CLONING ARTIFACT	UNP Q9NP68
U	91	SER	-	CLONING ARTIFACT	UNP Q9NP68
V	90	GLY	-	CLONING ARTIFACT	UNP Q9NP68
V	91	SER	-	CLONING ARTIFACT	UNP Q9NP68
W	90	GLY	-	CLONING ARTIFACT	UNP Q9NP68
W	91	SER	-	CLONING ARTIFACT	UNP Q9NP68
X	90	GLY	-	CLONING ARTIFACT	UNP Q9NP68
X	91	SER	-	CLONING ARTIFACT	UNP Q9NP68

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

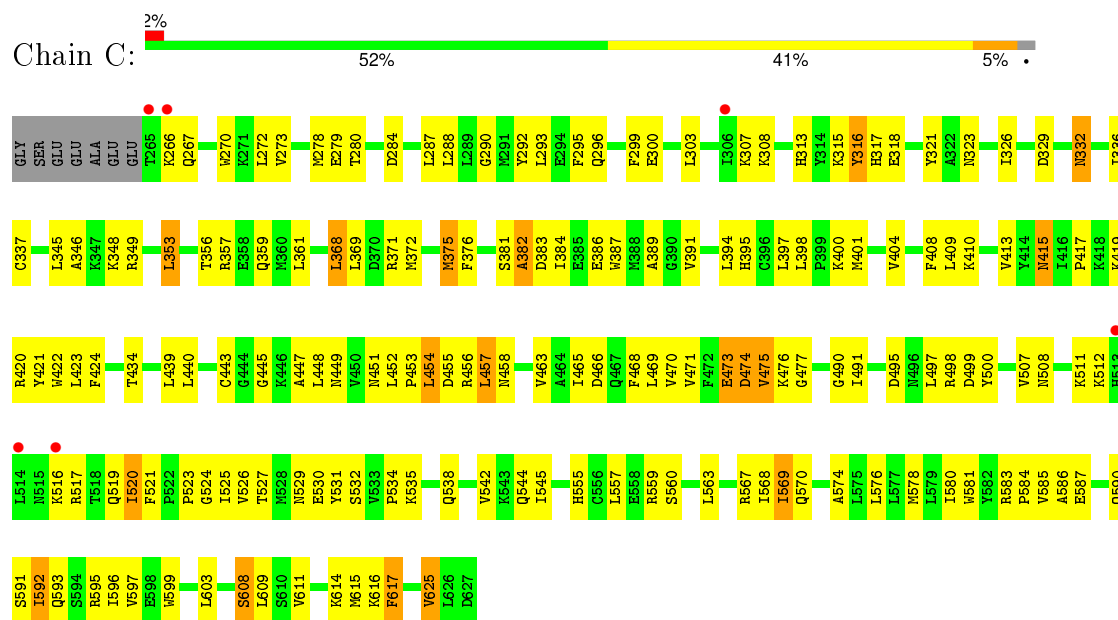
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	P	1	Total	Zn	0	0
			1	1		
3	K	1	Total	Zn	0	0
			1	1		
3	B	1	Total	Zn	0	0
			1	1		

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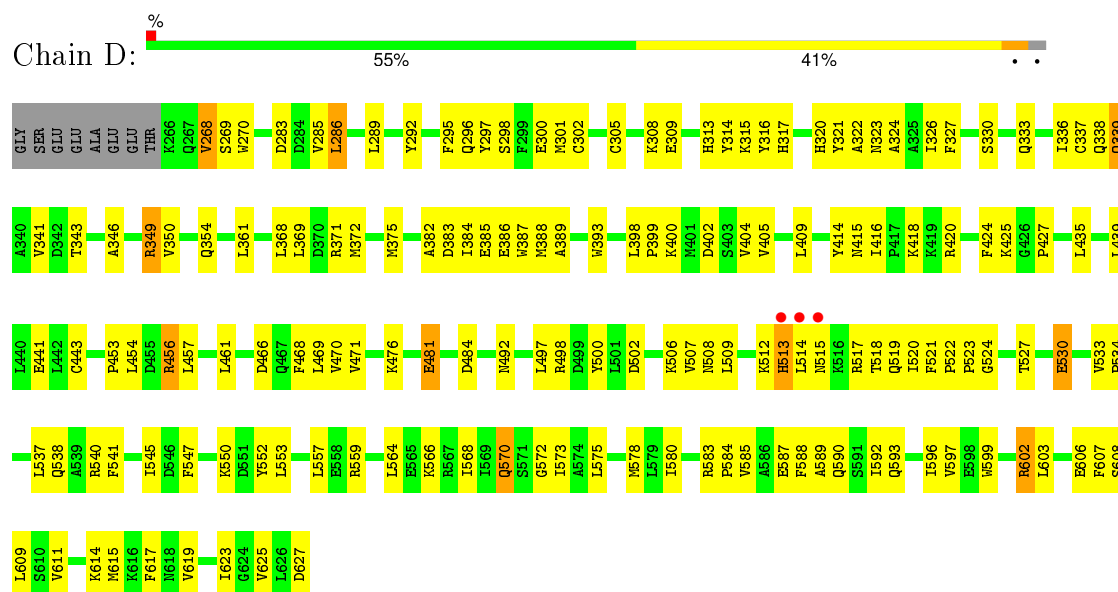
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	W	1	Total 1	Zn 1	0	0
3	N	1	Total 1	Zn 1	0	0
3	X	1	Total 1	Zn 1	0	0
3	S	1	Total 1	Zn 1	0	0
3	J	1	Total 1	Zn 1	0	0
3	E	1	Total 1	Zn 1	0	0
3	V	1	Total 1	Zn 1	0	0
3	A	1	Total 1	Zn 1	0	0
3	R	1	Total 1	Zn 1	0	0
3	M	1	Total 1	Zn 1	0	0
3	D	1	Total 1	Zn 1	0	0
3	I	1	Total 1	Zn 1	0	0
3	U	1	Total 1	Zn 1	0	0
3	L	1	Total 1	Zn 1	0	0
3	G	1	Total 1	Zn 1	0	0
3	Q	1	Total 1	Zn 1	0	0
3	H	1	Total 1	Zn 1	0	0
3	C	1	Total 1	Zn 1	0	0
3	T	1	Total 1	Zn 1	0	0
3	O	1	Total 1	Zn 1	0	0
3	F	1	Total 1	Zn 1	0	0

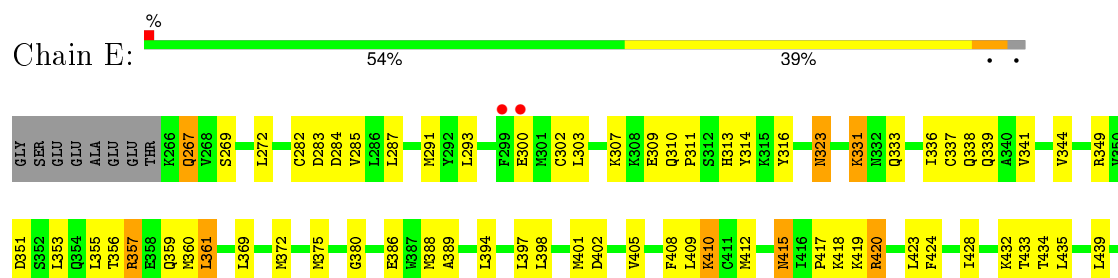
- Molecule 1: Large T antigen

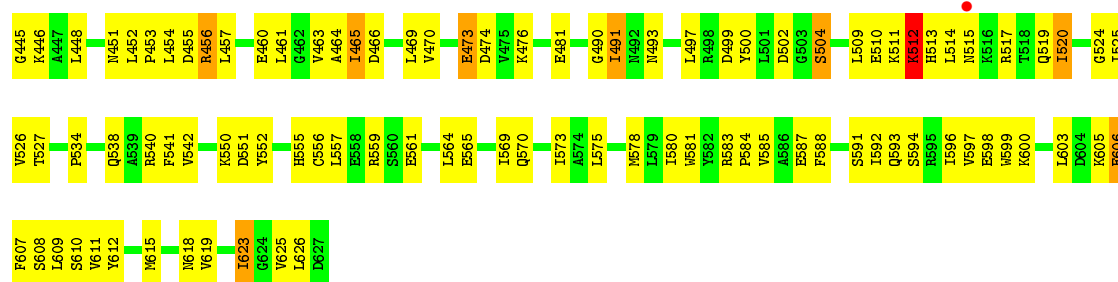


- Molecule 1: Large T antigen

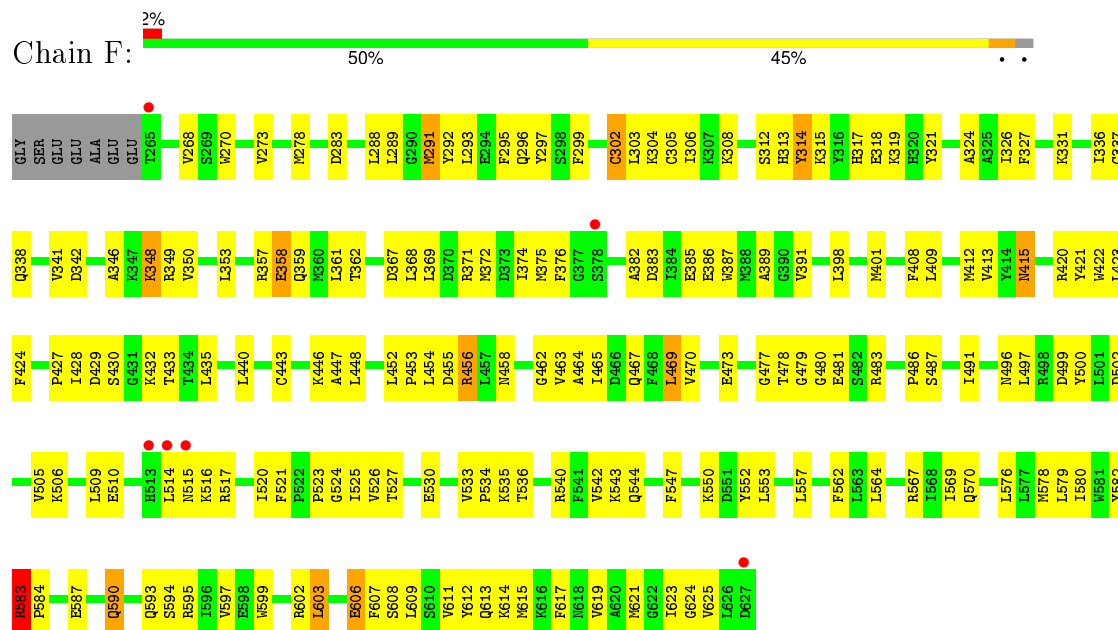


- Molecule 1: Large T antigen

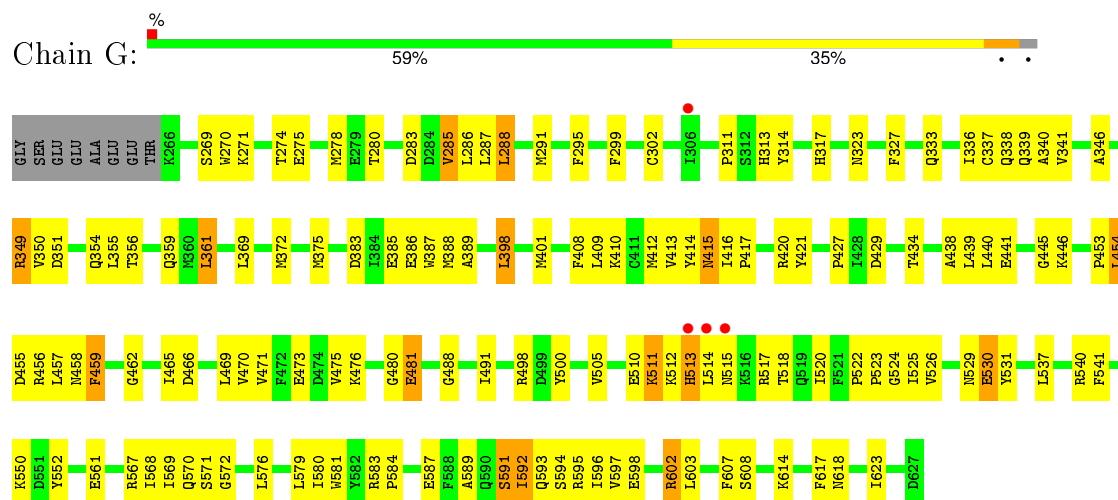




• Molecule 1: Large T antigen

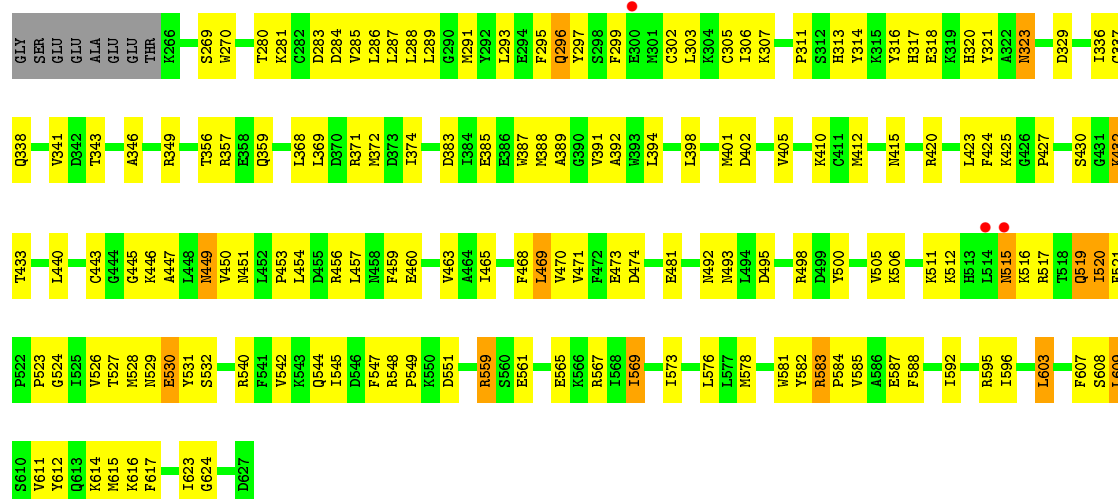


• Molecule 1: Large T antigen

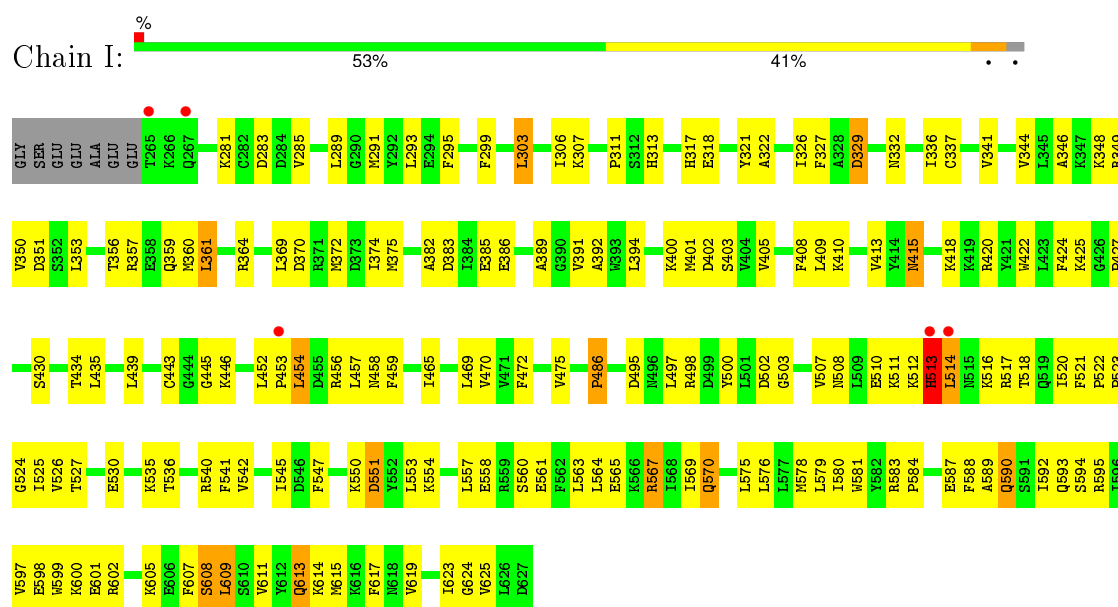


• Molecule 1: Large T antigen





• Molecule 1: Large T antigen

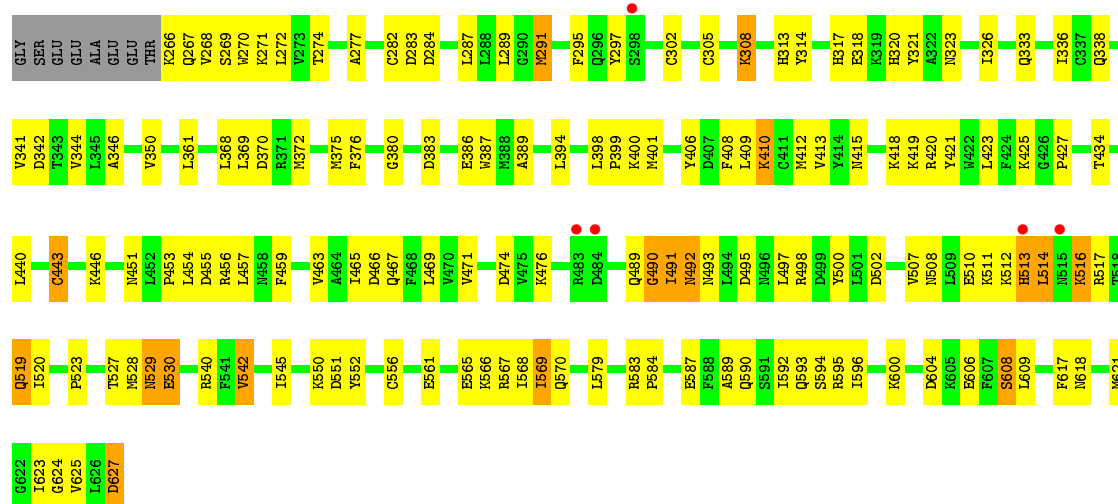


• Molecule 1: Large T antigen

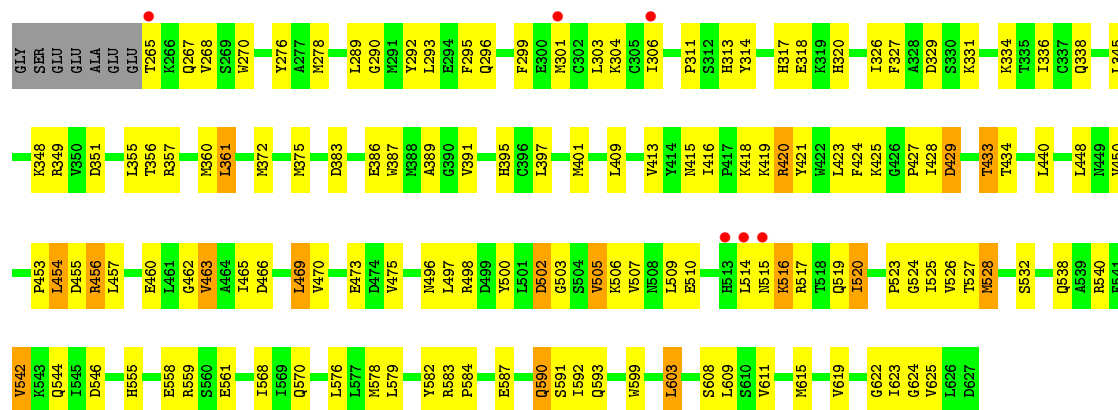




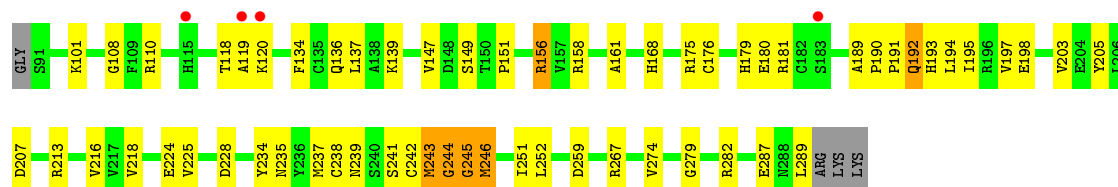
• Molecule 1: Large T antigen



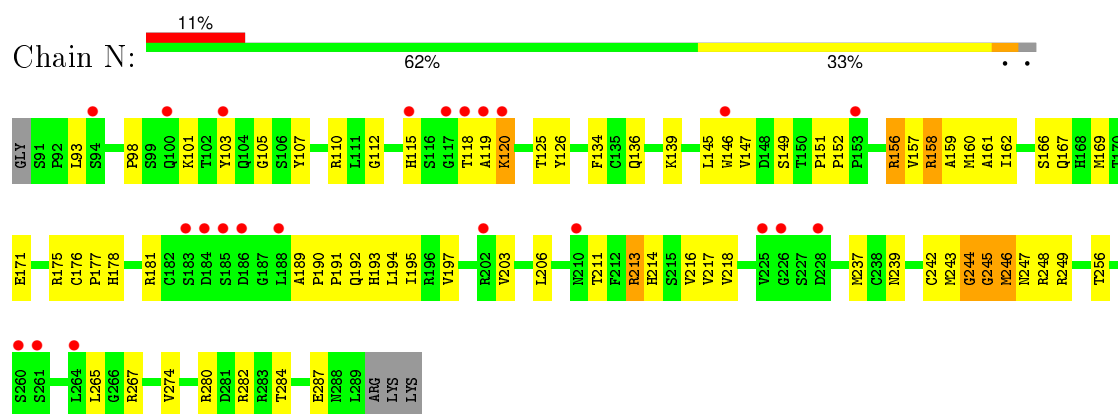
• Molecule 1: Large T antigen



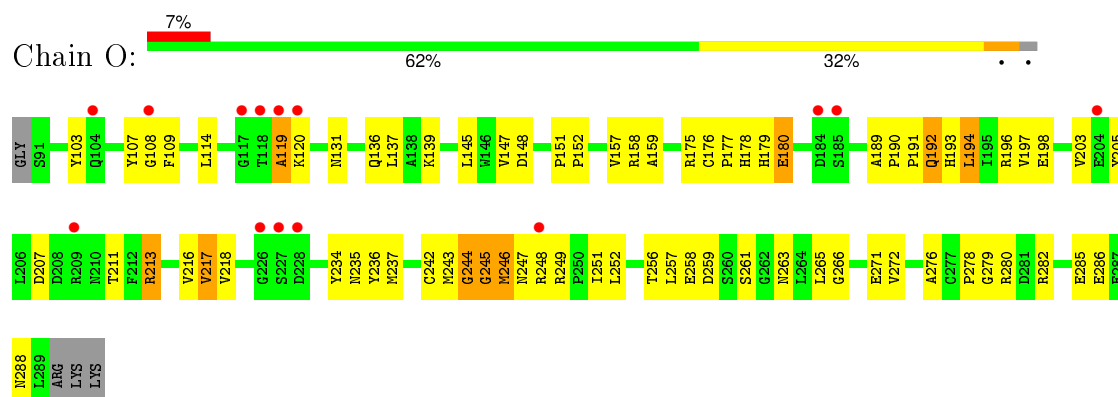
• Molecule 2: Cellular tumor antigen p53



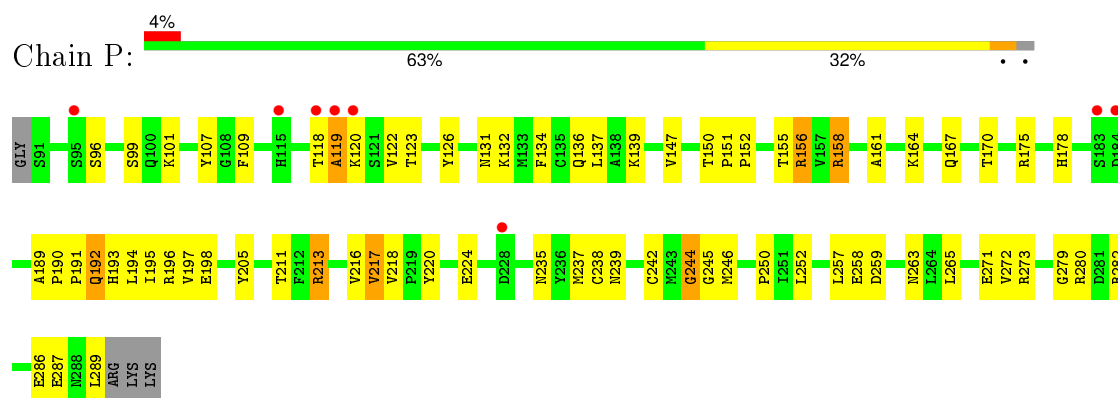
• Molecule 2: Cellular tumor antigen p53



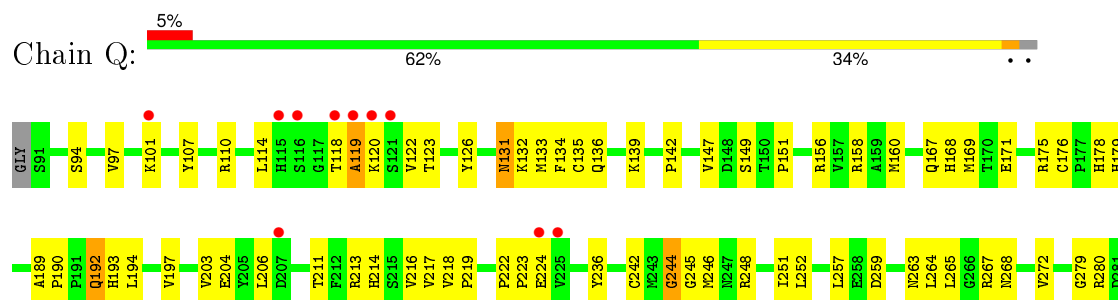
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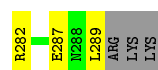


• Molecule 2: Cellular tumor antigen p53

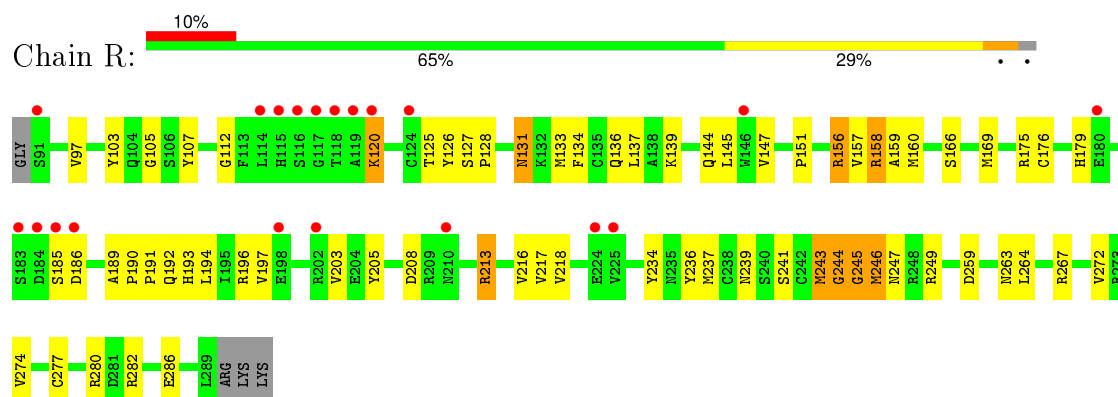


• Molecule 2: Cellular tumor antigen p53

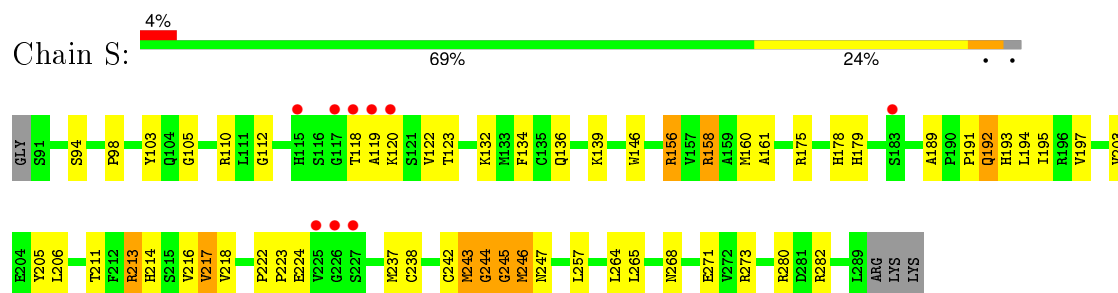




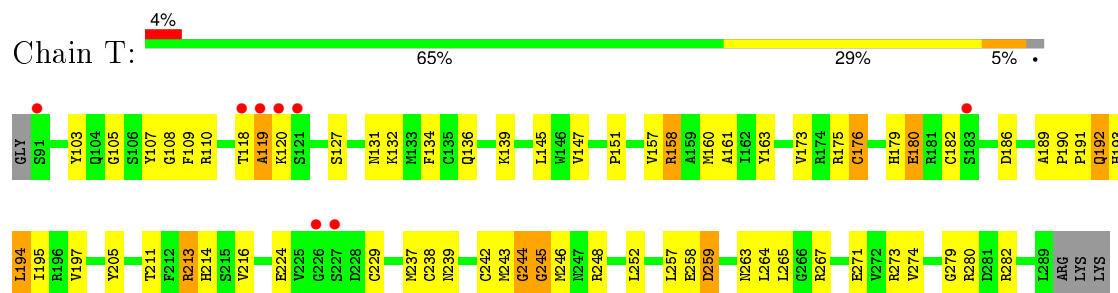
• Molecule 2: Cellular tumor antigen p53



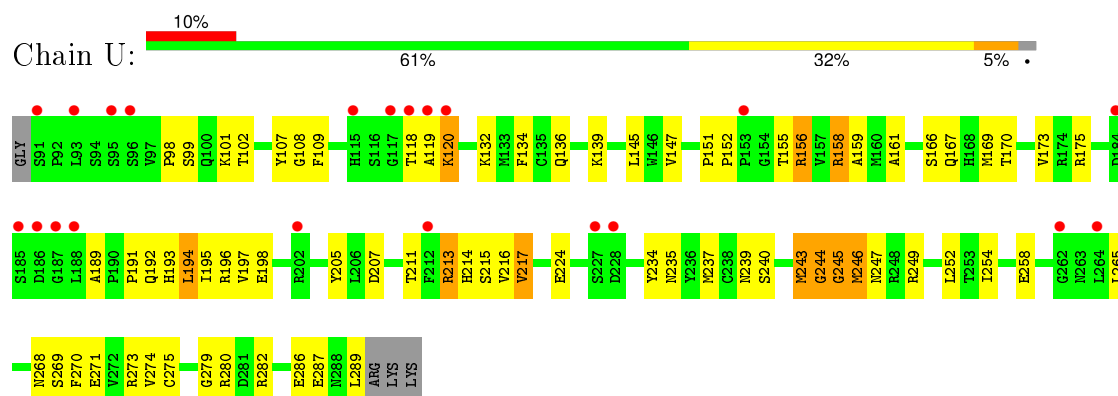
• Molecule 2: Cellular tumor antigen p53



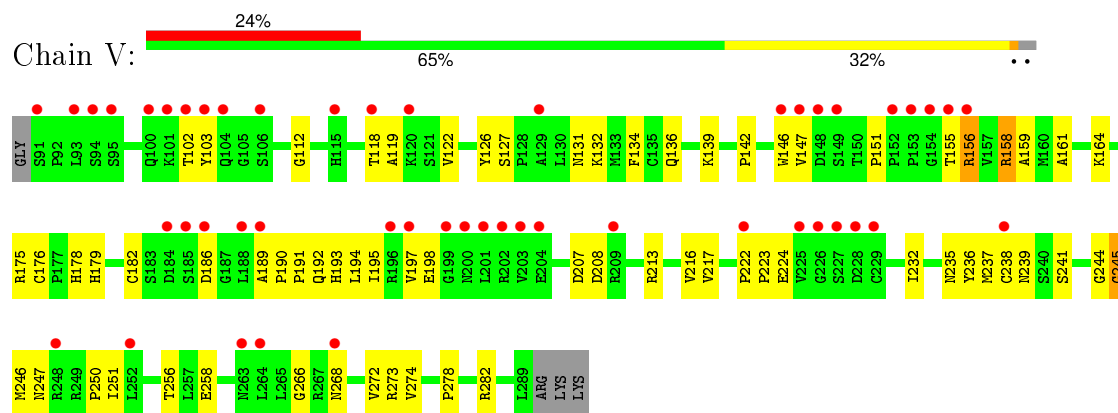
• Molecule 2: Cellular tumor antigen p53



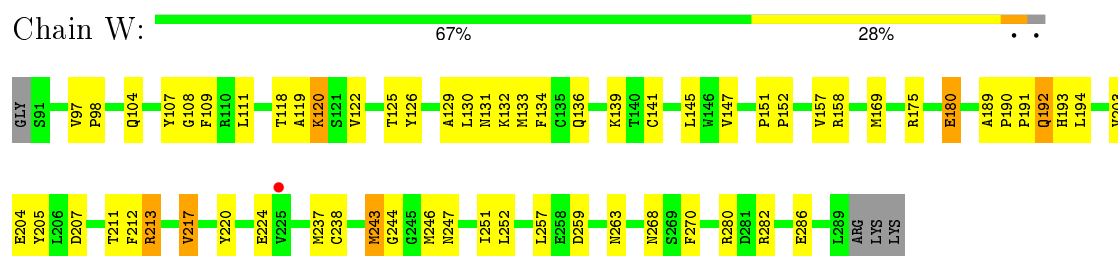
• Molecule 2: Cellular tumor antigen p53



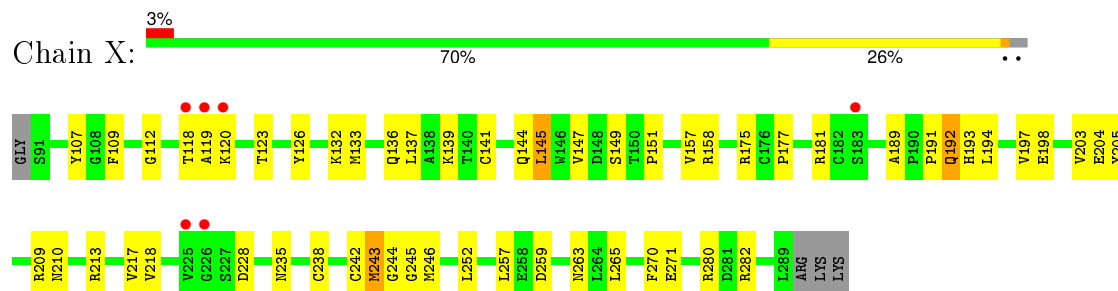
- Molecule 2: Cellular tumor antigen p53



- Molecule 2: Cellular tumor antigen p53



- Molecule 2: Cellular tumor antigen p53



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	127.02Å 182.72Å 262.06Å 90.00° 92.08° 90.00°	Depositor
Resolution (Å)	20.00 – 3.16 19.98 – 3.16	Depositor EDS
% Data completeness (in resolution range)	97.3 (20.00-3.16) 97.0 (19.98-3.16)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.13 (at 3.15Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.261 , 0.308 0.252 , 0.256	Depositor DCC
R_{free} test set	15189 reflections (7.69%)	DCC
Wilson B-factor (Å ²)	81.4	Xtriage
Anisotropy	0.595	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 42.6	EDS
Estimated twinning fraction	0.000 for h,-k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.36$	Xtriage
Outliers	0 of 197864 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	53920	wwPDB-VP
Average B, all atoms (Å ²)	75.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.15% 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, 0.375 respectively for untwinned datasets, and 0.333, 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	0.57	0/2992	0.67	0/4030
1	B	0.55	0/2992	0.63	0/4030
1	C	0.56	0/2999	0.63	0/4040
1	D	0.59	0/2992	0.68	0/4030
1	E	0.61	0/2992	0.67	0/4030
1	F	0.59	0/2999	0.67	0/4040
1	G	0.64	0/2992	0.70	1/4030 (0.0%)
1	H	0.60	0/2992	0.68	0/4030
1	I	0.55	0/2999	0.63	0/4040
1	J	0.52	0/2992	0.61	0/4030
1	K	0.62	0/2992	0.68	0/4030
1	L	0.67	0/2999	0.71	0/4040
2	M	0.55	1/1592 (0.1%)	0.64	0/2160
2	N	0.44	0/1592	0.57	0/2160
2	O	0.47	0/1592	0.61	0/2160
2	P	0.51	0/1592	0.62	0/2160
2	Q	0.47	0/1592	0.60	0/2160
2	R	0.46	1/1592 (0.1%)	0.59	0/2160
2	S	0.52	1/1592 (0.1%)	0.63	0/2160
2	T	0.49	1/1592 (0.1%)	0.61	0/2160
2	U	0.49	1/1592 (0.1%)	0.59	0/2160
2	V	0.46	1/1592 (0.1%)	0.54	0/2160
2	W	0.60	1/1592 (0.1%)	0.65	0/2160
2	X	0.62	1/1592 (0.1%)	0.68	0/2160
All	All	0.56	8/55036 (0.0%)	0.65	1/74320 (0.0%)

The worst 5 of 8 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	M	243	MET	SD-CE	7.60	2.20	1.77
2	U	243	MET	SD-CE	6.29	2.13	1.77

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	X	243	MET	SD-CE	6.17	2.12	1.77
2	T	176	CYS	CB-SG	-5.89	1.72	1.81
2	W	243	MET	SD-CE	5.65	2.09	1.77

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	349	ARG	NE-CZ-NH1	-5.27	117.66	120.30

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	2933	0	2982	129	0
1	B	2933	0	2982	163	0
1	C	2940	0	2989	177	0
1	D	2933	0	2982	144	0
1	E	2933	0	2982	142	0
1	F	2940	0	2989	162	0
1	G	2933	0	2982	129	0
1	H	2933	0	2982	144	0
1	I	2940	0	2989	150	0
1	J	2933	0	2982	152	0
1	K	2933	0	2982	130	0
1	L	2940	0	2989	133	0
2	M	1556	0	1514	40	0
2	N	1556	0	1515	59	0
2	O	1556	0	1513	65	0
2	P	1556	0	1513	55	0
2	Q	1556	0	1513	48	0
2	R	1556	0	1513	52	0
2	S	1556	0	1513	54	0
2	T	1556	0	1513	56	0
2	U	1556	0	1514	64	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	V	1556	0	1515	40	0
2	W	1556	0	1513	41	0
2	X	1556	0	1513	43	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
3	E	1	0	0	0	0
3	F	1	0	0	0	0
3	G	1	0	0	0	0
3	H	1	0	0	0	0
3	I	1	0	0	0	0
3	J	1	0	0	0	0
3	K	1	0	0	0	0
3	L	1	0	0	0	0
3	M	1	0	0	0	0
3	N	1	0	0	0	0
3	O	1	0	0	0	0
3	P	1	0	0	0	0
3	Q	1	0	0	0	0
3	R	1	0	0	0	0
3	S	1	0	0	0	0
3	T	1	0	0	0	0
3	U	1	0	0	0	0
3	V	1	0	0	0	0
3	W	1	0	0	0	0
3	X	1	0	0	0	0
All	All	53920	0	53974	2203	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 20.

The worst 5 of 2203 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:G:291:MET:SD	1:G:291:MET:CE	2.04	1.46
2:T:243:MET:CE	2:T:243:MET:SD	2.04	1.46
2:N:243:MET:SD	2:N:243:MET:CE	2.03	1.45
2:R:243:MET:CE	2:R:243:MET:SD	2.06	1.43
2:W:243:MET:CE	2:W:243:MET:SD	2.09	1.41

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	360/370 (97%)	320 (89%)	35 (10%)	5 (1%)	14	55
1	B	360/370 (97%)	310 (86%)	44 (12%)	6 (2%)	11	50
1	C	361/370 (98%)	306 (85%)	44 (12%)	11 (3%)	5	34
1	D	360/370 (97%)	312 (87%)	43 (12%)	5 (1%)	14	55
1	E	360/370 (97%)	316 (88%)	38 (11%)	6 (2%)	11	50
1	F	361/370 (98%)	309 (86%)	46 (13%)	6 (2%)	11	50
1	G	360/370 (97%)	326 (91%)	29 (8%)	5 (1%)	14	55
1	H	360/370 (97%)	317 (88%)	39 (11%)	4 (1%)	17	61
1	I	361/370 (98%)	321 (89%)	36 (10%)	4 (1%)	17	61
1	J	360/370 (97%)	319 (89%)	37 (10%)	4 (1%)	17	61
1	K	360/370 (97%)	329 (91%)	26 (7%)	5 (1%)	14	55
1	L	361/370 (98%)	330 (91%)	29 (8%)	2 (1%)	30	74
2	M	197/203 (97%)	174 (88%)	19 (10%)	4 (2%)	9	46
2	N	197/203 (97%)	175 (89%)	17 (9%)	5 (2%)	7	39
2	O	197/203 (97%)	177 (90%)	15 (8%)	5 (2%)	7	39
2	P	197/203 (97%)	174 (88%)	20 (10%)	3 (2%)	13	53
2	Q	197/203 (97%)	172 (87%)	22 (11%)	3 (2%)	13	53
2	R	197/203 (97%)	172 (87%)	21 (11%)	4 (2%)	9	46
2	S	197/203 (97%)	173 (88%)	20 (10%)	4 (2%)	9	46
2	T	197/203 (97%)	178 (90%)	15 (8%)	4 (2%)	9	46
2	U	197/203 (97%)	172 (87%)	22 (11%)	3 (2%)	13	53
2	V	197/203 (97%)	180 (91%)	14 (7%)	3 (2%)	13	53

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	W	197/203 (97%)	180 (91%)	16 (8%)	1 (0%)	34	76
2	X	197/203 (97%)	176 (89%)	19 (10%)	2 (1%)	19	63
All	All	6688/6876 (97%)	5918 (88%)	666 (10%)	104 (2%)	12	52

5 of 104 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	473	GLU
1	C	475	VAL
1	D	530	GLU
1	E	514	LEU
1	H	530	GLU

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	323/329 (98%)	305 (94%)	18 (6%)	26	66
1	B	323/329 (98%)	297 (92%)	26 (8%)	15	50
1	C	324/329 (98%)	308 (95%)	16 (5%)	31	71
1	D	323/329 (98%)	307 (95%)	16 (5%)	30	70
1	E	323/329 (98%)	298 (92%)	25 (8%)	16	52
1	F	324/329 (98%)	307 (95%)	17 (5%)	29	68
1	G	323/329 (98%)	307 (95%)	16 (5%)	30	70
1	H	323/329 (98%)	305 (94%)	18 (6%)	26	66
1	I	324/329 (98%)	305 (94%)	19 (6%)	24	63
1	J	323/329 (98%)	307 (95%)	16 (5%)	30	70
1	K	323/329 (98%)	301 (93%)	22 (7%)	20	57
1	L	324/329 (98%)	300 (93%)	24 (7%)	17	54
2	M	179/182 (98%)	168 (94%)	11 (6%)	23	62
2	N	179/182 (98%)	172 (96%)	7 (4%)	39	76

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	O	179/182 (98%)	171 (96%)	8 (4%)	34	73
2	P	179/182 (98%)	169 (94%)	10 (6%)	26	66
2	Q	179/182 (98%)	171 (96%)	8 (4%)	34	73
2	R	179/182 (98%)	173 (97%)	6 (3%)	44	80
2	S	179/182 (98%)	172 (96%)	7 (4%)	39	76
2	T	179/182 (98%)	171 (96%)	8 (4%)	34	73
2	U	179/182 (98%)	168 (94%)	11 (6%)	23	62
2	V	179/182 (98%)	174 (97%)	5 (3%)	51	83
2	W	179/182 (98%)	171 (96%)	8 (4%)	34	73
2	X	179/182 (98%)	171 (96%)	8 (4%)	34	73
All	All	6028/6132 (98%)	5698 (94%)	330 (6%)	27	66

5 of 330 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	H	559	ARG
1	J	562	PHE
2	U	194	LEU
1	I	291	MET
1	I	602	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 203 such sidechains are listed below:

Mol	Chain	Res	Type
1	I	449	ASN
1	K	451	ASN
2	V	131	ASN
1	I	529	ASN
1	J	458	ASN

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 24 ligands modelled in this entry, 24 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	362/370 (97%)	-0.31	7 (1%) 70 54	38, 60, 99, 131	0
1	B	362/370 (97%)	-0.29	1 (0%) 94 92	43, 70, 103, 139	0
1	C	363/370 (98%)	-0.15	6 (1%) 73 59	40, 74, 110, 136	0
1	D	362/370 (97%)	-0.29	3 (0%) 87 79	39, 62, 102, 148	0
1	E	362/370 (97%)	-0.33	3 (0%) 87 79	35, 61, 97, 124	0
1	F	363/370 (98%)	-0.22	6 (1%) 73 59	36, 66, 110, 133	0
1	G	362/370 (97%)	-0.45	4 (1%) 82 71	28, 52, 91, 129	0
1	H	362/370 (97%)	-0.34	3 (0%) 87 79	33, 57, 98, 133	0
1	I	363/370 (98%)	-0.25	5 (1%) 78 64	41, 68, 103, 135	0
1	J	362/370 (97%)	-0.21	6 (1%) 73 59	44, 74, 106, 123	0
1	K	362/370 (97%)	-0.39	5 (1%) 78 64	32, 56, 91, 127	0
1	L	363/370 (98%)	-0.42	6 (1%) 73 59	28, 48, 97, 134	0
2	M	199/203 (98%)	-0.01	4 (2%) 68 52	50, 75, 120, 137	0
2	N	199/203 (98%)	0.70	23 (11%) 6 3	80, 110, 144, 151	0
2	O	199/203 (98%)	0.39	14 (7%) 19 10	67, 96, 134, 147	0
2	P	199/203 (98%)	0.15	8 (4%) 42 26	55, 87, 131, 144	0
2	Q	199/203 (98%)	0.18	10 (5%) 32 17	64, 94, 126, 137	0
2	R	199/203 (98%)	0.51	20 (10%) 9 4	69, 104, 142, 152	0
2	S	199/203 (98%)	0.21	9 (4%) 37 21	53, 88, 129, 143	0
2	T	199/203 (98%)	0.26	8 (4%) 42 26	60, 91, 128, 138	0
2	U	199/203 (98%)	0.65	21 (10%) 8 4	72, 103, 138, 147	0
2	V	199/203 (98%)	1.17	49 (24%) 1 1	85, 123, 147, 154	0
2	W	199/203 (98%)	-0.11	1 (0%) 91 87	41, 70, 109, 123	0
2	X	199/203 (98%)	-0.14	6 (3%) 54 37	35, 65, 112, 136	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å ²)	Q<0.9
All	All	6736/6876 (97%)	-0.08	228 (3%)	49	32	28, 71, 126, 154	0

The worst 5 of 228 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	N	119	ALA	8.9
2	S	119	ALA	8.6
2	P	120	LYS	8.2
2	S	118	THR	7.6
1	C	513	HIS	7.5

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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
3	ZN	C	6	1/1	0.96	0.15	-0.66	79,79,79,79	0
3	ZN	B	2	1/1	0.99	0.12	-0.81	71,71,71,71	0
3	ZN	K	11	1/1	0.98	0.09	-0.98	59,59,59,59	0
3	ZN	E	4	1/1	0.95	0.13	-1.06	57,57,57,57	0
3	ZN	A	1	1/1	0.95	0.12	-1.06	66,66,66,66	0
3	ZN	H	9	1/1	0.98	0.13	-1.07	60,60,60,60	0
3	ZN	D	3	1/1	0.95	0.13	-1.09	84,84,84,84	0
3	ZN	N	14	1/1	0.98	0.13	-1.12	81,81,81,81	0
3	ZN	I	8	1/1	0.98	0.13	-1.18	60,60,60,60	0
3	ZN	G	7	1/1	0.98	0.09	-1.27	66,66,66,66	0
3	ZN	V	22	1/1	0.94	0.15	-1.27	85,85,85,85	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	ZN	U	21	1/1	0.98	0.13	-1.29	78,78,78,78	0
3	ZN	R	18	1/1	0.98	0.12	-1.33	74,74,74,74	0
3	ZN	X	24	1/1	0.99	0.12	-1.39	38,38,38,38	0
3	ZN	S	19	1/1	0.99	0.11	-1.73	45,45,45,45	0
3	ZN	W	23	1/1	0.98	0.10	-1.81	45,45,45,45	0
3	ZN	L	12	1/1	0.99	0.13	-1.85	57,57,57,57	0
3	ZN	Q	17	1/1	0.98	0.11	-1.99	44,44,44,44	0
3	ZN	J	10	1/1	0.98	0.10	-2.12	76,76,76,76	0
3	ZN	T	20	1/1	0.99	0.09	-2.14	51,51,51,51	0
3	ZN	M	13	1/1	0.95	0.12	-2.33	55,55,55,55	0
3	ZN	O	15	1/1	0.99	0.11	-2.50	61,61,61,61	0
3	ZN	P	16	1/1	0.97	0.12	-	58,58,58,58	0
3	ZN	F	5	1/1	0.98	0.12	-	68,68,68,68	0

6.5 Other polymers [i](#)

There are no such residues in this entry.