



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 10, 2023 – 03:53 AM EDT

PDB ID : 7SUV  
Title : APE1 exonuclease substrate complex with 8oxoG opposite A  
Authors : Whitaker, A.W.; Freudenthal, B.D.  
Deposited on : 2021-11-18  
Resolution : 1.99 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](mailto: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>.

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

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtrriage (Phenix) : 1.13  
EDS : 2.35.1  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.35.1

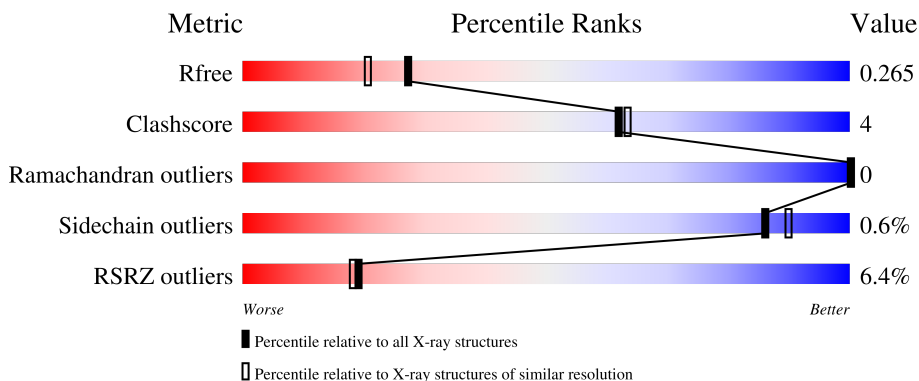
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.99 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))
$R_{free}$	130704	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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  $\geq 3$ , 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	C	11	
2	D	10	
3	E	21	
4	A	276	
4	B	276	

## 2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 5553 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 DNA chain called DNA (5'-D(P\*TP\*CP\*GP\*AP\*CP\*GP\*GP\*AP\*TP\*CP\*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
1	C	11	225	106	41	67	11	0	0	0

- Molecule 2 is a DNA chain called DNA (5'-D(\*GP\*CP\*TP\*GP\*AP\*TP\*GP\*CP\*GP\*(80G))-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
2	D	10	207	98	40	60	9	0	0	0

- Molecule 3 is a DNA chain called DNA (5'-D(\*GP\*GP\*AP\*TP\*CP\*CP\*GP\*TP\*CP\*GP\*AP\*AP\*CP\*GP\*CP\*AP\*TP\*CP\*AP\*GP\*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
3	E	21	427	203	82	122	20	0	0	0

- Molecule 4 is a protein called DNA-(apurinic or apyrimidinic site) lyase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
4	A	276	2207	1410	383	406	8	0	3	0
4	B	276	2224	1419	389	408	8	0	4	0

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	96	GLN	GLU	conflict	UNP P27695
A	138	ALA	CYS	conflict	UNP P27695
A	210	ASN	ASP	conflict	UNP P27695

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Chain	Residue	Modelled	Actual	Comment	Reference
B	96	GLN	GLU	conflict	UNP P27695
B	138	ALA	CYS	conflict	UNP P27695
B	210	ASN	ASP	conflict	UNP P27695

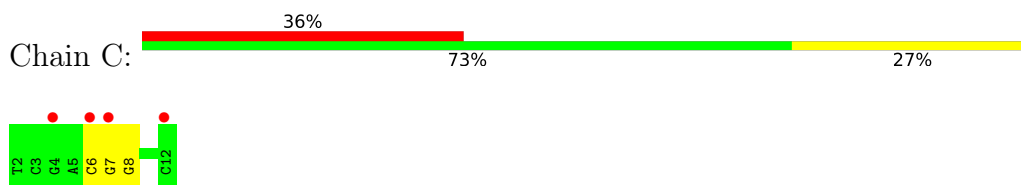
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	C	1	Total O 1 1	0	0
5	D	10	Total O 10 10	0	0
5	E	12	Total O 12 12	0	0
5	A	101	Total O 101 101	0	0
5	B	139	Total O 139 139	0	0

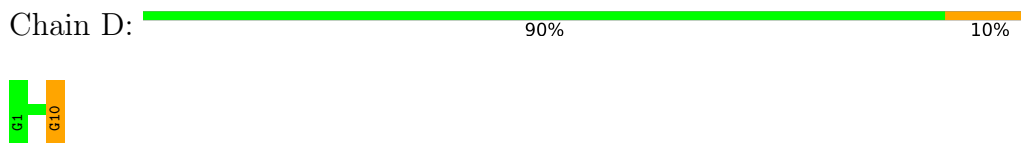
### 3 Residue-property plots [i](#)

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.

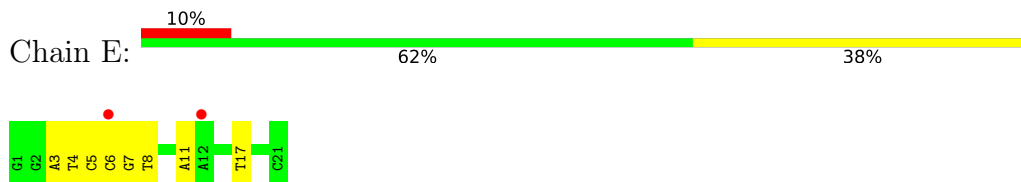
- Molecule 1: DNA (5'-D(P\*TP\*CP\*GP\*AP\*CP\*GP\*GP\*AP\*TP\*CP\*C)-3')



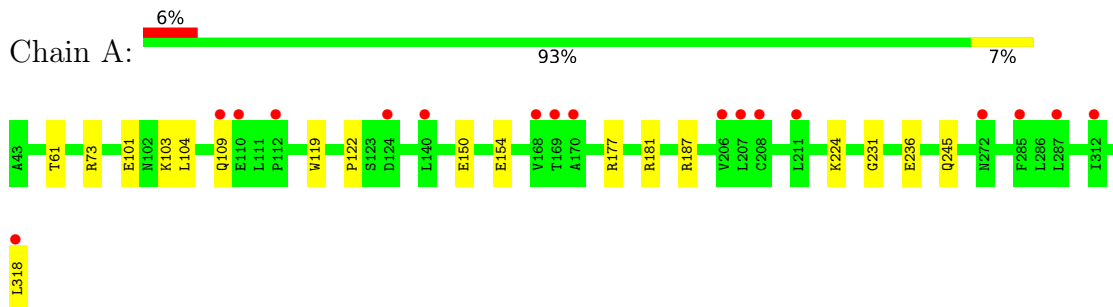
- Molecule 2: DNA (5'-D(\*GP\*CP\*TP\*GP\*AP\*TP\*GP\*CP\*GP\*(8OG))-3')



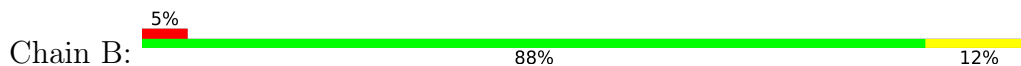
- Molecule 3: DNA (5'-D(\*GP\*GP\*AP\*TP\*CP\*CP\*GP\*TP\*CP\*GP\*AP\*AP\*CP\*GP\*CP\*AP\*TP\*CP\*AP\*GP\*C)-3')

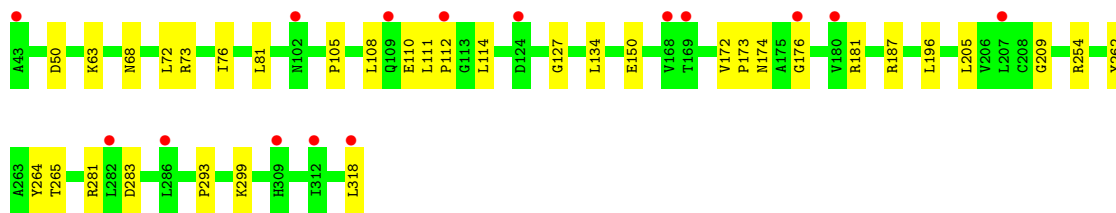


- Molecule 4: DNA-(apurinic or apyrimidinic site) lyase



- Molecule 4: DNA-(apurinic or apyrimidinic site) lyase





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	71.11Å 66.49Å 91.98Å 90.00° 110.69° 90.00°	Depositor
Resolution (Å)	24.96 – 1.99 24.96 – 1.99	Depositor EDS
% Data completeness (in resolution range)	89.1 (24.96-1.99) 73.4 (24.96-1.99)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.78 (at 1.99Å)	Xtrriage
Refinement program	PHENIX 1.19.2_4158	Depositor
R, $R_{free}$	0.226 , 0.265 0.226 , 0.265	Depositor DCC
$R_{free}$ test set	1871 reflections (3.72%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	30.0	Xtrriage
Anisotropy	0.514	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 36.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.023 for h,-k,-h-l	Xtrriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	5553	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	45.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.21% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: 8OG

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	C	0.74	0/251	0.84	0/383
2	D	0.95	0/206	0.97	0/317
3	E	0.93	0/479	0.99	0/737
4	A	0.40	0/2271	0.61	0/3079
4	B	0.39	0/2285	0.63	0/3096
All	All	0.51	0/5492	0.69	0/7612

There are no bond length outliers.

There are no bond angle outliers.

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	C	225	0	124	3	0
2	D	207	0	114	1	0
3	E	427	0	236	11	0
4	A	2207	0	2187	13	0
4	B	2224	0	2204	20	0
5	A	101	0	0	1	0
5	B	139	0	0	1	0
5	C	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	D	10	0	0	0	0
5	E	12	0	0	0	0
All	All	5553	0	4865	43	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 4.

All (43) 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)
3:E:17:DT:H5'	4:B:127:GLY:HA3	1.72	0.71
1:C:6:DC:H42	3:E:6:DC:H42	1.40	0.70
4:A:245:GLN:HG3	4:B:293:PRO:HG2	1.76	0.66
4:A:101:GLU:HA	4:A:104:LEU:HD13	1.78	0.64
3:E:5:DC:H6	3:E:5:DC:H5''	1.64	0.62
4:A:177:ARG:NH1	5:A:401:HOH:O	2.26	0.62
4:A:61:THR:HB	4:A:318:LEU:H	1.66	0.60
3:E:5:DC:H2'	3:E:6:DC:C6	2.39	0.58
3:E:6:DC:H2'	3:E:7:DG:C8	2.40	0.56
1:C:6:DC:N4	3:E:6:DC:H42	2.01	0.56
4:B:76:ILE:HD12	4:B:105:PRO:HG2	1.88	0.55
2:D:10:8OG:OP2	4:B:174:ASN:HB2	2.08	0.53
4:A:154:GLU:OE2	4:A:181:ARG:NH2	2.42	0.52
4:A:73:ARG:NH1	4:A:103:LYS:HD2	2.25	0.51
4:B:174:ASN:OD1	4:B:176:GLY:N	2.39	0.50
4:A:104:LEU:HD21	4:A:119:TRP:CE2	2.47	0.49
3:E:3:DA:H2''	3:E:4:DT:H5'	1.95	0.48
4:B:110:GLU:O	4:B:112:PRO:HD3	2.14	0.47
3:E:7:DG:H2''	3:E:8:DT:H5''	1.97	0.47
4:A:150:GLU:CD	4:A:187:ARG:HH21	2.18	0.47
4:B:262:TYR:HA	4:B:264:TYR:CZ	2.49	0.47
3:E:4:DT:H2''	3:E:5:DC:H5''	1.97	0.46
4:B:150:GLU:OE1	4:B:187:ARG:NH2	2.49	0.46
4:A:122:PRO:HB3	4:A:154:GLU:HA	1.96	0.45
1:C:7:DG:C6	1:C:8:DG:C6	3.04	0.45
4:B:81:LEU:HD22	4:B:111:LEU:HD11	1.98	0.45
4:A:73:ARG:HH12	4:A:103:LYS:HD2	1.82	0.45
4:B:50:ASP:OD1	4:B:299:LYS:HE2	2.17	0.45
4:B:63:LYS:NZ	5:B:413:HOH:O	2.51	0.44
4:B:172:VAL:HG13	4:B:173:PRO:HD2	1.99	0.44
4:A:318:LEU:HD23	4:A:318:LEU:HA	1.78	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:B:114:LEU:HB3	4:B:134:LEU:HB3	2.00	0.43
4:A:231:GLY:H	4:A:236:GLU:CD	2.23	0.42
4:B:72:LEU:HD21	4:B:108:LEU:HD11	2.02	0.42
4:B:265:THR:HG23	4:B:281:ARG:HG3	2.01	0.41
3:E:5:DC:H2'	3:E:6:DC:H6	1.84	0.41
4:B:254:ARG:HD3	4:B:254:ARG:HA	1.90	0.41
4:B:73:ARG:HG2	4:B:105:PRO:HG3	2.03	0.41
4:B:318:LEU:HA	4:B:318:LEU:HD23	1.82	0.40
3:E:11:DA:H2'	3:E:11:DA:OP2	2.21	0.40
4:A:224:LYS:HB2	4:A:224:LYS:HE2	1.82	0.40
4:B:196:LEU:HD22	4:B:205:LEU:HD21	2.03	0.40
4:B:209:GLY:O	4:B:283:ASP:HA	2.21	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	
4	A	277/276 (100%)	268 (97%)	9 (3%)	0	100	100
4	B	278/276 (101%)	269 (97%)	9 (3%)	0	100	100
All	All	555/552 (100%)	537 (97%)	18 (3%)	0	100	100

There are no Ramachandran outliers to report.

### 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	
4	A	237/235 (101%)	236 (100%)	1 (0%)	91	93
4	B	238/235 (101%)	236 (99%)	2 (1%)	81	86
All	All	475/470 (101%)	472 (99%)	3 (1%)	86	90

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
4	A	109	GLN
4	B	68	ASN
4	B	181	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

1 non-standard protein/DNA/RNA residue is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
2	8OG	D	10	2	22,25,26	2.24	6 (27%)	30,37,40	2.53	9 (30%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '2' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	8OG	D	10	2	-	1/7/21/22	0/3/3/3

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	10	8OG	C5-N7	5.86	1.47	1.37
2	D	10	8OG	C8-N9	4.79	1.49	1.40
2	D	10	8OG	O8-C8	-3.86	1.16	1.23
2	D	10	8OG	C6-N1	-3.10	1.33	1.38
2	D	10	8OG	C8-N7	-2.33	1.33	1.38
2	D	10	8OG	C5-C4	2.28	1.40	1.37

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	10	8OG	C5-C4-N3	-7.87	119.67	127.80
2	D	10	8OG	N9-C4-N3	5.30	131.86	125.81
2	D	10	8OG	C2-N1-C6	-4.59	116.73	125.10
2	D	10	8OG	O4'-C1'-N9	4.34	112.65	108.29
2	D	10	8OG	O6-C6-C5	-3.49	119.23	127.24
2	D	10	8OG	C5-C6-N1	3.04	121.36	112.31
2	D	10	8OG	O8-C8-N7	-2.92	121.35	126.64
2	D	10	8OG	O8-C8-N9	2.80	129.90	125.99
2	D	10	8OG	C2-N3-C4	2.30	116.39	112.30

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	10	8OG	O4'-C4'-C5'-O5'

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	10	8OG	1	0

## 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

There are no ligands in this entry.

## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	C	11/11 (100%)	1.86	4 (36%) 0 0	61, 92, 101, 101	0
2	D	9/10 (90%)	-0.19	0 100 100	40, 43, 48, 50	0
3	E	21/21 (100%)	0.68	2 (9%) 8 7	40, 57, 91, 100	0
4	A	276/276 (100%)	0.53	17 (6%) 20 19	29, 41, 60, 76	0
4	B	276/276 (100%)	0.50	15 (5%) 25 24	29, 39, 56, 73	0
All	All	593/594 (99%)	0.53	38 (6%) 19 18	29, 40, 64, 101	0

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	A	112	PRO	4.2
4	A	109	GLN	3.5
4	A	208	CYS	3.4
4	B	180	VAL	3.3
1	C	6	DC	3.2
1	C	12	DC	3.2
3	E	12	DA	3.1
4	B	43	ALA	3.0
1	C	7	DG	3.0
4	A	169	THR	2.8
4	B	176	GLY	2.8
4	B	207	LEU	2.7
4	A	207	LEU	2.6
4	B	124	ASP	2.6
3	E	6	DC	2.6
4	B	102	ASN	2.5
4	B	282	LEU	2.5
4	B	318	LEU	2.5
4	A	110	GLU	2.4
1	C	4	DG	2.4

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Mol	Chain	Res	Type	RSRZ
4	A	140	LEU	2.4
4	A	318	LEU	2.4
4	A	287	LEU	2.3
4	A	272	ASN	2.3
4	B	112	PRO	2.3
4	A	312	ILE	2.3
4	A	168	VAL	2.3
4	A	124	ASP	2.3
4	B	168	VAL	2.2
4	B	169	THR	2.2
4	A	211	LEU	2.2
4	A	170	ALA	2.2
4	A	206	VAL	2.2
4	B	309	HIS	2.2
4	A	285	PHE	2.2
4	B	312	ILE	2.1
4	B	109	GLN	2.1
4	B	286	LEU	2.1

## 6.2 Non-standard residues in protein, DNA, RNA chains [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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
2	8OG	D	10	23/24	0.80	0.32	41,57,68,72	0

## 6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 6.4 Ligands [i](#)

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.