



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 19, 2016 – 10:13 PM GMT

PDB ID : 5CO8  
Title : Crystal structure of the Holliday junction-resolving enzyme GEN1 (WT) in complex with product DNA and Mg<sup>2+</sup> ion  
Authors : Liu, Y.J.; Freeman, A.D.J.; Declais, A.C.; Wilson, T.J.; Gartner, A.; Lilley, D.M.J.  
Deposited on : 2015-07-20  
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](mailto: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.

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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	:	unknown
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026982
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)	:	rb-20026982

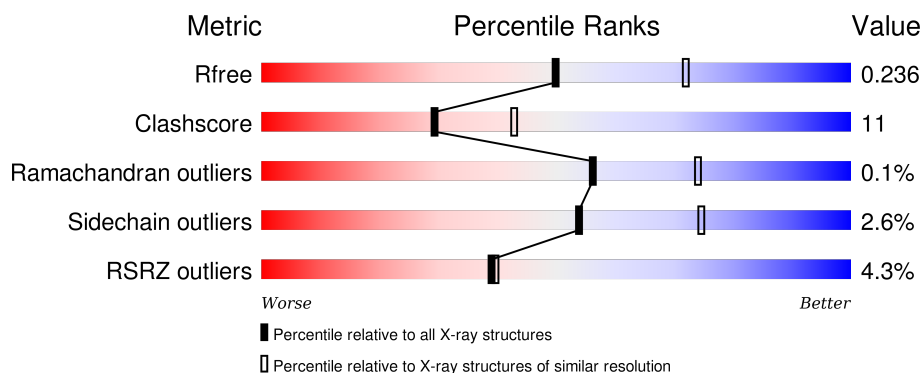
# 1 Overall quality at a glance

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))
$R_{free}$	91344	2919 (2.40-2.40)
Clashscore	102246	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)
RSRZ outliers	91569	2928 (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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	R	31	<div> <div>6%</div> <div>39%</div> <div>61%</div> </div>
2	C	464	<div> <div>4%</div> <div>71%</div> <div>13%</div> <div>•</div> <div>14%</div> </div>
3	X	16	<div> <div>6%</div> <div>63%</div> <div>38%</div> </div>
4	A	464	<div> <div>3%</div> <div>68%</div> <div>15%</div> <div>•</div> <div>15%</div> </div>
5	H	15	<div> <div>27%</div> <div>73%</div> </div>

## 2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 7414 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 (31-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	R	31	Total	C	N	O	P	0	0	1
			602	287	106	179	30			

- Molecule 2 is a protein called Nuclease-like protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	C	397	Total	C	N	O	S	Se	0	0	0
			3109	1984	542	570	6	7			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	X	16	Total	C	N	O	P	0	0	1
			296	138	54	89	15			

- Molecule 4 is a protein called Nuclease-like protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
4	A	393	Total	C	N	O	S	Se	0	0	0
			3078	1962	538	565	8	5			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	H	15	Total	C	N	O	P	0	0	1
			298	139	59	86	14			

- Molecule 6 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	1	Total	Mg	0	0
			1	1		
6	C	1	Total	Mg	0	0
			1	1		

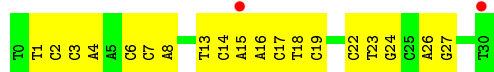
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	C	14	Total	O	0	0
			14	14		
7	X	1	Total	O	0	0
			1	1		
7	A	11	Total	O	0	0
			11	11		
7	H	3	Total	O	0	0
			3	3		

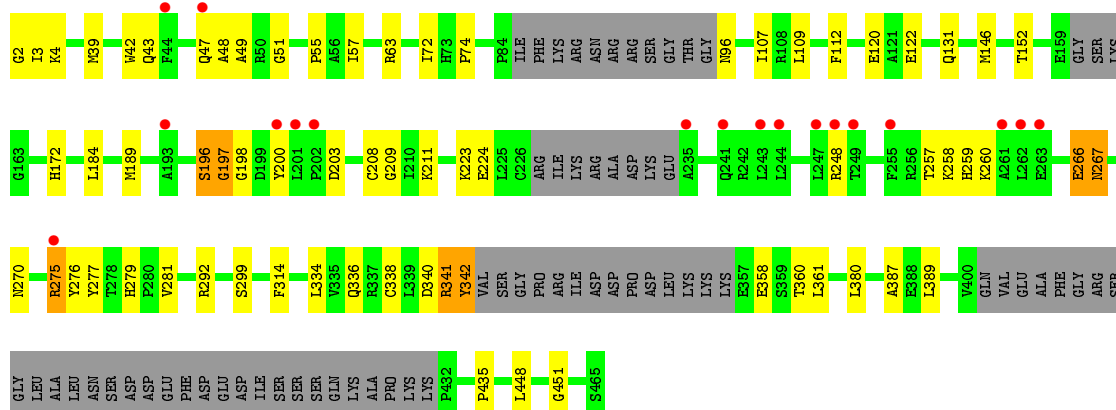
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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 ( $\text{RSRZ} > 2$ ). 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 (31-MER)



- Molecule 2: Nuclease-like protein





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	98.07Å 98.07Å 119.57Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	49.04 – 2.40 69.24 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.2 (49.04-2.40) 99.0 (69.24-2.10)	Depositor EDS
$R_{merge}$	0.14	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.02 (at 2.10Å)	Xtriage
Refinement program	PHENIX	Depositor
R, $R_{free}$	0.239 , 0.255 0.248 , 0.236	Depositor DCC
$R_{free}$ test set	1353 reflections (5.40%)	DCC
Wilson B-factor (Å <sup>2</sup> )	53.6	Xtriage
Anisotropy	0.165	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 52.1	EDS
Estimated twinning fraction	0.036 for -h,-k,l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 39180 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	7414	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.92% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: MG

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	R	0.78	0/672	0.92	1/1032 (0.1%)
2	C	0.24	0/3174	0.46	3/4286 (0.1%)
3	X	0.48	0/331	0.93	0/512
4	A	0.38	1/3142 (0.0%)	0.49	3/4243 (0.1%)
5	H	0.73	1/335 (0.3%)	0.88	0/519
All	All	0.42	2/7654 (0.0%)	0.59	7/10592 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	297	PRO	C-O	14.64	1.52	1.23
5	H	17	DG	O3'-P	-7.71	1.51	1.61

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	297	PRO	O-C-N	-10.05	106.61	122.70
2	C	196	SER	N-CA-C	7.69	131.77	111.00
4	A	297	PRO	CA-C-O	7.08	137.20	120.20
2	C	196	SER	CB-CA-C	-6.67	97.42	110.10
2	C	197	GLY	N-CA-C	-6.21	97.57	113.10
4	A	243	LEU	CA-CB-CG	5.55	128.07	115.30
1	R	4	DA	O5'-P-OP1	-5.24	100.99	105.70

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	R	602	0	337	32	0
2	C	3109	0	3095	65	16
3	X	296	0	158	16	0
4	A	3078	0	3066	60	16
5	H	298	0	158	18	0
6	A	1	0	0	0	0
6	C	1	0	0	0	0
7	A	11	0	0	4	0
7	C	14	0	0	2	0
7	H	3	0	0	0	0
7	X	1	0	0	4	0
All	All	7414	0	6814	155	16

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 11.

All (155) 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)
4:A:261:ALA:HB2	5:H:27:DT:P	1.91	1.10
2:C:341:ARG:NH2	2:C:358:GLU:OE2	1.88	1.07
2:C:203:ASP:CG	2:C:267:ASN:HB3	1.78	1.04
1:R:6:DC:H42	5:H:25:DG:H1	1.06	0.97
2:C:122:GLU:OE1	7:C:602:HOH:O	1.82	0.97
1:R:6:DC:N4	5:H:25:DG:H1	1.66	0.94
1:R:26:DA:N1	3:X:5:DT:N3	2.15	0.94
4:A:261:ALA:HB2	5:H:27:DT:OP1	1.67	0.94
4:A:261:ALA:CB	5:H:27:DT:OP1	2.18	0.92
1:R:24:DG:H1	3:X:7:DC:H42	1.20	0.90
1:R:23:DT:OP1	2:C:211:LYS:N	2.05	0.90
4:A:341:ARG:HH11	4:A:341:ARG:HG2	1.41	0.83
2:C:257:THR:HB	3:X:12:DG:O3'	1.80	0.81
1:R:8:DA:OP1	4:A:209:GLY:HA2	1.81	0.80
4:A:357:GLU:HG3	4:A:358:GLU:H	1.47	0.79
3:X:4:DC:C5	7:X:101:HOH:O	2.35	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:275:ARG:HG2	2:C:279:HIS:HB2	1.64	0.79
2:C:299:SER:O	2:C:336:GLN:NE2	2.18	0.77
4:A:122:GLU:OE2	7:A:602:HOH:O	2.00	0.77
1:R:1:DT:O4	5:H:30:DA:N1	2.20	0.75
1:R:24:DG:H1	3:X:7:DC:N4	1.84	0.74
4:A:140:GLU:OE1	7:A:603:HOH:O	2.05	0.73
1:R:27:DG:N2	3:X:4:DC:O2	2.20	0.73
2:C:203:ASP:OD2	2:C:270:ASN:ND2	2.20	0.73
1:R:24:DG:N2	3:X:7:DC:N3	2.36	0.73
3:X:4:DC:N4	7:X:101:HOH:O	2.07	0.72
2:C:2:GLY:N	2:C:200:TYR:HH	1.88	0.71
2:C:248:ARG:NH1	2:C:259:HIS:O	2.24	0.71
2:C:258:LYS:O	3:X:12:DG:H5''	1.91	0.71
3:X:2:DG:H1'	3:X:3:DA:H5'	1.73	0.70
1:R:6:DC:N3	5:H:25:DG:N2	2.40	0.70
1:R:22:DC:H3'	2:C:211:LYS:HG3	1.74	0.69
2:C:275:ARG:HG3	2:C:279:HIS:ND1	2.09	0.67
4:A:341:ARG:HG2	4:A:341:ARG:NH1	2.03	0.66
4:A:244:LEU:HD13	4:A:264:ILE:HD13	1.78	0.65
2:C:203:ASP:CB	2:C:267:ASN:HB3	2.27	0.65
4:A:319:ARG:CZ	4:A:399:GLU:HB3	2.27	0.64
2:C:387:ALA:HB2	2:C:435:PRO:HB2	1.79	0.64
4:A:261:ALA:CB	5:H:27:DT:P	2.78	0.64
4:A:387:ALA:HB2	4:A:435:PRO:HB2	1.79	0.64
1:R:1:DT:H3	5:H:30:DA:H2	1.46	0.63
4:A:141:ASP:OD1	7:A:604:HOH:O	2.15	0.63
2:C:334:LEU:HB2	2:C:389:LEU:HD21	1.81	0.63
1:R:13:DT:H2''	1:R:14:DC:O5'	1.99	0.62
2:C:223:LYS:HA	2:C:224:GLU:HB3	1.82	0.62
1:R:2:DC:H2''	1:R:3:DC:O5'	1.98	0.62
2:C:358:GLU:HB3	2:C:360:THR:HG22	1.82	0.61
2:C:203:ASP:CG	2:C:267:ASN:CB	2.64	0.61
4:A:258:LYS:HB2	5:H:28:DG:OP1	2.00	0.61
4:A:261:ALA:N	5:H:27:DT:OP1	2.34	0.60
1:R:17:DC:OP1	4:A:63:ARG:NH1	2.35	0.60
1:R:23:DT:OP1	2:C:211:LYS:HB2	2.02	0.59
4:A:146:MSE:HG3	4:A:189:MSE:HE2	1.84	0.58
2:C:275:ARG:CG	2:C:279:HIS:ND1	2.67	0.58
4:A:336:GLN:O	4:A:340:ASP:HB2	2.02	0.58
4:A:54:ASN:HB3	4:A:57:ILE:HG22	1.85	0.58
4:A:463:GLN:C	4:A:465:SER:H	2.08	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:248:ARG:O	4:A:258:LYS:NZ	2.34	0.57
2:C:63:ARG:NH2	7:C:605:HOH:O	2.32	0.56
2:C:43:GLN:HA	2:C:96:ASN:ND2	2.21	0.55
3:X:4:DC:H5	7:X:101:HOH:O	1.79	0.55
1:R:6:DC:H4'	4:A:256:ARG:CZ	2.38	0.54
4:A:340:ASP:C	4:A:342:TYR:H	2.11	0.53
2:C:43:GLN:HA	2:C:96:ASN:HD22	1.73	0.53
2:C:266:GLU:O	2:C:266:GLU:HG2	2.08	0.53
2:C:341:ARG:HG3	2:C:341:ARG:O	2.09	0.53
2:C:248:ARG:HD3	2:C:260:LYS:HA	1.91	0.53
2:C:360:THR:HG23	2:C:361:LEU:HD13	1.90	0.52
4:A:275:ARG:HA	4:A:278:THR:OG1	2.09	0.52
1:R:8:DA:OP1	4:A:209:GLY:CA	2.56	0.52
2:C:72:ILE:O	2:C:74:PRO:HD3	2.10	0.52
4:A:236:ILE:HD13	4:A:271:MSE:HE3	1.92	0.52
4:A:261:ALA:HB2	5:H:26:DT:O3'	2.09	0.51
2:C:146:MSE:HE2	2:C:189:MSE:HB3	1.92	0.51
4:A:14:GLU:OE2	7:A:605:HOH:O	2.19	0.51
4:A:297:PRO:O	4:A:298:SER:C	2.48	0.51
4:A:57:ILE:HG23	4:A:314:PHE:HE1	1.75	0.51
4:A:72:ILE:O	4:A:74:PRO:HD3	2.11	0.51
2:C:203:ASP:HB3	2:C:267:ASN:HB3	1.93	0.50
4:A:463:GLN:O	4:A:465:SER:N	2.43	0.50
4:A:223:LYS:HA	4:A:224:GLU:HB2	1.92	0.50
2:C:340:ASP:C	2:C:342:TYR:H	2.14	0.50
1:R:23:DT:OP1	2:C:211:LYS:CB	2.60	0.49
4:A:326:ILE:HD11	4:A:392:LEU:HB2	1.95	0.49
2:C:257:THR:OG1	3:X:12:DG:H4'	2.12	0.49
4:A:342:TYR:CD1	4:A:342:TYR:C	2.86	0.49
2:C:342:TYR:C	2:C:342:TYR:CD1	2.85	0.49
2:C:275:ARG:C	2:C:277:TYR:N	2.65	0.48
1:R:1:DT:C4	5:H:30:DA:N1	2.81	0.48
2:C:257:THR:CB	3:X:12:DG:H4'	2.43	0.48
1:R:15:DA:H4'	1:R:16:DA:OP1	2.13	0.48
4:A:341:ARG:CG	4:A:341:ARG:NH1	2.73	0.48
4:A:334:LEU:HB2	4:A:389:LEU:HD21	1.94	0.48
2:C:203:ASP:OD1	2:C:267:ASN:HB3	2.12	0.47
1:R:26:DA:C2	3:X:5:DT:N3	2.75	0.47
4:A:7:TYR:HE1	5:H:18:DA:H5''	1.79	0.47
2:C:39:MSE:HE3	2:C:107:ILE:HD11	1.97	0.46
5:H:21:DG:H1'	5:H:22:DG:H5'	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:341:ARG:NH2	2:C:358:GLU:CD	2.64	0.46
1:R:19:DC:H42	3:X:12:DG:H1	1.62	0.46
2:C:2:GLY:HA2	2:C:3:ILE:HA	1.71	0.46
2:C:146:MSE:HG3	2:C:189:MSE:HE2	1.97	0.46
2:C:196:SER:O	2:C:208:CYS:O	2.34	0.46
2:C:47:GLN:HG2	2:C:48:ALA:HB2	1.98	0.46
2:C:4:LYS:NZ	7:X:101:HOH:O	2.12	0.46
1:R:17:DC:H2'	1:R:18:DT:H72	1.98	0.46
4:A:239:TRP:HD1	4:A:242:ARG:HH21	1.64	0.46
2:C:39:MSE:CE	2:C:107:ILE:HD11	2.46	0.45
4:A:208:CYS:HA	4:A:209:GLY:HA3	1.62	0.45
2:C:120:GLU:OE2	2:C:276:TYR:OH	2.19	0.45
1:R:17:DC:H2'	1:R:18:DT:C7	2.47	0.45
2:C:49:ALA:O	2:C:55:PRO:HB2	2.17	0.44
4:A:243:LEU:O	4:A:247:LEU:N	2.46	0.44
1:R:27:DG:N2	3:X:4:DC:N3	2.65	0.44
2:C:208:CYS:HA	2:C:209:GLY:HA3	1.64	0.44
4:A:437:GLN:HG3	4:A:438:PRO:HD2	1.99	0.44
2:C:338:CYS:SG	2:C:451:GLY:HA3	2.58	0.44
4:A:42:TRP:O	4:A:46:ILE:HG12	2.18	0.44
4:A:248:ARG:HH21	4:A:263:GLU:HA	1.83	0.44
1:R:23:DT:P	2:C:211:LYS:HB2	2.58	0.43
2:C:39:MSE:SE	2:C:42:TRP:CD1	3.22	0.43
4:A:96:ASN:HA	4:A:97:GLY:HA2	1.72	0.43
2:C:131:GLN:O	2:C:292:ARG:NH1	2.51	0.43
4:A:221:PHE:CE1	4:A:243:LEU:HB3	2.53	0.43
4:A:270:ASN:OD1	4:A:272:GLU:HG2	2.19	0.43
2:C:203:ASP:HB2	2:C:270:ASN:HB2	2.01	0.43
5:H:20:DC:H2'	5:H:21:DG:C8	2.54	0.42
4:A:247:LEU:O	4:A:259:HIS:N	2.49	0.42
2:C:203:ASP:HB3	2:C:267:ASN:O	2.19	0.42
4:A:182:SER:HB2	4:A:184:LEU:HD23	2.00	0.42
4:A:192:VAL:HA	4:A:213:ALA:HB1	2.01	0.42
2:C:275:ARG:O	2:C:276:TYR:C	2.58	0.42
4:A:261:ALA:HB3	5:H:27:DT:OP1	2.14	0.42
2:C:275:ARG:HD2	2:C:279:HIS:ND1	2.35	0.42
5:H:22:DG:H1'	5:H:23:DT:H5'	2.01	0.42
4:A:339:LEU:HA	4:A:339:LEU:HD23	1.87	0.41
2:C:51:GLY:O	2:C:55:PRO:HB3	2.20	0.41
1:R:14:DC:C4	1:R:15:DA:N6	2.88	0.41
4:A:463:GLN:C	4:A:465:SER:N	2.73	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:269:PRO:HG2	4:A:271:MSE:HE1	2.03	0.41
4:A:184:LEU:HA	4:A:184:LEU:HD13	1.91	0.41
1:R:2:DC:H2"	1:R:3:DC:C5'	2.50	0.41
1:R:1:DT:H2"	1:R:2:DC:H6	1.86	0.41
2:C:47:GLN:HA	2:C:48:ALA:HA	1.47	0.41
4:A:223:LYS:HB3	4:A:225:LEU:HG	2.02	0.41
2:C:380:LEU:HB2	2:C:448:LEU:HD12	2.02	0.41
2:C:107:ILE:HG23	2:C:112:PHE:HB2	2.03	0.41
2:C:197:GLY:HA2	2:C:198:GLY:HA3	1.63	0.40
2:C:152:THR:HG23	2:C:172:HIS:HB2	2.02	0.40
2:C:341:ARG:CG	2:C:341:ARG:O	2.70	0.40
2:C:57:ILE:HG23	2:C:314:PHE:HE1	1.86	0.40
4:A:131:GLN:O	4:A:292:ARG:NH2	2.48	0.40
1:R:7:DC:H2"	1:R:8:DA:C8	2.57	0.40
4:A:397:GLU:HA	4:A:398:PRO:HD3	1.93	0.40
4:A:340:ASP:C	4:A:342:TYR:N	2.73	0.40

All (16) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:342:TYR:OH	4:A:342:TYR:OH[5_444]	0.19	2.01
2:C:342:TYR:CZ	4:A:342:TYR:CZ[5_444]	0.23	1.97
2:C:342:TYR:CE2	4:A:342:TYR:CE2[5_444]	0.23	1.97
2:C:342:TYR:CD2	4:A:342:TYR:CD2[5_444]	0.30	1.90
2:C:342:TYR:CE1	4:A:342:TYR:CE1[5_444]	0.31	1.89
2:C:342:TYR:CG	4:A:342:TYR:CG[5_444]	0.38	1.82
2:C:342:TYR:CD1	4:A:342:TYR:CD1[5_444]	0.39	1.81
2:C:342:TYR:CB	4:A:342:TYR:CB[5_444]	0.48	1.72
2:C:342:TYR:CB	4:A:342:TYR:CG[5_444]	1.69	0.51
2:C:342:TYR:CG	4:A:342:TYR:CD1[5_444]	1.73	0.47
2:C:342:TYR:CB	4:A:342:TYR:CA[5_444]	1.89	0.31
2:C:342:TYR:C	4:A:342:TYR:CB[5_444]	2.06	0.14
2:C:342:TYR:CD1	4:A:342:TYR:CD2[5_444]	2.09	0.11
2:C:342:TYR:CE1	4:A:342:TYR:CG[5_444]	2.14	0.06
2:C:342:TYR:CE1	4:A:342:TYR:CE2[5_444]	2.16	0.04
2:C:342:TYR:CZ	4:A:342:TYR:CD2[5_444]	2.18	0.02

## 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	
2	C	385/464 (83%)	359 (93%)	26 (7%)	0	100	100
4	A	379/464 (82%)	358 (94%)	20 (5%)	1 (0%)	46	63
All	All	764/928 (82%)	717 (94%)	46 (6%)	1 (0%)	56	74

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
4	A	464	ARG

### 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	
2	C	332/383 (87%)	324 (98%)	8 (2%)	57	76
4	A	329/385 (86%)	320 (97%)	9 (3%)	52	73
All	All	661/768 (86%)	644 (97%)	17 (3%)	54	74

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	C	109	LEU
2	C	184	LEU
2	C	266	GLU
2	C	267	ASN
2	C	275	ARG

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Mol	Chain	Res	Type
2	C	281	VAL
2	C	341	ARG
2	C	342	TYR
4	A	128	LEU
4	A	184	LEU
4	A	243	LEU
4	A	281	VAL
4	A	337	ARG
4	A	338	CYS
4	A	341	ARG
4	A	342	TYR
4	A	464	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
2	C	43	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

Of 2 ligands modelled in this entry, 2 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

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, 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	R	31/31 (100%)	-0.20	2 (6%) 22 22	56, 70, 98, 110	31 (100%)
2	C	390/464 (84%)	0.19	18 (4%) 36 37	27, 47, 82, 147	390 (100%)
3	X	16/16 (100%)	-0.25	1 (6%) 23 24	46, 67, 77, 91	16 (100%)
4	A	388/464 (83%)	0.16	15 (3%) 43 44	27, 48, 81, 147	388 (100%)
5	H	15/15 (100%)	-0.39	0 100 100	55, 65, 76, 86	15 (100%)
All	All	840/990 (84%)	0.14	36 (4%) 39 40	27, 49, 84, 147	840 (100%)

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	A	44	PHE	6.9
2	C	44	PHE	6.7
4	A	243	LEU	5.1
2	C	201	LEU	3.6
2	C	248	ARG	3.6
2	C	241	GLN	3.5
4	A	262	LEU	3.4
2	C	249	THR	3.3
2	C	200	TYR	3.3
4	A	264	ILE	3.2
2	C	202	PRO	3.1
2	C	243	LEU	3.1
2	C	262	LEU	3.0
2	C	235	ALA	2.9
2	C	255	PHE	2.8
2	C	193	ALA	2.8
4	A	263	GLU	2.7
2	C	275	ARG	2.7
4	A	275	ARG	2.6
4	A	235	ALA	2.6

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Mol	Chain	Res	Type	RSRZ
2	C	263	GLU	2.6
4	A	205	ILE	2.5
3	X	1	DA	2.5
1	R	15	DA	2.5
2	C	47	GLN	2.5
1	R	30	DT	2.4
4	A	248	ARG	2.4
2	C	247	LEU	2.3
2	C	244	LEU	2.3
4	A	239	TRP	2.2
4	A	256	ARG	2.2
4	A	41	ILE	2.1
4	A	241	GLN	2.1
4	A	342	TYR	2.1
2	C	261	ALA	2.0
4	A	461	GLU	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 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, 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	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
6	MG	C	501	1/1	0.94	0.06	-2.66	52,52,52,52	1
6	MG	A	501	1/1	0.96	0.04	-2.95	62,62,62,62	1

## 6.5 Other polymers

There are no such residues in this entry.