



# Full wwPDB X-ray Structure Validation Report ⓘ

May 30, 2016 – 01:21 PM EDT

PDB ID : 5AYS  
Title : Crystal structure of SAUGI/HSV UDG complex  
Authors : Wang, H.C.; Ko, T.P.; Huang, M.F.; Wang, A.H.J.  
Deposited on : 2015-09-02  
Resolution : 2.09 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-20027674  
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-20027674

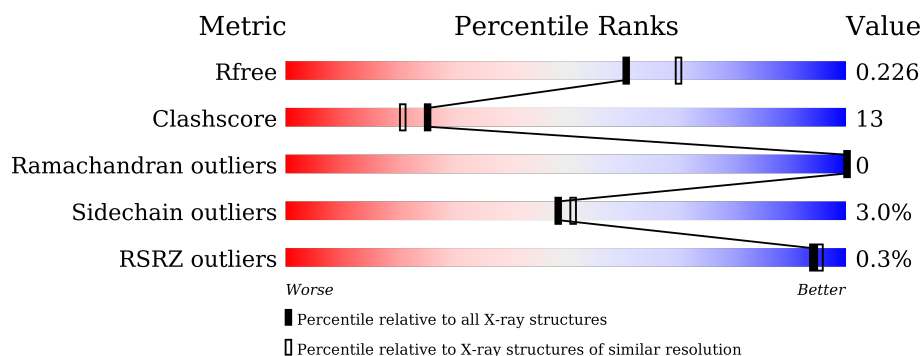
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.09 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	3939 (2.10-2.10)
Clashscore	102246	4460 (2.10-2.10)
Ramachandran outliers	100387	4413 (2.10-2.10)
Sidechain outliers	100360	4414 (2.10-2.10)
RSRZ outliers	91569	3948 (2.10-2.10)

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	A	256	
1	B	256	
2	C	112	
2	D	112	

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6232 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 Uracil-DNA glycosylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	230	Total	C	N	O	S	0	0	0
			1835	1170	344	312	9			
1	B	230	Total	C	N	O	S	0	0	0
			1835	1170	344	312	9			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	245	LEU	-	expression tag	UNP P10186
A	246	GLU	-	expression tag	UNP P10186
A	247	GLU	-	expression tag	UNP P10186
A	248	LEU	-	expression tag	UNP P10186
A	249	LEU	-	expression tag	UNP P10186
A	250	GLU	-	expression tag	UNP P10186
A	251	HIS	-	expression tag	UNP P10186
A	252	HIS	-	expression tag	UNP P10186
A	253	HIS	-	expression tag	UNP P10186
A	254	HIS	-	expression tag	UNP P10186
A	255	HIS	-	expression tag	UNP P10186
A	256	HIS	-	expression tag	UNP P10186
B	245	LEU	-	expression tag	UNP P10186
B	246	GLU	-	expression tag	UNP P10186
B	247	GLU	-	expression tag	UNP P10186
B	248	LEU	-	expression tag	UNP P10186
B	249	LEU	-	expression tag	UNP P10186
B	250	GLU	-	expression tag	UNP P10186
B	251	HIS	-	expression tag	UNP P10186
B	252	HIS	-	expression tag	UNP P10186
B	253	HIS	-	expression tag	UNP P10186
B	254	HIS	-	expression tag	UNP P10186
B	255	HIS	-	expression tag	UNP P10186
B	256	HIS	-	expression tag	UNP P10186

- Molecule 2 is a protein called Uncharacterized protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	109	Total	C	N	O	S	0	0	0
			925	603	141	177	4			
2	D	109	Total	C	N	O	S	0	0	0
			925	603	141	177	4			

- Molecule 3 is water.

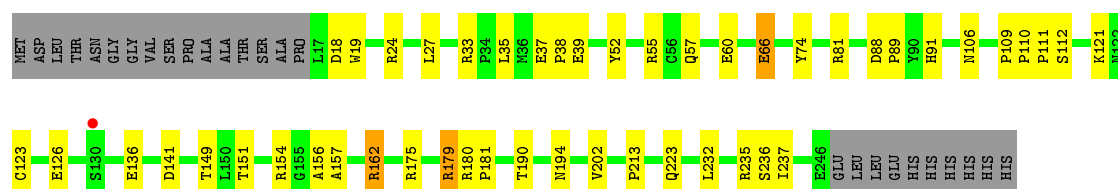
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	220	Total	O	0	0
			220	220		
3	C	129	Total	O	0	0
			129	129		
3	B	227	Total	O	0	0
			227	227		
3	D	136	Total	O	0	0
			136	136		

### 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 ( $RSRZ > 2$ ). 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.

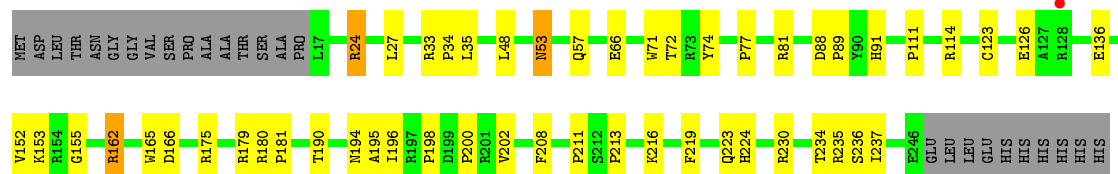
- Molecule 1: Uracil-DNA glycosylase

Chain A: 



- Molecule 1: Uracil-DNA glycosylase

Chain B: 




- Molecule 2: Uncharacterized protein

Chain C: 



- Molecule 2: Uncharacterized protein

Chain D: 



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	53.30Å 63.01Å 65.51Å 65.10° 89.58° 88.38°	Depositor
Resolution (Å)	20.00 – 2.09 19.88 – 2.09	Depositor EDS
% Data completeness (in resolution range)	94.4 (20.00-2.09) 81.3 (19.88-2.09)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.48 (at 2.09Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, $R_{free}$	0.195 , 0.223 0.197 , 0.226	Depositor DCC
$R_{free}$ test set	2190 reflections (5.31%)	DCC
Wilson B-factor (Å <sup>2</sup> )	22.6	Xtriage
Anisotropy	0.158	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 49.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.025 for h,-k,-l 0.000 for -h,l,k 0.000 for -h,-l,-k	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	6232	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.78% 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 [i](#)

### 5.1 Standard geometry [i](#)

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	0.48	0/1892	0.71	0/2584
1	B	0.48	0/1892	0.71	0/2584
2	C	0.46	0/948	0.76	0/1290
2	D	0.49	0/948	0.77	0/1290
All	All	0.48	0/5680	0.73	0/7748

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	A	1835	0	1817	42	0
1	B	1835	0	1817	43	0
2	C	925	0	903	36	0
2	D	925	0	903	26	0
3	A	220	0	0	13	1
3	B	227	0	0	11	3
3	C	129	0	0	21	1
3	D	136	0	0	9	1
All	All	6232	0	5440	138	3

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 13.

All (138) 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)
2:C:13:ASN:HB3	3:C:290:HOH:O	1.34	1.25
2:D:85:GLU:HG3	3:D:209:HOH:O	1.61	0.97
1:B:230:ARG:NE	3:B:301:HOH:O	1.84	0.96
2:C:9:HIS:HB3	3:C:213:HOH:O	1.67	0.93
2:C:85:GLU:HG3	3:C:207:HOH:O	1.68	0.91
1:A:18:ASP:OD1	3:A:301:HOH:O	1.94	0.86
1:A:106:ASN:ND2	3:A:302:HOH:O	1.95	0.85
2:D:65:ASN:HB2	2:D:85:GLU:HG2	1.58	0.84
2:C:65:ASN:HB2	2:C:85:GLU:HG2	1.60	0.81
1:A:35:LEU:HD21	1:A:179:ARG:HD3	1.62	0.81
2:D:53:THR:HG23	3:D:232:HOH:O	1.82	0.79
1:A:88:ASP:HB2	1:A:89:PRO:CD	2.12	0.78
1:B:35:LEU:HD21	1:B:179:ARG:HD3	1.65	0.77
1:B:166:ASP:HB3	3:B:343:HOH:O	1.84	0.76
1:A:88:ASP:HB2	1:A:89:PRO:HD3	1.66	0.74
1:B:181:PRO:HB3	1:B:202:VAL:HG13	1.71	0.72
2:C:53:THR:CG2	3:C:221:HOH:O	2.38	0.72
2:C:85:GLU:CG	3:C:207:HOH:O	2.29	0.72
1:B:126:GLU:HB2	3:B:390:HOH:O	1.91	0.70
2:C:88:ASP:OD2	3:C:202:HOH:O	2.08	0.70
2:C:28:ILE:HA	3:C:265:HOH:O	1.92	0.70
1:B:190:THR:HG22	1:B:194:ASN:HD21	1.57	0.69
2:D:83:ILE:HD11	2:D:89:MET:HE3	1.75	0.68
2:C:53:THR:HG23	3:C:221:HOH:O	1.93	0.68
1:A:126:GLU:HB2	3:A:359:HOH:O	1.94	0.68
1:B:190:THR:HG22	1:B:194:ASN:ND2	2.08	0.67
2:C:28:ILE:O	2:C:53:THR:HB	1.96	0.66
2:D:85:GLU:OE1	3:D:201:HOH:O	2.13	0.66
1:B:180:ARG:O	3:B:302:HOH:O	2.13	0.66
1:B:136:GLU:OE1	3:B:303:HOH:O	2.14	0.65
1:A:74:TYR:CE1	1:A:136:GLU:HG2	2.31	0.65
2:D:83:ILE:HG12	2:D:89:MET:HG2	1.78	0.63
2:D:31:ILE:HG22	2:D:33:ASP:H	1.64	0.62
2:D:53:THR:CG2	3:D:232:HOH:O	2.42	0.62
1:B:88:ASP:HB2	1:B:89:PRO:CD	2.31	0.61
1:B:66:GLU:CD	1:B:66:GLU:H	2.03	0.61
2:C:87:HIS:NE2	3:C:203:HOH:O	2.10	0.60
1:A:121:LYS:NZ	3:A:306:HOH:O	2.30	0.60
2:D:83:ILE:HD11	2:D:89:MET:CE	2.31	0.60
2:C:104:GLN:NE2	3:C:201:HOH:O	2.01	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:96:ASN:ND2	2:C:99:MET:H	1.99	0.59
1:A:37:GLU:HB3	3:A:341:HOH:O	2.02	0.59
1:B:74:TYR:CE1	1:B:136:GLU:HG2	2.38	0.59
1:A:88:ASP:HB3	3:A:333:HOH:O	2.02	0.58
1:A:190:THR:HG22	1:A:194:ASN:ND2	2.18	0.58
1:A:88:ASP:HB3	1:A:149:THR:HB	1.85	0.58
2:C:62:HIS:NE2	3:C:204:HOH:O	2.13	0.58
2:D:75:LYS:NZ	3:D:202:HOH:O	2.20	0.58
2:D:96:ASN:ND2	2:D:98:ILE:H	2.03	0.57
1:B:162:ARG:NH2	3:B:309:HOH:O	2.37	0.56
1:A:181:PRO:HB3	1:A:202:VAL:HG13	1.88	0.55
1:A:123:CYS:SG	1:A:223:GLN:HA	2.46	0.55
2:C:83:ILE:HG12	2:C:89:MET:HG2	1.90	0.54
1:B:48:LEU:HD21	1:B:165:TRP:CE3	2.43	0.54
2:C:109:THR:HG23	3:C:268:HOH:O	2.06	0.54
1:A:91:HIS:HE1	3:C:208:HOH:O	1.90	0.54
2:D:65:ASN:CB	2:D:85:GLU:HG2	2.35	0.54
1:B:91:HIS:HD2	1:B:155:GLY:CA	2.22	0.53
2:D:28:ILE:O	2:D:53:THR:HB	2.09	0.53
1:A:60:GLU:OE2	1:A:154:ARG:NH1	2.39	0.53
1:B:216:LYS:NZ	3:B:308:HOH:O	2.37	0.53
1:A:235:ARG:O	1:A:236:SER:HB2	2.08	0.53
1:A:27:LEU:HD12	1:A:66:GLU:HB3	1.89	0.53
1:B:230:ARG:O	1:B:234:THR:HG23	2.09	0.53
1:B:33:ARG:HB3	1:B:34:PRO:HD3	1.91	0.52
2:C:96:ASN:ND2	2:C:98:ILE:H	2.08	0.51
2:C:27:SER:HB3	3:C:208:HOH:O	2.10	0.51
1:A:66:GLU:CD	1:A:66:GLU:H	2.13	0.51
3:B:366:HOH:O	2:D:53:THR:HG21	2.11	0.50
2:D:96:ASN:ND2	2:D:99:MET:H	2.10	0.50
1:A:180:ARG:O	3:A:304:HOH:O	2.19	0.50
2:C:31:ILE:HG22	2:C:33:ASP:H	1.77	0.49
2:C:96:ASN:HD21	2:C:98:ILE:HB	1.77	0.49
1:B:53:ASN:ND2	3:B:312:HOH:O	2.46	0.49
2:C:96:ASN:HD21	2:C:99:MET:H	1.59	0.49
1:A:91:HIS:CE1	3:C:208:HOH:O	2.64	0.49
1:A:156:ALA:HB1	3:A:397:HOH:O	2.13	0.48
1:A:111:PRO:HD3	2:C:57:TYR:CD1	2.48	0.48
1:B:111:PRO:HD2	2:D:26:GLU:HB2	1.95	0.48
1:B:91:HIS:HD2	1:B:155:GLY:HA2	1.77	0.48
1:A:141:ASP:OD2	3:A:305:HOH:O	2.20	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:35:LEU:HD11	1:B:77:PRO:HB3	1.96	0.48
1:B:88:ASP:HB2	1:B:89:PRO:HD3	1.96	0.48
2:C:1:MET:HA	3:C:238:HOH:O	2.13	0.48
1:B:81:ARG:HD2	1:B:237:ILE:HG21	1.96	0.47
1:B:175:ARG:HD3	3:B:481:HOH:O	2.14	0.47
1:B:111:PRO:HB3	1:B:213:PRO:HG3	1.95	0.47
1:A:190:THR:HG22	1:A:194:ASN:HD21	1.80	0.47
1:B:114:ARG:NH1	2:D:62:HIS:CE1	2.82	0.47
1:A:111:PRO:HB3	1:A:213:PRO:HG3	1.97	0.47
1:B:91:HIS:CD2	1:B:155:GLY:CA	2.98	0.47
2:D:27:SER:HA	2:D:53:THR:O	2.16	0.46
1:A:81:ARG:HD2	1:A:237:ILE:HG21	1.97	0.46
1:A:111:PRO:HD3	2:C:57:TYR:CE1	2.50	0.46
1:B:71:TRP:CZ3	1:B:72:THR:HG22	2.51	0.46
2:C:36:LEU:CD1	2:C:46:LEU:HD13	2.45	0.46
1:A:112:SER:HB2	3:C:210:HOH:O	2.16	0.46
1:B:216:LYS:HE3	2:D:36:LEU:O	2.16	0.45
1:B:216:LYS:HE2	3:D:210:HOH:O	2.15	0.45
1:B:91:HIS:HE1	2:D:27:SER:HB3	1.81	0.45
2:D:96:ASN:HD21	2:D:99:MET:H	1.63	0.45
2:D:85:GLU:CG	3:D:209:HOH:O	2.38	0.45
1:A:55:ARG:NH2	3:A:303:HOH:O	2.19	0.45
1:B:123:CYS:SG	1:B:223:GLN:HA	2.56	0.44
1:A:35:LEU:CD2	1:A:179:ARG:HD3	2.40	0.44
1:B:24:ARG:NH2	3:B:316:HOH:O	2.50	0.44
1:B:195:ALA:O	1:B:196:ILE:HD13	2.18	0.44
1:B:235:ARG:HB2	1:B:237:ILE:HD12	1.99	0.44
1:A:74:TYR:CZ	1:A:136:GLU:HG2	2.53	0.44
1:B:27:LEU:HD12	1:B:66:GLU:HB3	1.99	0.43
1:B:152:VAL:HG22	1:B:153:LYS:O	2.18	0.43
1:A:91:HIS:HE1	2:C:27:SER:HB3	1.83	0.43
1:A:19:TRP:CE2	1:A:33:ARG:HG3	2.53	0.43
1:B:198:PRO:O	1:B:200:PRO:HD3	2.19	0.43
2:C:46:LEU:HD23	2:C:46:LEU:HA	1.87	0.43
1:B:235:ARG:O	1:B:236:SER:HB2	2.18	0.43
1:A:157:ALA:HA	3:A:404:HOH:O	2.18	0.43
2:D:96:ASN:C	2:D:96:ASN:HD22	2.22	0.43
2:D:108:LEU:HB2	3:D:255:HOH:O	2.18	0.42
2:C:50:ILE:HD12	3:C:308:HOH:O	2.19	0.42
2:C:87:HIS:NE2	3:C:206:HOH:O	2.37	0.42
1:B:211:PRO:HA	1:B:219:PHE:HB2	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:208:PHE:HB2	1:B:224:HIS:CE1	2.54	0.42
1:A:232:LEU:HB3	1:A:237:ILE:HB	2.02	0.42
1:A:162:ARG:NH1	2:D:90:ASP:OD1	2.53	0.42
1:A:109:PRO:HA	1:A:110:PRO:HD3	1.92	0.41
3:A:377:HOH:O	2:C:62:HIS:HB2	2.20	0.41
1:B:74:TYR:CZ	1:B:136:GLU:HG2	2.55	0.41
1:A:39:GLU:HG2	1:A:175:ARG:NH2	2.36	0.41
2:C:92:LEU:HD23	2:C:92:LEU:HA	1.85	0.41
1:A:38:PRO:HD3	3:A:341:HOH:O	2.20	0.41
2:C:85:GLU:N	3:C:207:HOH:O	2.40	0.41
1:A:52:TYR:CE1	1:A:151:THR:HG21	2.56	0.41
1:A:190:THR:OG1	2:C:34:ASP:OD2	2.38	0.41
2:C:2:THR:OG1	2:C:5:LEU:HG	2.21	0.41
2:C:87:HIS:CE1	3:C:206:HOH:O	2.73	0.40
2:C:47:SER:O	2:C:48:ASN:HB2	2.21	0.40
2:D:109:THR:N	3:D:207:HOH:O	2.52	0.40

All (3) 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)
3:B:301:HOH:O	3:D:227:HOH:O[1_455]	1.84	0.36
3:A:375:HOH:O	3:B:356:HOH:O[1_565]	2.11	0.09
3:C:202:HOH:O	3:B:381:HOH:O[1_556]	2.18	0.02

## 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	228/256 (89%)	223 (98%)	5 (2%)	0	100	100
1	B	228/256 (89%)	222 (97%)	6 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	C	107/112 (96%)	101 (94%)	6 (6%)	0	100	100
2	D	107/112 (96%)	103 (96%)	4 (4%)	0	100	100
All	All	670/736 (91%)	649 (97%)	21 (3%)	0	100	100

There are no Ramachandran outliers to report.

### 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	196/219 (90%)	191 (97%)	5 (3%)	54	58
1	B	196/219 (90%)	192 (98%)	4 (2%)	63	68
2	C	104/105 (99%)	99 (95%)	5 (5%)	31	29
2	D	104/105 (99%)	100 (96%)	4 (4%)	40	40
All	All	600/648 (93%)	582 (97%)	18 (3%)	48	51

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	24	ARG
1	A	57	GLN
1	A	66	GLU
1	A	162	ARG
1	A	179	ARG
2	C	33	ASP
2	C	53	THR
2	C	90	ASP
2	C	96	ASN
2	C	109	THR
1	B	24	ARG
1	B	53	ASN
1	B	57	GLN
1	B	162	ARG

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Mol	Chain	Res	Type
2	D	53	THR
2	D	90	ASP
2	D	96	ASN
2	D	109	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	91	HIS
1	A	92	HIS
1	A	194	ASN
2	C	9	HIS
2	C	96	ASN
1	B	91	HIS
1	B	92	HIS
1	B	191	HIS
1	B	194	ASN
2	D	9	HIS
2	D	74	GLN
2	D	87	HIS
2	D	96	ASN
2	D	104	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 ⓘ

There are no ligands in this entry.

## 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 [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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	A	230/256 (89%)	-0.24	1 (0%) 93 94	12, 23, 42, 54	0
1	B	230/256 (89%)	-0.23	1 (0%) 93 94	13, 23, 41, 61	0
2	C	109/112 (97%)	-0.21	0 100 100	15, 25, 40, 46	0
2	D	109/112 (97%)	-0.18	0 100 100	14, 24, 39, 46	0
All	All	678/736 (92%)	-0.22	2 (0%) 94 95	12, 24, 41, 61	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	130	SER	2.2
1	B	128	ARG	2.1

### 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](#)

There are no ligands in this entry.

### 6.5 Other polymers [i](#)

There are no such residues in this entry.