



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 03:54 PM GMT

PDB ID : 4DOY  
Title : Crystal structure of Dibenzothiophene desulfurization enzyme C  
Authors : Liu, S.; Zhang, C.; Zhu, D.; Gu, L.  
Deposited on : 2012-02-12  
Resolution : 1.79 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 : 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

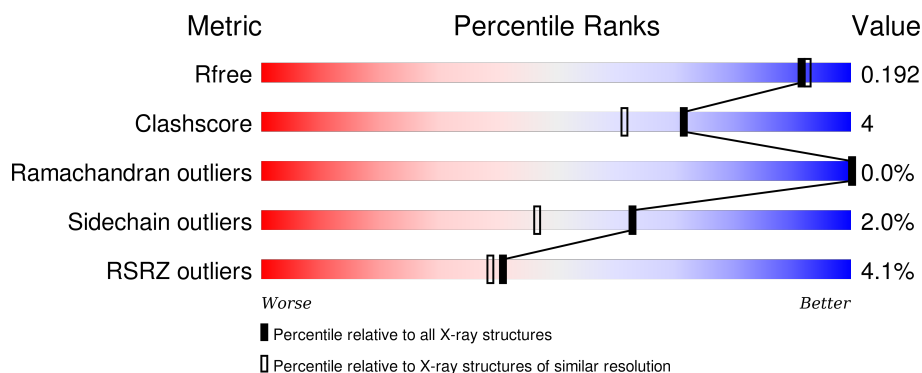
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.79 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	6655 (1.80-1.76)
Clashscore	102246	7658 (1.80-1.76)
Ramachandran outliers	100387	7570 (1.80-1.76)
Sidechain outliers	100360	7569 (1.80-1.76)
RSRZ outliers	91569	6671 (1.80-1.76)

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	437	 85% 6% 8%
1	B	437	 4% 81% 10% 8%
1	C	437	 % 84% 7% 8%
1	D	437	 % 85% 6% 8%
1	E	437	 2% 85% 6% 8%

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Mol	Chain	Length	Quality of chain
1	F	437	
1	G	437	
1	H	437	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	GOL	A	500	-	-	-	X
2	GOL	C	500	-	-	-	X
2	GOL	D	501	-	-	-	X
2	GOL	D	502	-	-	-	X
2	GOL	E	502	-	-	-	X
2	GOL	G	501	-	-	-	X
2	GOL	G	502	-	-	-	X
2	GOL	H	502	-	-	-	X

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 27568 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 Dibenzothiophene desulfurization enzyme C.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	400	Total	C	N	O	S	0	7	0
			3109	1949	555	601	4			
1	B	400	Total	C	N	O	S	0	8	0
			3119	1957	555	603	4			
1	C	400	Total	C	N	O	S	0	4	0
			3083	1935	549	595	4			
1	D	400	Total	C	N	O	S	0	6	0
			3099	1943	553	599	4			
1	E	400	Total	C	N	O	S	0	6	0
			3105	1949	554	598	4			
1	F	400	Total	C	N	O	S	0	6	0
			3103	1948	552	599	4			
1	G	400	Total	C	N	O	S	0	5	0
			3092	1940	549	599	4			
1	H	400	Total	C	N	O	S	0	2	0
			3066	1926	544	592	4			

There are 160 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	EXPRESSION TAG	UNP Q6WNP1
A	-18	GLY	-	EXPRESSION TAG	UNP Q6WNP1
A	-17	SER	-	EXPRESSION TAG	UNP Q6WNP1
A	-16	SER	-	EXPRESSION TAG	UNP Q6WNP1
A	-15	HIS	-	EXPRESSION TAG	UNP Q6WNP1
A	-14	HIS	-	EXPRESSION TAG	UNP Q6WNP1
A	-13	HIS	-	EXPRESSION TAG	UNP Q6WNP1
A	-12	HIS	-	EXPRESSION TAG	UNP Q6WNP1
A	-11	HIS	-	EXPRESSION TAG	UNP Q6WNP1
A	-10	HIS	-	EXPRESSION TAG	UNP Q6WNP1
A	-9	SER	-	EXPRESSION TAG	UNP Q6WNP1
A	-8	SER	-	EXPRESSION TAG	UNP Q6WNP1
A	-7	GLY	-	EXPRESSION TAG	UNP Q6WNP1

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-6	LEU	-	EXPRESSION TAG	UNP Q6WNP1
A	-5	VAL	-	EXPRESSION TAG	UNP Q6WNP1
A	-4	PRO	-	EXPRESSION TAG	UNP Q6WNP1
A	-3	ARG	-	EXPRESSION TAG	UNP Q6WNP1
A	-2	GLY	-	EXPRESSION TAG	UNP Q6WNP1
A	-1	SER	-	EXPRESSION TAG	UNP Q6WNP1
A	0	HIS	-	EXPRESSION TAG	UNP Q6WNP1
B	-19	MET	-	EXPRESSION TAG	UNP Q6WNP1
B	-18	GLY	-	EXPRESSION TAG	UNP Q6WNP1
B	-17	SER	-	EXPRESSION TAG	UNP Q6WNP1
B	-16	SER	-	EXPRESSION TAG	UNP Q6WNP1
B	-15	HIS	-	EXPRESSION TAG	UNP Q6WNP1
B	-14	HIS	-	EXPRESSION TAG	UNP Q6WNP1
B	-13	HIS	-	EXPRESSION TAG	UNP Q6WNP1
B	-12	HIS	-	EXPRESSION TAG	UNP Q6WNP1
B	-11	HIS	-	EXPRESSION TAG	UNP Q6WNP1
B	-10	HIS	-	EXPRESSION TAG	UNP Q6WNP1
B	-9	SER	-	EXPRESSION TAG	UNP Q6WNP1
B	-8	SER	-	EXPRESSION TAG	UNP Q6WNP1
B	-7	GLY	-	EXPRESSION TAG	UNP Q6WNP1
B	-6	LEU	-	EXPRESSION TAG	UNP Q6WNP1
B	-5	VAL	-	EXPRESSION TAG	UNP Q6WNP1
B	-4	PRO	-	EXPRESSION TAG	UNP Q6WNP1
B	-3	ARG	-	EXPRESSION TAG	UNP Q6WNP1
B	-2	GLY	-	EXPRESSION TAG	UNP Q6WNP1
B	-1	SER	-	EXPRESSION TAG	UNP Q6WNP1
B	0	HIS	-	EXPRESSION TAG	UNP Q6WNP1
C	-19	MET	-	EXPRESSION TAG	UNP Q6WNP1
C	-18	GLY	-	EXPRESSION TAG	UNP Q6WNP1
C	-17	SER	-	EXPRESSION TAG	UNP Q6WNP1
C	-16	SER	-	EXPRESSION TAG	UNP Q6WNP1
C	-15	HIS	-	EXPRESSION TAG	UNP Q6WNP1
C	-14	HIS	-	EXPRESSION TAG	UNP Q6WNP1
C	-13	HIS	-	EXPRESSION TAG	UNP Q6WNP1
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C	-9	SER	-	EXPRESSION TAG	UNP Q6WNP1
C	-8	SER	-	EXPRESSION TAG	UNP Q6WNP1
C	-7	GLY	-	EXPRESSION TAG	UNP Q6WNP1
C	-6	LEU	-	EXPRESSION TAG	UNP Q6WNP1
C	-5	VAL	-	EXPRESSION TAG	UNP Q6WNP1

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-4	PRO	-	EXPRESSION TAG	UNP Q6WNP1
C	-3	ARG	-	EXPRESSION TAG	UNP Q6WNP1
C	-2	GLY	-	EXPRESSION TAG	UNP Q6WNP1
C	-1	SER	-	EXPRESSION TAG	UNP Q6WNP1
C	0	HIS	-	EXPRESSION TAG	UNP Q6WNP1
D	-19	MET	-	EXPRESSION TAG	UNP Q6WNP1
D	-18	GLY	-	EXPRESSION TAG	UNP Q6WNP1
D	-17	SER	-	EXPRESSION TAG	UNP Q6WNP1
D	-16	SER	-	EXPRESSION TAG	UNP Q6WNP1
D	-15	HIS	-	EXPRESSION TAG	UNP Q6WNP1
D	-14	HIS	-	EXPRESSION TAG	UNP Q6WNP1
D	-13	HIS	-	EXPRESSION TAG	UNP Q6WNP1
D	-12	HIS	-	EXPRESSION TAG	UNP Q6WNP1
D	-11	HIS	-	EXPRESSION TAG	UNP Q6WNP1
D	-10	HIS	-	EXPRESSION TAG	UNP Q6WNP1
D	-9	SER	-	EXPRESSION TAG	UNP Q6WNP1
D	-8	SER	-	EXPRESSION TAG	UNP Q6WNP1
D	-7	GLY	-	EXPRESSION TAG	UNP Q6WNP1
D	-6	LEU	-	EXPRESSION TAG	UNP Q6WNP1
D	-5	VAL	-	EXPRESSION TAG	UNP Q6WNP1
D	-4	PRO	-	EXPRESSION TAG	UNP Q6WNP1
D	-3	ARG	-	EXPRESSION TAG	UNP Q6WNP1
D	-2	GLY	-	EXPRESSION TAG	UNP Q6WNP1
D	-1	SER	-	EXPRESSION TAG	UNP Q6WNP1
D	0	HIS	-	EXPRESSION TAG	UNP Q6WNP1
E	-19	MET	-	EXPRESSION TAG	UNP Q6WNP1
E	-18	GLY	-	EXPRESSION TAG	UNP Q6WNP1
E	-17	SER	-	EXPRESSION TAG	UNP Q6WNP1
E	-16	SER	-	EXPRESSION TAG	UNP Q6WNP1
E	-15	HIS	-	EXPRESSION TAG	UNP Q6WNP1
E	-14	HIS	-	EXPRESSION TAG	UNP Q6WNP1
E	-13	HIS	-	EXPRESSION TAG	UNP Q6WNP1
E	-12	HIS	-	EXPRESSION TAG	UNP Q6WNP1
E	-11	HIS	-	EXPRESSION TAG	UNP Q6WNP1
E	-10	HIS	-	EXPRESSION TAG	UNP Q6WNP1
E	-9	SER	-	EXPRESSION TAG	UNP Q6WNP1
E	-8	SER	-	EXPRESSION TAG	UNP Q6WNP1
E	-7	GLY	-	EXPRESSION TAG	UNP Q6WNP1
E	-6	LEU	-	EXPRESSION TAG	UNP Q6WNP1
E	-5	VAL	-	EXPRESSION TAG	UNP Q6WNP1
E	-4	PRO	-	EXPRESSION TAG	UNP Q6WNP1
E	-3	ARG	-	EXPRESSION TAG	UNP Q6WNP1

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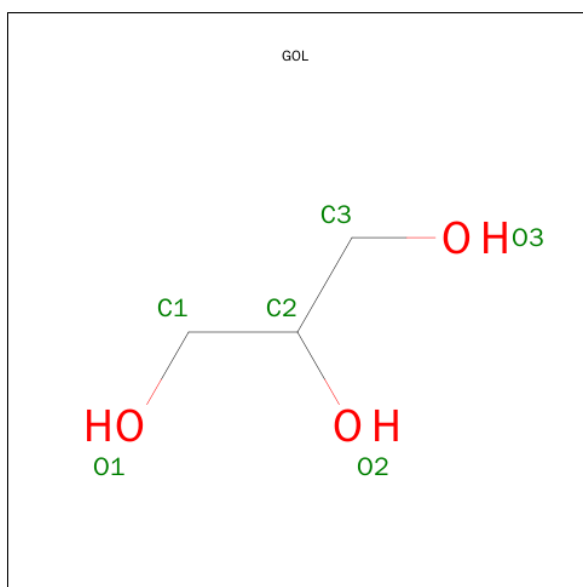
Chain	Residue	Modelled	Actual	Comment	Reference
E	-2	GLY	-	EXPRESSION TAG	UNP Q6WNP1
E	-1	SER	-	EXPRESSION TAG	UNP Q6WNP1
E	0	HIS	-	EXPRESSION TAG	UNP Q6WNP1
F	-19	MET	-	EXPRESSION TAG	UNP Q6WNP1
F	-18	GLY	-	EXPRESSION TAG	UNP Q6WNP1
F	-17	SER	-	EXPRESSION TAG	UNP Q6WNP1
F	-16	SER	-	EXPRESSION TAG	UNP Q6WNP1
F	-15	HIS	-	EXPRESSION TAG	UNP Q6WNP1
F	-14	HIS	-	EXPRESSION TAG	UNP Q6WNP1
F	-13	HIS	-	EXPRESSION TAG	UNP Q6WNP1
F	-12	HIS	-	EXPRESSION TAG	UNP Q6WNP1
F	-11	HIS	-	EXPRESSION TAG	UNP Q6WNP1
F	-10	HIS	-	EXPRESSION TAG	UNP Q6WNP1
F	-9	SER	-	EXPRESSION TAG	UNP Q6WNP1
F	-8	SER	-	EXPRESSION TAG	UNP Q6WNP1
F	-7	GLY	-	EXPRESSION TAG	UNP Q6WNP1
F	-6	LEU	-	EXPRESSION TAG	UNP Q6WNP1
F	-5	VAL	-	EXPRESSION TAG	UNP Q6WNP1
F	-4	PRO	-	EXPRESSION TAG	UNP Q6WNP1
F	-3	ARG	-	EXPRESSION TAG	UNP Q6WNP1
F	-2	GLY	-	EXPRESSION TAG	UNP Q6WNP1
F	-1	SER	-	EXPRESSION TAG	UNP Q6WNP1
F	0	HIS	-	EXPRESSION TAG	UNP Q6WNP1
G	-19	MET	-	EXPRESSION TAG	UNP Q6WNP1
G	-18	GLY	-	EXPRESSION TAG	UNP Q6WNP1
G	-17	SER	-	EXPRESSION TAG	UNP Q6WNP1
G	-16	SER	-	EXPRESSION TAG	UNP Q6WNP1
G	-15	HIS	-	EXPRESSION TAG	UNP Q6WNP1
G	-14	HIS	-	EXPRESSION TAG	UNP Q6WNP1
G	-13	HIS	-	EXPRESSION TAG	UNP Q6WNP1
G	-12	HIS	-	EXPRESSION TAG	UNP Q6WNP1
G	-11	HIS	-	EXPRESSION TAG	UNP Q6WNP1
G	-10	HIS	-	EXPRESSION TAG	UNP Q6WNP1
G	-9	SER	-	EXPRESSION TAG	UNP Q6WNP1
G	-8	SER	-	EXPRESSION TAG	UNP Q6WNP1
G	-7	GLY	-	EXPRESSION TAG	UNP Q6WNP1
G	-6	LEU	-	EXPRESSION TAG	UNP Q6WNP1
G	-5	VAL	-	EXPRESSION TAG	UNP Q6WNP1
G	-4	PRO	-	EXPRESSION TAG	UNP Q6WNP1
G	-3	ARG	-	EXPRESSION TAG	UNP Q6WNP1
G	-2	GLY	-	EXPRESSION TAG	UNP Q6WNP1
G	-1	SER	-	EXPRESSION TAG	UNP Q6WNP1

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Chain	Residue	Modelled	Actual	Comment	Reference
G	0	HIS	-	EXPRESSION TAG	UNP Q6WNP1
H	-19	MET	-	EXPRESSION TAG	UNP Q6WNP1
H	-18	GLY	-	EXPRESSION TAG	UNP Q6WNP1
H	-17	SER	-	EXPRESSION TAG	UNP Q6WNP1
H	-16	SER	-	EXPRESSION TAG	UNP Q6WNP1
H	-15	HIS	-	EXPRESSION TAG	UNP Q6WNP1
H	-14	HIS	-	EXPRESSION TAG	UNP Q6WNP1
H	-13	HIS	-	EXPRESSION TAG	UNP Q6WNP1
H	-12	HIS	-	EXPRESSION TAG	UNP Q6WNP1
H	-11	HIS	-	EXPRESSION TAG	UNP Q6WNP1
H	-10	HIS	-	EXPRESSION TAG	UNP Q6WNP1
H	-9	SER	-	EXPRESSION TAG	UNP Q6WNP1
H	-8	SER	-	EXPRESSION TAG	UNP Q6WNP1
H	-7	GLY	-	EXPRESSION TAG	UNP Q6WNP1
H	-6	LEU	-	EXPRESSION TAG	UNP Q6WNP1
H	-5	VAL	-	EXPRESSION TAG	UNP Q6WNP1
H	-4	PRO	-	EXPRESSION TAG	UNP Q6WNP1
H	-3	ARG	-	EXPRESSION TAG	UNP Q6WNP1
H	-2	GLY	-	EXPRESSION TAG	UNP Q6WNP1
H	-1	SER	-	EXPRESSION TAG	UNP Q6WNP1
H	0	HIS	-	EXPRESSION TAG	UNP Q6WNP1

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			6	3	3		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total C O 6 3 3	0	0
2	B	1	Total C O 6 3 3	0	0
2	C	1	Total C O 6 3 3	0	0
2	D	1	Total O 1 1	0	0
2	D	1	Total C O 5 3 2	0	0
2	E	1	Total C O 6 3 3	0	0
2	F	1	Total C O 6 3 3	0	0
2	G	1	Total C O 6 3 3	0	0
2	E	1	Total C O 6 3 3	0	0
2	H	1	Total C O 6 3 3	0	0
2	G	1	Total C O 6 3 3	0	0
2	H	1	Total C O 6 3 3	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	413	Total O 413 413	0	0
3	B	384	Total O 384 384	0	0
3	C	432	Total O 432 432	0	0
3	D	387	Total O 387 387	0	0
3	E	343	Total O 343 343	0	0
3	F	310	Total O 310 310	0	0
3	G	227	Total O 227 227	0	0

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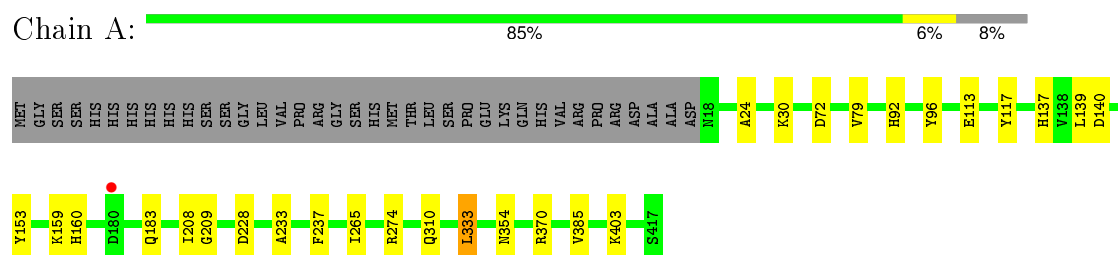
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	H	224	Total 224	O 224	0	0

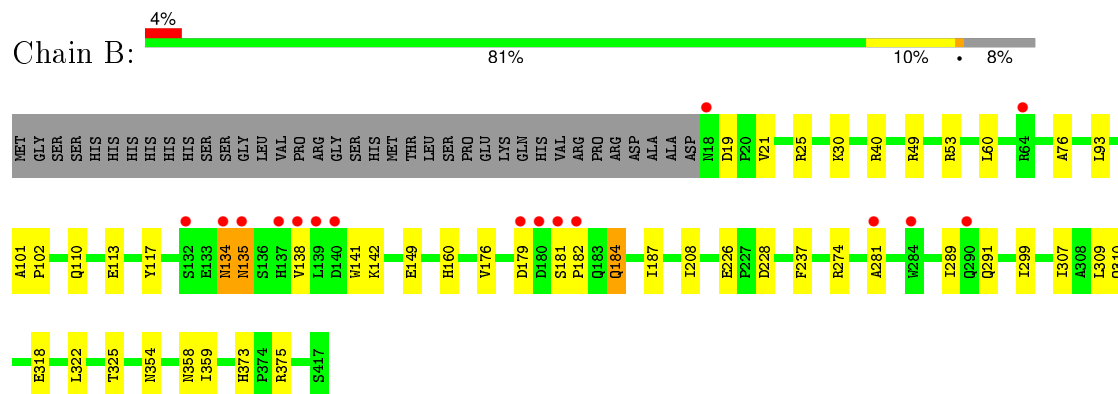
### 3 Residue-property plots [i](#)

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.

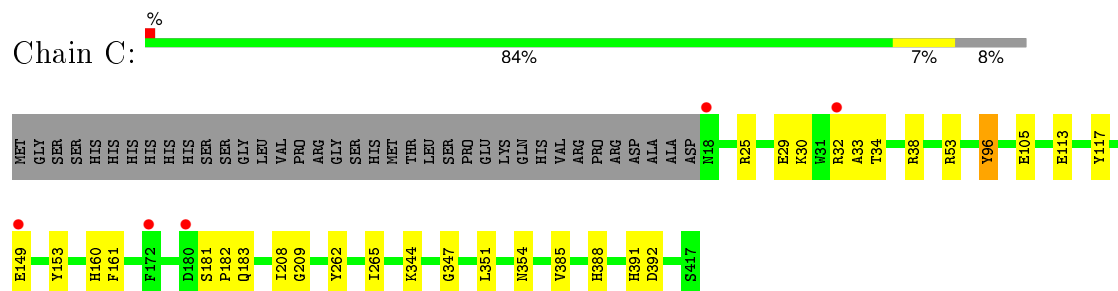
- Molecule 1: Dibenzothiophene desulfurization enzyme C



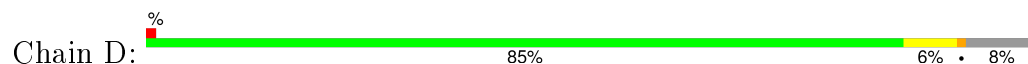
- Molecule 1: Dibenzothiophene desulfurization enzyme C

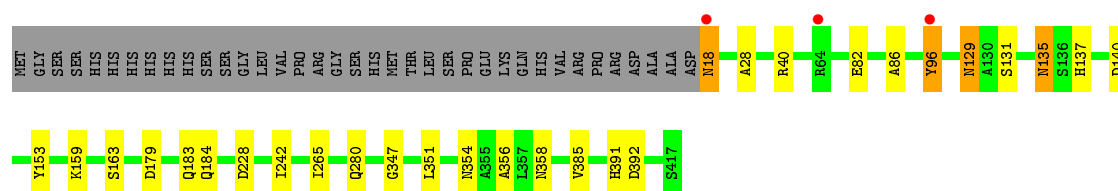


- Molecule 1: Dibenzothiophene desulfurization enzyme C

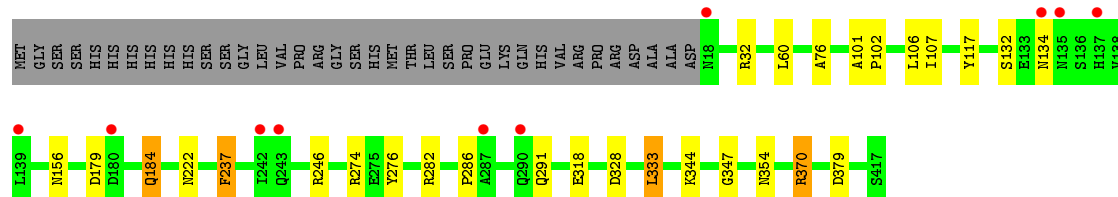
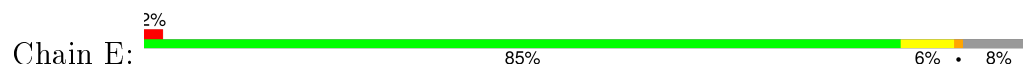


- Molecule 1: Dibenzothiophene desulfurization enzyme C

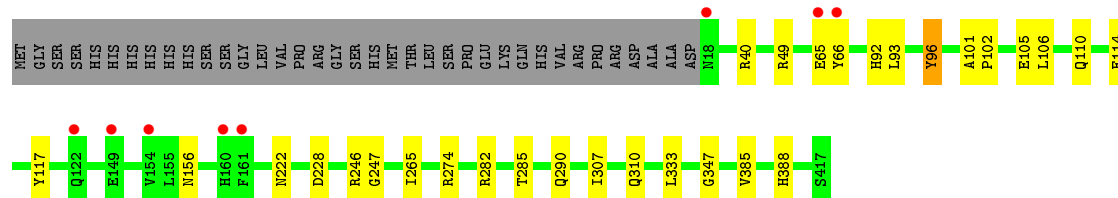
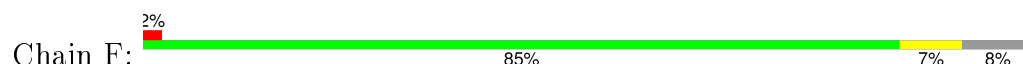




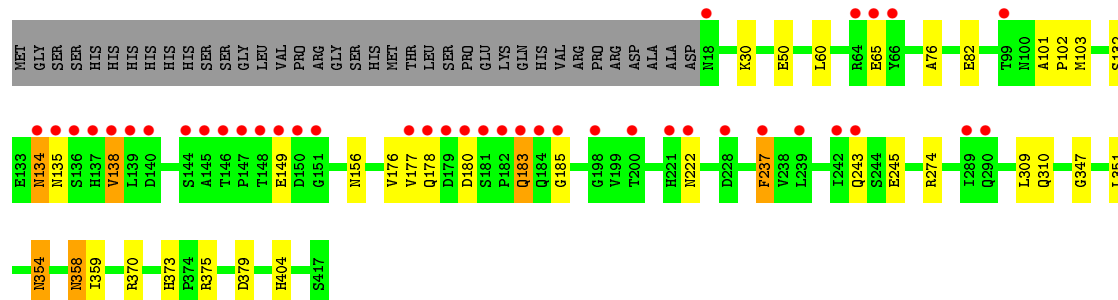
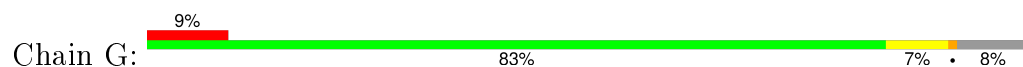
- Molecule 1: Dibenzothiophene desulfurization enzyme C



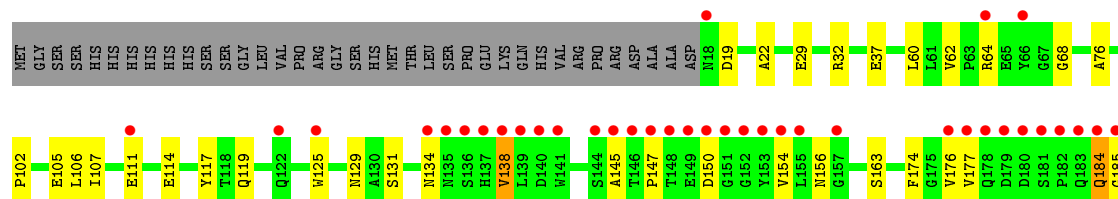
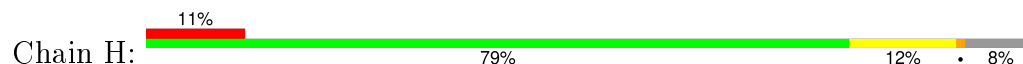
- Molecule 1: Dibenzothiophene desulfurization enzyme C

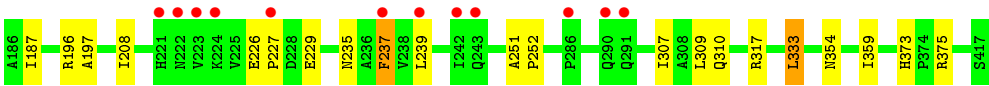


- Molecule 1: Dibenzothiophene desulfurization enzyme C



- Molecule 1: Dibenzothiophene desulfurization enzyme C





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	95.52Å 98.82Å 111.47Å 98.46° 106.99° 107.07°	Depositor
Resolution (Å)	35.42 – 1.79 35.42 – 1.79	Depositor EDS
% Data completeness (in resolution range)	89.7 (35.42-1.79) 83.7 (35.42-1.79)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.83 (at 1.79Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6.4_486)	Depositor
R, $R_{free}$	0.168 , 0.194 0.164 , 0.192	Depositor DCC
$R_{free}$ test set	15280 reflections (5.01%)	DCC
Wilson B-factor (Å <sup>2</sup> )	21.2	Xtriage
Anisotropy	0.225	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 44.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 324571 reflections	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	27568	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.05% 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL

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	0.36	0/3186	0.53	1/4344 (0.0%)
1	B	0.37	0/3195	0.52	0/4355
1	C	0.38	0/3159	0.53	0/4307
1	D	0.36	0/3174	0.51	0/4326
1	E	0.33	0/3182	0.48	0/4336
1	F	0.33	0/3179	0.48	0/4333
1	G	0.29	0/3167	0.43	0/4317
1	H	0.30	0/3141	0.45	0/4282
All	All	0.34	0/25383	0.49	1/34600 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	333	LEU	CA-CB-CG	5.12	127.08	115.30

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	3109	0	2962	21	0
1	B	3119	0	2978	33	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	3083	0	2948	25	0
1	D	3099	0	2965	28	0
1	E	3105	0	2966	31	0
1	F	3103	0	2965	22	0
1	G	3092	0	2956	31	0
1	H	3066	0	2939	32	0
2	A	6	0	8	1	0
2	B	12	0	16	0	0
2	C	6	0	8	0	0
2	D	6	0	5	1	0
2	E	12	0	16	0	0
2	F	6	0	8	0	0
2	G	12	0	16	0	0
2	H	12	0	16	0	0
3	A	413	0	0	7	0
3	B	384	0	0	10	0
3	C	432	0	0	10	0
3	D	387	0	0	7	0
3	E	343	0	0	10	0
3	F	310	0	0	4	0
3	G	227	0	0	9	0
3	H	224	0	0	4	0
All	All	27568	0	23772	207	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 (207) 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:D:135:ASN:HD21	1:F:285:THR:H	1.20	0.88
1:H:373:HIS:HD2	1:H:375:ARG:H	1.20	0.87
1:E:276:TYR:OH	1:E:370:ARG:HD2	1.75	0.87
1:G:373:HIS:HD2	1:G:375:ARG:H	1.25	0.85
1:E:274[B]:ARG:HG3	1:E:274[B]:ARG:HH11	1.42	0.82
1:F:49:ARG:NH1	3:F:843:HOH:O	2.12	0.82
1:E:107:ILE:HD13	1:E:237:PHE:HD2	1.47	0.80
1:B:373:HIS:HD2	1:B:375:ARG:H	1.27	0.79
1:B:226:GLU:OE1	3:B:827:HOH:O	2.03	0.76
1:E:328:ASP:OD1	3:E:831:HOH:O	2.03	0.75
1:E:274[B]:ARG:HH11	1:E:274[B]:ARG:CG	2.00	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:373:HIS:CD2	1:H:375:ARG:H	2.05	0.73
1:D:129:ASN:ND2	1:D:131:SER:OG	2.22	0.73
1:G:132:SER:OG	3:G:780:HOH:O	2.08	0.71
1:G:370:ARG:NH1	3:G:754:HOH:O	2.24	0.70
1:G:373:HIS:CD2	1:G:375:ARG:H	2.09	0.70
1:F:156:ASN:ND2	1:F:222:ASN:H	1.90	0.70
1:E:179:ASP:O	1:E:184:GLN:HG2	1.92	0.70
1:C:25:ARG:O	1:C:29:GLU:HG3	1.93	0.69
1:A:265:ILE:HG21	1:A:385[B]:VAL:HG13	1.75	0.69
1:D:354:ASN:ND2	3:D:700:HOH:O	2.24	0.68
1:E:132:SER:OG	3:E:887:HOH:O	2.12	0.68
1:C:53:ARG:NH1	3:C:865:HOH:O	2.26	0.67
1:C:344:LYS:HE2	3:C:950:HOH:O	1.94	0.67
1:C:265:ILE:HG21	1:C:385[A]:VAL:HG23	1.77	0.67
1:E:107:ILE:HD13	1:E:237:PHE:CD2	2.30	0.66
1:B:373:HIS:CD2	1:B:375:ARG:H	2.12	0.66
2:D:501:GOL:O1	2:D:502:GOL:C1	2.46	0.64
1:C:53:ARG:HD3	3:C:863:HOH:O	1.99	0.63
1:H:226:GLU:HB3	1:H:227:PRO:HD2	1.80	0.62
1:A:265:ILE:HG21	1:A:385[A]:VAL:HG23	1.81	0.62
1:A:92:HIS:HE1	3:A:707:HOH:O	1.83	0.62
1:B:310:GLN:HE22	1:C:347:GLY:HA3	1.65	0.62
1:D:179:ASP:O	1:D:184[B]:GLN:NE2	2.33	0.62
1:B:113[A]:GLU:CD	3:B:901:HOH:O	2.39	0.62
1:B:179:ASP:O	1:B:184:GLN:HG2	2.01	0.61
1:D:347:GLY:HA3	1:H:310:GLN:HE22	1.65	0.61
1:H:187:ILE:H	1:H:235:ASN:HD22	1.47	0.61
1:E:32:ARG:NH2	3:E:859:HOH:O	2.30	0.60
1:E:354:ASN:HB2	3:E:706:HOH:O	2.02	0.60
1:F:156:ASN:HD21	1:F:222:ASN:H	1.49	0.60
1:C:391:HIS:HD2	1:C:392:ASP:OD1	1.84	0.60
1:B:318[B]:GLU:OE2	3:B:626:HOH:O	2.16	0.60
1:G:243:GLN:HB2	1:G:245:GLU:HG3	1.84	0.59
1:D:135:ASN:ND2	1:F:285:THR:H	1.96	0.58
1:H:105:GLU:HG3	1:H:117:TYR:OH	2.03	0.58
1:H:60:LEU:HD21	1:H:76:ALA:HA	1.84	0.58
1:C:354:ASN:HB2	3:C:936:HOH:O	2.02	0.58
1:C:53:ARG:HD3	3:C:793:HOH:O	2.03	0.57
1:G:176:VAL:HG12	1:G:178:GLN:HG3	1.85	0.57
1:G:177:VAL:HG11	1:G:183:GLN:HB3	1.87	0.57
1:B:113[A]:GLU:HG3	1:B:117:TYR:CE2	2.39	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:60:LEU:HD21	1:E:76:ALA:HA	1.87	0.57
1:G:50:GLU:HG2	3:G:686:HOH:O	2.04	0.57
1:G:156:ASN:ND2	1:G:222:ASN:H	2.03	0.57
1:F:310:GLN:HE22	1:G:347:GLY:HA3	1.70	0.56
1:D:184[B]:GLN:HA	1:D:184[B]:GLN:HE21	1.70	0.56
1:C:391:HIS:HE1	3:C:794:HOH:O	1.89	0.56
1:G:358[A]:ASN:N	1:G:358[A]:ASN:HD22	2.03	0.56
1:D:391:HIS:HD2	1:D:392:ASP:OD1	1.89	0.56
1:B:53:ARG:NH2	3:B:772:HOH:O	2.39	0.55
1:A:310:GLN:HE22	1:E:347:GLY:HA3	1.71	0.55
1:E:156:ASN:ND2	1:E:222:ASN:H	2.03	0.55
1:A:30:LYS:NZ	3:A:826:HOH:O	2.39	0.55
1:F:66:TYR:CE1	1:F:114:GLU:HG2	2.43	0.54
1:G:138:VAL:HG13	1:G:176:VAL:HG11	1.90	0.54
1:G:404:HIS:HD2	3:G:679:HOH:O	1.90	0.54
1:H:107:ILE:HD13	1:H:237:PHE:CD2	2.44	0.53
1:A:403:LYS:NZ	3:A:903:HOH:O	2.18	0.53
1:H:145:ALA:O	1:H:177:VAL:HG22	2.08	0.53
1:F:347:GLY:HA3	1:G:310:GLN:HE22	1.73	0.53
1:C:105:GLU:OE2	3:C:790:HOH:O	2.18	0.53
1:B:181:SER:HB2	1:B:182:PRO:HD2	1.90	0.53
1:F:66:TYR:HE1	1:F:114:GLU:HG2	1.72	0.53
1:H:32:ARG:NH1	3:H:661:HOH:O	2.41	0.53
1:A:137[B]:HIS:HE1	1:A:139:LEU:HB2	1.75	0.52
1:G:379:ASP:HB3	1:H:208:ILE:HD13	1.91	0.52
1:A:113:GLU:HG3	1:A:117:TYR:CE2	2.45	0.52
1:D:18:ASN:N	1:D:18:ASN:HD22	2.08	0.51
1:G:354:ASN:HB2	3:G:651:HOH:O	2.10	0.51
1:C:208:ILE:HG13	1:C:209:GLY:N	2.25	0.51
1:A:274:ARG:NH2	3:A:951:HOH:O	2.43	0.51
1:B:135:ASN:N	1:B:135:ASN:OD1	2.43	0.51
1:E:106:LEU:HD22	1:E:246:ARG:HG2	1.93	0.51
1:H:235:ASN:O	1:H:239:LEU:HG	2.11	0.51
1:G:82:GLU:HG2	3:G:755:HOH:O	2.11	0.51
1:G:274:ARG:NH2	3:G:645:HOH:O	2.40	0.50
1:D:82:GLU:HG2	3:D:972:HOH:O	2.10	0.50
1:D:129:ASN:OD1	1:D:163:SER:HB2	2.12	0.50
1:B:291:GLN:HG2	3:B:925:HOH:O	2.12	0.50
1:D:265:ILE:HG21	1:D:385[B]:VAL:HG13	1.92	0.50
1:G:134:ASN:HD22	1:G:135:ASN:N	2.09	0.50
1:F:92:HIS:HE1	3:F:737:HOH:O	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:333:LEU:HD13	1:E:333:LEU:O	2.11	0.49
1:D:358[B]:ASN:OD1	3:D:914:HOH:O	2.20	0.49
1:F:282:ARG:NH2	3:F:635:HOH:O	2.18	0.49
1:D:28:ALA:HB1	1:D:86:ALA:HB2	1.94	0.49
1:H:129:ASN:HB3	1:H:163:SER:HB2	1.94	0.49
1:E:274[B]:ARG:NH1	1:E:274[B]:ARG:CG	2.68	0.49
1:D:391:HIS:HE1	3:D:832:HOH:O	1.95	0.49
1:C:160[A]:HIS:HE1	3:C:772:HOH:O	1.96	0.49
1:B:60:LEU:HD21	1:B:76:ALA:HA	1.93	0.49
1:A:160[A]:HIS:HE1	3:A:932:HOH:O	1.94	0.48
1:D:137:HIS:CE1	1:E:286:PRO:HA	2.48	0.48
1:C:32:ARG:HG3	1:C:33:ALA:N	2.27	0.48
1:A:137[B]:HIS:CE1	1:A:139:LEU:HB2	2.48	0.48
1:F:106:LEU:HB3	1:F:247:GLY:HA2	1.95	0.48
1:C:113:GLU:O	1:C:117:TYR:HB2	2.14	0.48
1:H:138:VAL:O	1:H:176:VAL:HG11	2.13	0.48
1:C:153:TYR:OH	1:C:183:GLN:NE2	2.41	0.47
1:C:262:TYR:OH	1:C:388:HIS:HD2	1.96	0.47
1:B:281:ALA:HB2	1:B:289:ILE:HD11	1.96	0.47
1:H:119:GLN:HB3	1:H:125:TRP:CZ3	2.50	0.47
1:A:159:LYS:HD3	3:A:677:HOH:O	2.15	0.47
1:C:181:SER:HB2	1:C:182:PRO:HD2	1.97	0.47
1:A:72:ASP:HA	2:A:500:GOL:H11	1.96	0.47
1:B:354[B]:ASN:ND2	1:B:358[B]:ASN:ND2	2.63	0.47
1:B:134:ASN:HB2	3:B:801:HOH:O	2.15	0.47
1:B:176:VAL:HG22	1:B:187:ILE:CD1	2.45	0.47
1:H:196:ARG:NH2	1:H:229:GLU:OE1	2.48	0.46
1:E:286:PRO:HG2	3:E:771:HOH:O	2.15	0.46
1:G:103:MET:HG3	1:G:237:PHE:CZ	2.50	0.46
1:B:19:ASP:OD1	3:B:787:HOH:O	2.21	0.46
1:H:64:ARG:HB3	1:H:64:ARG:HH11	1.81	0.46
1:F:92:HIS:CD2	1:F:388:HIS:NE2	2.83	0.46
1:G:183:GLN:HE21	1:G:183:GLN:HA	1.80	0.46
1:H:196:ARG:HG3	1:H:197:ALA:N	2.30	0.46
1:B:208:ILE:HD13	1:E:379:ASP:HB3	1.97	0.46
1:A:137[B]:HIS:CD2	1:A:140:ASP:OD2	2.69	0.45
1:A:24:ALA:HB2	1:A:79:VAL:HG13	1.99	0.45
1:D:129:ASN:HD22	1:D:131:SER:HB3	1.81	0.45
1:G:30:LYS:HE2	3:G:728:HOH:O	2.16	0.45
1:B:274:ARG:HD3	3:B:942:HOH:O	2.16	0.45
1:B:142:LYS:HG3	1:B:160:HIS:CE1	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:106:LEU:HD22	1:F:246:ARG:O	2.17	0.45
1:B:176:VAL:HG22	1:B:187:ILE:HD12	1.97	0.45
1:E:156:ASN:HD21	1:E:222:ASN:H	1.65	0.45
1:G:60:LEU:HD21	1:G:76:ALA:HA	1.98	0.45
1:D:351:LEU:HG	1:H:307:ILE:HG23	1.99	0.45
1:E:318[B]:GLU:HG2	3:E:615:HOH:O	2.16	0.45
1:A:233:ALA:HB2	1:H:29:GLU:HG3	1.97	0.45
1:H:317:ARG:HD3	3:H:648:HOH:O	2.15	0.45
1:E:246:ARG:HD3	3:E:929:HOH:O	2.18	0.44
1:H:163:SER:HA	3:H:736:HOH:O	2.16	0.44
1:A:354[A]:ASN:ND2	3:A:964:HOH:O	2.50	0.44
1:C:30:LYS:NZ	3:C:969:HOH:O	2.51	0.44
1:H:131:SER:HB2	1:H:174:PHE:CD2	2.52	0.44
1:F:274[B]:ARG:HD2	3:F:767:HOH:O	2.17	0.44
1:D:242:ILE:HG22	1:E:274[A]:ARG:HH12	1.82	0.44
1:H:251:ALA:HB3	1:H:252:PRO:HD3	2.00	0.44
1:B:21:VAL:HG12	1:B:25:ARG:NH1	2.33	0.44
1:D:96:TYR:HE2	1:D:163:SER:HG	1.64	0.43
1:B:113[A]:GLU:HG3	1:B:117:TYR:CD2	2.53	0.43
1:B:322:LEU:HA	1:B:325[A]:THR:HG22	1.99	0.43
1:E:274[B]:ARG:CB	1:E:274[B]:ARG:NH1	2.81	0.43
1:H:354:ASN:HB2	3:H:778:HOH:O	2.19	0.43
1:H:102:PRO:O	1:H:106:LEU:HG	2.18	0.43
1:D:140:ASP:OD2	3:D:923:HOH:O	2.21	0.43
1:D:159:LYS:HD3	3:D:661:HOH:O	2.18	0.43
1:D:129:ASN:HD22	1:D:131:SER:CB	2.31	0.43
1:G:103:MET:SD	1:G:237:PHE:HE2	2.42	0.43
1:B:101:ALA:N	1:B:102:PRO:CD	2.81	0.43
1:B:53:ARG:CZ	3:B:772:HOH:O	2.66	0.43
1:A:137[B]:HIS:NE2	1:A:140:ASP:OD2	2.52	0.43
1:C:34:THR:OG1	1:C:38:ARG:NH1	2.52	0.43
1:F:96:TYR:HD1	1:F:96:TYR:HA	1.66	0.43
1:D:358[B]:ASN:CG	3:D:914:HOH:O	2.57	0.43
1:B:30:LYS:NZ	3:B:792:HOH:O	2.45	0.42
1:B:49:ARG:NH2	1:B:93:LEU:HD11	2.34	0.42
1:H:19:ASP:OD2	1:H:22:ALA:HB2	2.19	0.42
1:C:96:TYR:HA	1:C:96:TYR:HD1	1.66	0.42
1:E:282:ARG:NH2	1:F:290:GLN:HE22	2.16	0.42
1:E:102:PRO:O	1:E:106:LEU:HG	2.19	0.42
1:C:181:SER:HB2	1:C:182:PRO:CD	2.49	0.42
1:B:307:ILE:HG23	1:C:351:LEU:HG	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:49:ARG:HH11	1:F:93:LEU:HD21	1.84	0.42
1:H:184:GLN:HG3	1:H:185:GLY:N	2.35	0.42
1:A:153:TYR:OH	1:A:183:GLN:NE2	2.43	0.42
1:B:309:LEU:HG	1:B:359:ILE:CD1	2.50	0.42
1:D:184[B]:GLN:HE21	1:D:184[B]:GLN:CA	2.30	0.42
1:G:370:ARG:HB3	3:G:637:HOH:O	2.20	0.41
1:F:105:GLU:HG3	1:F:117:TYR:OH	2.19	0.41
1:F:101:ALA:N	1:F:102:PRO:CD	2.83	0.41
1:D:356:ALA:HB1	1:D:385[B]:VAL:HG11	2.02	0.41
1:H:333:LEU:O	1:H:333:LEU:HD13	2.20	0.41
1:F:265:ILE:HG21	1:F:385[B]:VAL:HG13	2.02	0.41
1:C:53:ARG:NH2	3:C:863:HOH:O	2.52	0.41
1:A:370:ARG:NH2	1:C:161:PHE:HE1	2.18	0.41
1:E:134:ASN:N	3:E:814:HOH:O	2.53	0.41
1:H:154:VAL:HG12	1:H:156:ASN:ND2	2.35	0.41
1:E:274[B]:ARG:CB	1:E:274[B]:ARG:HH11	2.33	0.41
1:G:309:LEU:HG	1:G:359:ILE:CD1	2.51	0.41
1:H:309:LEU:HG	1:H:359:ILE:CD1	2.50	0.41
1:D:129:ASN:ND2	1:D:131:SER:CB	2.84	0.41
1:G:156:ASN:HD21	1:G:222:ASN:H	1.66	0.41
1:D:153:TYR:OH	1:D:183:GLN:NE2	2.40	0.41
1:A:208:ILE:HG13	1:A:209:GLY:N	2.35	0.41
1:F:307:ILE:HG23	1:G:351:LEU:HG	2.03	0.41
1:B:299:ILE:HA	1:B:299:ILE:HD13	1.91	0.41
1:B:138:VAL:HA	1:B:141:TRP:CD1	2.56	0.41
1:E:344:LYS:NZ	3:E:740:HOH:O	2.54	0.41
1:G:101:ALA:N	1:G:102:PRO:CD	2.85	0.40
1:G:138:VAL:HG11	1:G:185:GLY:O	2.20	0.40
1:G:134:ASN:HD22	1:G:135:ASN:H	1.68	0.40
1:E:101:ALA:HB1	1:E:117:TYR:HE1	1.87	0.40
1:E:318[B]:GLU:CG	3:E:615:HOH:O	2.70	0.40
1:H:62:VAL:O	1:H:68:GLY:HA3	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	
1	A	405/437 (93%)	401 (99%)	4 (1%)	0	100	100
1	B	406/437 (93%)	397 (98%)	9 (2%)	0	100	100
1	C	402/437 (92%)	397 (99%)	5 (1%)	0	100	100
1	D	404/437 (92%)	399 (99%)	5 (1%)	0	100	100
1	E	404/437 (92%)	396 (98%)	8 (2%)	0	100	100
1	F	404/437 (92%)	400 (99%)	4 (1%)	0	100	100
1	G	403/437 (92%)	390 (97%)	13 (3%)	0	100	100
1	H	400/437 (92%)	385 (96%)	14 (4%)	1 (0%)	46	28
All	All	3228/3496 (92%)	3165 (98%)	62 (2%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	H	147	PRO

### 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	315/340 (93%)	311 (99%)	4 (1%)	76	66
1	B	316/340 (93%)	308 (98%)	8 (2%)	55	37
1	C	312/340 (92%)	310 (99%)	2 (1%)	90	87
1	D	314/340 (92%)	307 (98%)	7 (2%)	60	42
1	E	314/340 (92%)	309 (98%)	5 (2%)	70	57
1	F	314/340 (92%)	308 (98%)	6 (2%)	65	49
1	G	313/340 (92%)	303 (97%)	10 (3%)	46	26
1	H	310/340 (91%)	301 (97%)	9 (3%)	50	31

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	2508/2720 (92%)	2457 (98%)	51 (2%)	63 47

All (51) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	96	TYR
1	A	228	ASP
1	A	237	PHE
1	A	333	LEU
1	B	40	ARG
1	B	110	GLN
1	B	134	ASN
1	B	135	ASN
1	B	149	GLU
1	B	184	GLN
1	B	228	ASP
1	B	237	PHE
1	C	96	TYR
1	C	149	GLU
1	D	18	ASN
1	D	40	ARG
1	D	96	TYR
1	D	129	ASN
1	D	135	ASN
1	D	228	ASP
1	D	280	GLN
1	E	184	GLN
1	E	237	PHE
1	E	291	GLN
1	E	333	LEU
1	E	370	ARG
1	F	40	ARG
1	F	65	GLU
1	F	96	TYR
1	F	110	GLN
1	F	228	ASP
1	F	333	LEU
1	G	65	GLU
1	G	134	ASN
1	G	138	VAL
1	G	149	GLU
1	G	180	ASP

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Mol	Chain	Res	Type
1	G	183	GLN
1	G	237	PHE
1	G	354	ASN
1	G	358[A]	ASN
1	G	358[B]	ASN
1	H	37	GLU
1	H	111	GLU
1	H	114	GLU
1	H	134	ASN
1	H	138	VAL
1	H	150	ASP
1	H	184	GLN
1	H	237	PHE
1	H	333	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (61) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	92	HIS
1	A	134	ASN
1	A	183	GLN
1	A	243	GLN
1	A	280	GLN
1	A	310	GLN
1	A	404	HIS
1	B	122	GLN
1	B	134	ASN
1	B	160	HIS
1	B	184	GLN
1	B	290	GLN
1	B	310	GLN
1	B	373	HIS
1	C	122	GLN
1	C	134	ASN
1	C	183	GLN
1	C	388	HIS
1	C	391	HIS
1	D	100	ASN
1	D	129	ASN
1	D	134	ASN
1	D	135	ASN
1	D	137	HIS

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Mol	Chain	Res	Type
1	D	183	GLN
1	D	391	HIS
1	E	124	ASN
1	E	156	ASN
1	E	202	ASN
1	E	243	GLN
1	E	290	GLN
1	E	354	ASN
1	F	92	HIS
1	F	110	GLN
1	F	122	GLN
1	F	134	ASN
1	F	156	ASN
1	F	183	GLN
1	F	280	GLN
1	F	290	GLN
1	F	310	GLN
1	G	100	ASN
1	G	119	GLN
1	G	122	GLN
1	G	129	ASN
1	G	134	ASN
1	G	156	ASN
1	G	183	GLN
1	G	202	ASN
1	G	310	GLN
1	G	373	HIS
1	G	404	HIS
1	H	100	ASN
1	H	122	GLN
1	H	134	ASN
1	H	202	ASN
1	H	222	ASN
1	H	235	ASN
1	H	243	GLN
1	H	310	GLN
1	H	373	HIS

### 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 13 ligands modelled in this entry, 1 is modelled with single atom - leaving 12 for Mogul analysis.

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	GOL	A	500	-	5,5,5	0.28	0	5,5,5	0.30	0
2	GOL	B	501	-	5,5,5	0.29	0	5,5,5	0.30	0
2	GOL	B	502	-	5,5,5	0.91	0	5,5,5	1.40	1 (20%)
2	GOL	C	500	-	5,5,5	0.29	0	5,5,5	0.24	0
2	GOL	D	502	-	4,4,5	0.76	0	2,4,5	2.15	1 (50%)
2	GOL	E	501	-	5,5,5	0.32	0	5,5,5	0.23	0
2	GOL	E	502	-	5,5,5	0.98	0	5,5,5	1.24	0
2	GOL	F	500	-	5,5,5	0.30	0	5,5,5	0.24	0
2	GOL	G	501	-	5,5,5	0.32	0	5,5,5	0.23	0
2	GOL	G	502	-	5,5,5	0.98	0	5,5,5	1.38	1 (20%)
2	GOL	H	501	-	5,5,5	0.33	0	5,5,5	0.32	0
2	GOL	H	502	-	5,5,5	0.97	0	5,5,5	1.39	0

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. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GOL	A	500	-	-	0/4/4/4	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GOL	B	501	-	-	0/4/4/4	0/0/0/0
2	GOL	B	502	-	-	0/4/4/4	0/0/0/0
2	GOL	C	500	-	-	0/4/4/4	0/0/0/0
2	GOL	D	502	-	-	0/2/2/4	0/0/0/0
2	GOL	E	501	-	-	0/4/4/4	0/0/0/0
2	GOL	E	502	-	-	0/4/4/4	0/0/0/0
2	GOL	F	500	-	-	0/4/4/4	0/0/0/0
2	GOL	G	501	-	-	0/4/4/4	0/0/0/0
2	GOL	G	502	-	-	0/4/4/4	0/0/0/0
2	GOL	H	501	-	-	0/4/4/4	0/0/0/0
2	GOL	H	502	-	-	0/4/4/4	0/0/0/0

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	502	GOL	O1-C1-C2	2.04	120.10	110.18
2	G	502	GOL	O3-C3-C2	2.07	120.22	110.18
2	D	502	GOL	O3-C3-C2	2.91	117.77	110.87

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	500	GOL	1	0
2	D	502	GOL	1	0

## 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	A	400/437 (91%)	-0.24	1 (0%) 94 93	14, 22, 35, 57	0
1	B	400/437 (91%)	-0.06	16 (4%) 42 40	13, 23, 48, 88	0
1	C	400/437 (91%)	-0.28	5 (1%) 79 79	13, 22, 36, 60	0
1	D	400/437 (91%)	-0.29	3 (0%) 87 87	14, 23, 40, 66	0
1	E	400/437 (91%)	-0.09	10 (2%) 61 59	14, 28, 49, 83	0
1	F	400/437 (91%)	-0.14	8 (2%) 68 68	16, 28, 43, 69	0
1	G	400/437 (91%)	0.42	40 (10%) 9 8	16, 36, 77, 99	0
1	H	400/437 (91%)	0.51	49 (12%) 5 5	15, 36, 83, 103	0
All	All	3200/3496 (91%)	-0.02	132 (4%) 41 39	13, 25, 60, 103	0

All (132) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	139	LEU	11.4
1	G	139	LEU	9.7
1	B	135	ASN	8.0
1	G	135	ASN	8.0
1	H	135	ASN	7.8
1	H	151	GLY	6.4
1	H	182	PRO	6.4
1	H	137	HIS	6.3
1	H	180	ASP	6.2
1	G	180	ASP	6.1
1	G	18	ASN	5.8
1	G	137	HIS	5.8
1	H	140	ASP	5.8
1	H	149	GLU	5.7
1	B	139	LEU	5.7
1	H	181	SER	5.6

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Mol	Chain	Res	Type	RSRZ
1	G	182	PRO	5.6
1	F	18	ASN	5.5
1	H	177	VAL	5.2
1	G	184	GLN	4.9
1	G	149	GLU	4.9
1	G	140	ASP	4.9
1	H	227	PRO	4.9
1	G	138	VAL	4.8
1	G	181	SER	4.8
1	B	180	ASP	4.7
1	H	134	ASN	4.7
1	H	154	VAL	4.6
1	G	290[A]	GLN	4.6
1	B	137	HIS	4.6
1	D	18	ASN	4.6
1	H	185	GLY	4.5
1	H	150	ASP	4.4
1	E	137[A]	HIS	4.2
1	G	151	GLY	4.2
1	H	176	VAL	4.2
1	B	290	GLN	4.2
1	H	179	ASP	4.1
1	E	135	ASN	4.1
1	E	180	ASP	4.0
1	G	179	ASP	4.0
1	G	178	GLN	4.0
1	H	183	GLN	3.9
1	H	178	GLN	3.8
1	H	18	ASN	3.8
1	H	152	GLY	3.8
1	G	134	ASN	3.7
1	H	221	HIS	3.7
1	G	183	GLN	3.7
1	G	221	HIS	3.6
1	G	65	GLU	3.5
1	B	138	VAL	3.5
1	H	237	PHE	3.5
1	G	147	PRO	3.5
1	H	290	GLN	3.4
1	B	134	ASN	3.4
1	H	138	VAL	3.4
1	G	146	THR	3.4

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Mol	Chain	Res	Type	RSRZ
1	H	242	ILE	3.4
1	G	185	GLY	3.4
1	E	134	ASN	3.4
1	B	140	ASP	3.3
1	H	136	SER	3.3
1	G	222	ASN	3.3
1	G	237	PHE	3.3
1	C	18	ASN	3.2
1	G	242	ILE	3.2
1	F	65	GLU	3.2
1	H	144	SER	3.2
1	H	122	GLN	3.2
1	B	281	ALA	3.2
1	B	179	ASP	3.1
1	G	148	THR	3.1
1	G	239	LEU	3.1
1	H	147	PRO	3.1
1	G	289	ILE	3.1
1	E	139	LEU	3.1
1	B	181	SER	3.1
1	E	18	ASN	3.1
1	H	222	ASN	3.1
1	H	184	GLN	3.0
1	C	180	ASP	3.0
1	H	286	PRO	3.0
1	H	153	TYR	3.0
1	H	291	GLN	3.0
1	B	18	ASN	3.0
1	G	136	SER	2.9
1	H	148	THR	2.9
1	H	155	LEU	2.8
1	E	290	GLN	2.8
1	G	198	GLY	2.7
1	H	223	VAL	2.7
1	C	149	GLU	2.7
1	F	161[A]	PHE	2.6
1	H	146	THR	2.6
1	H	125	TRP	2.6
1	F	66	TYR	2.6
1	H	239	LEU	2.6
1	B	284	TRP	2.6
1	G	243	GLN	2.5

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Mol	Chain	Res	Type	RSRZ
1	H	111	GLU	2.5
1	H	64	ARG	2.5
1	G	66	TYR	2.4
1	H	243	GLN	2.4
1	H	224	LYS	2.3
1	B	182	PRO	2.3
1	B	64[A]	ARG	2.3
1	E	242	ILE	2.3
1	G	150	ASP	2.3
1	D	96	TYR	2.3
1	B	132	SER	2.3
1	F	149	GLU	2.3
1	H	145	ALA	2.3
1	H	66	TYR	2.3
1	F	160	HIS	2.2
1	E	243	GLN	2.2
1	F	122	GLN	2.2
1	G	228	ASP	2.2
1	F	154	VAL	2.2
1	G	145	ALA	2.1
1	G	64	ARG	2.1
1	H	141	TRP	2.1
1	E	287	ALA	2.1
1	H	157	GLY	2.1
1	D	64	ARG	2.1
1	G	177	VAL	2.0
1	G	200	THR	2.0
1	A	180	ASP	2.0
1	C	172	PHE	2.0
1	G	144	SER	2.0
1	C	32	ARG	2.0
1	G	99	THR	2.0

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

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

## 6.3 Carbohydrates ⓘ

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(Å <sup>2</sup> )	Q<0.9
2	GOL	G	502	6/6	0.72	0.20	6.69	38,42,45,53	0
2	GOL	D	501	1/6	0.90	0.20	4.60	31,31,31,31	0
2	GOL	A	500	6/6	0.94	0.13	4.04	25,27,33,36	0
2	GOL	E	502	6/6	0.74	0.15	3.99	36,40,45,47	0
2	GOL	C	500	6/6	0.93	0.15	3.62	27,32,35,38	0
2	GOL	H	502	6/6	0.73	0.17	3.35	40,44,46,52	0
2	GOL	D	502	5/6	0.91	0.13	2.08	34,38,41,42	0
2	GOL	G	501	6/6	0.82	0.14	2.05	40,45,48,54	0
2	GOL	B	502	6/6	0.79	0.15	1.84	30,34,41,43	0
2	GOL	H	501	6/6	0.90	0.12	1.30	40,47,53,55	0
2	GOL	E	501	6/6	0.94	0.09	0.54	33,39,43,48	0
2	GOL	F	500	6/6	0.84	0.09	-0.05	35,38,43,44	0
2	GOL	B	501	6/6	0.95	0.07	-0.11	26,31,33,33	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.