



Full wwPDB X-ray Structure Validation Report ⓘ

May 26, 2020 – 11:25 am BST

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

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
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) : 2.11

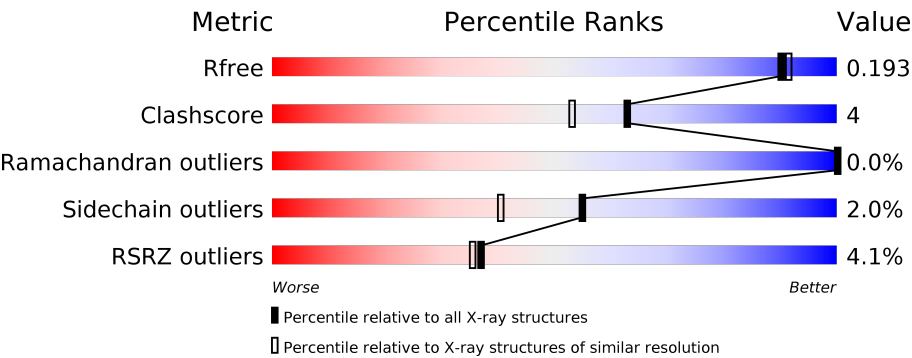
1 Overall quality at a glance i

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}	130704	9185 (1.80-1.76)
Clashscore	141614	10184 (1.80-1.76)
Ramachandran outliers	138981	10051 (1.80-1.76)
Sidechain outliers	138945	10050 (1.80-1.76)
RSRZ outliers	127900	9032 (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 ≥ 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	A	437	<div><div></div><div>85%6%8%</div></div>
1	B	437	<div>4%</div> <div><div></div><div>81%10%8%</div></div>
1	C	437	<div>%</div> <div><div></div><div>84%7%8%</div></div>
1	D	437	<div></div> <div><div></div><div>85%6%8%</div></div>
1	E	437	<div>2%</div> <div><div></div><div>85%6%8%</div></div>
1	F	437	<div>2%</div> <div><div></div><div>85%7%8%</div></div>

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Mol	Chain	Length	Quality of chain
1	G	437	<div><div></div><div>9%</div><div>83%</div><div>7%</div><div>8%</div></div>
1	H	437	<div><div></div><div>11%</div><div>79%</div><div>12%</div><div>8%</div></div>

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	-10	HIS	-	EXPRESSION TAG	UNP Q6WNP1
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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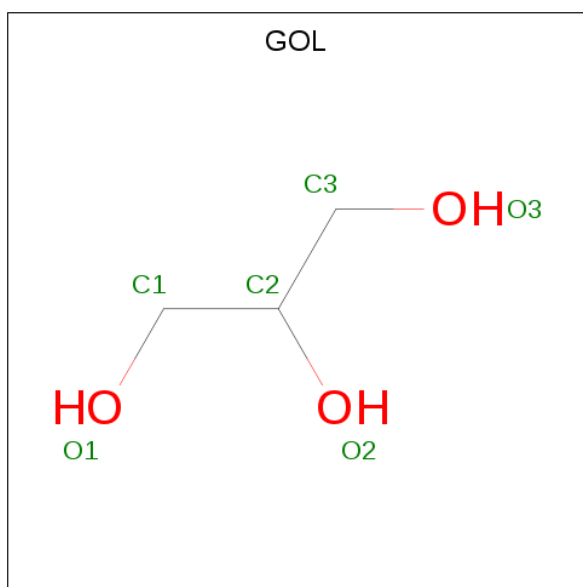
Chain	Residue	Modelled	Actual	Comment	Reference
E	-2	GLY	-	EXPRESSION TAG	UNP Q6WNP1
E	-1	SER	-	EXPRESSION TAG	UNP Q6WNP1
E	0	HIS	-	EXPRESSION TAG	UNP Q6WNP1
F	-19	MET	-	EXPRESSION TAG	UNP Q6WNP1
F	-18	GLY	-	EXPRESSION TAG	UNP Q6WNP1
F	-17	SER	-	EXPRESSION TAG	UNP Q6WNP1
F	-16	SER	-	EXPRESSION TAG	UNP 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: $C_3H_8O_3$).



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	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	G	1	Total C O 6 3 3	0	0
2	H	1	Total C O 6 3 3	0	0
2	H	1	Total C O 6 3 3	0	0

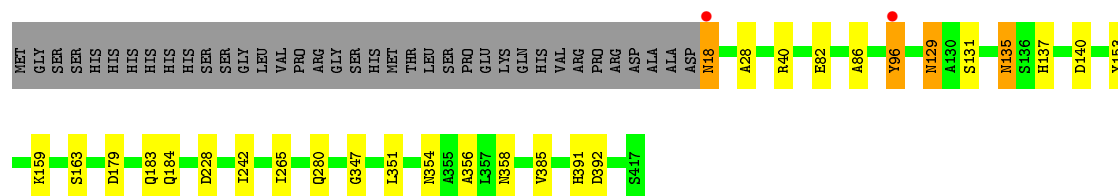
- Molecule 3 is water.

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

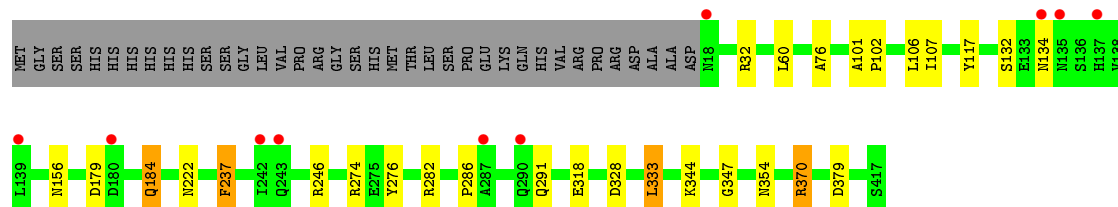
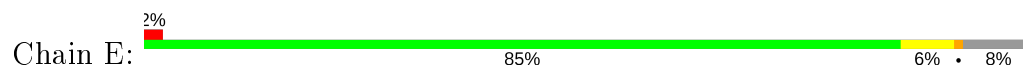
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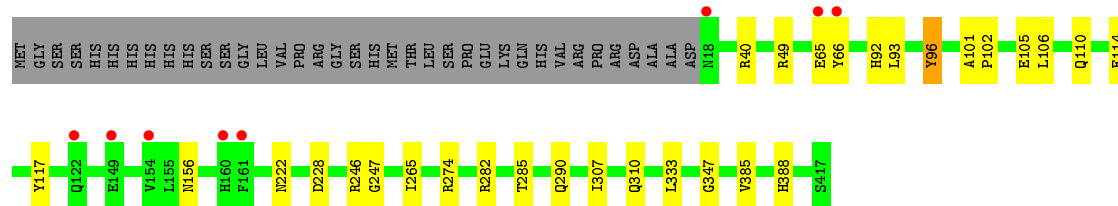
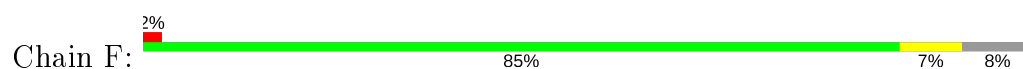
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	H	224	Total 224	O 224	0	0



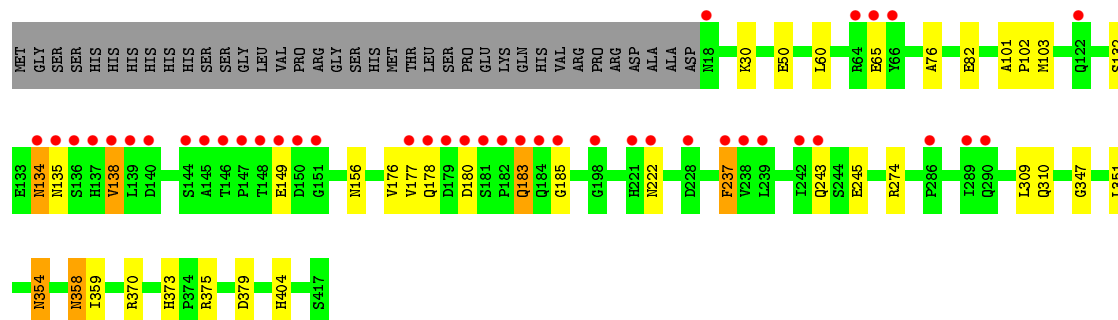
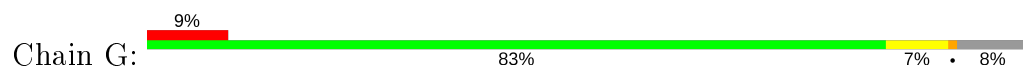
- Molecule 1: Dibenzothiophene desulfurization enzyme C



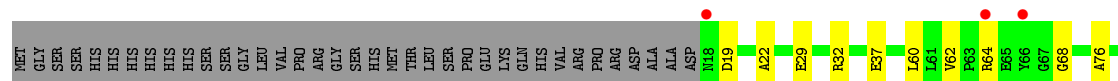
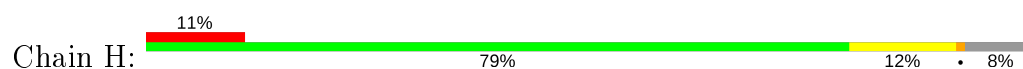
- Molecule 1: Dibenzothiophene desulfurization enzyme C

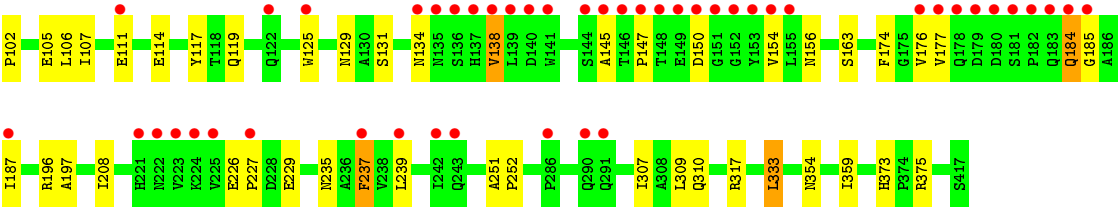


- Molecule 1: Dibenzothiophene desulfurization enzyme C



- Molecule 1: Dibenzothiophene desulfurization enzyme C





4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	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) 89.7 (35.42-1.79)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.83 (at 1.79Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6.4_486)	Depositor
R, R_{free}	0.168 , 0.194 0.167 , 0.193	Depositor DCC
R_{free} test set	16229 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	21.2	Xtriage
Anisotropy	0.225	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 44.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	27568	wwPDB-VP
Average B, all atoms (Å ²)	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.*

¹Intensities estimated from amplitudes.

²Theoretical 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: 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:B:373:HIS:CD2	1:B:375:ARG:H	2.12	0.66
1:E:107:ILE:HD13	1:E:237:PHE:CD2	2.30	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:B:113[A]:GLU:CD	3:B:901:HOH:O	2.39	0.62
1:D:179:ASP:O	1:D:184[B]:GLN:NE2	2.33	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:B:318[B]:GLU:OE2	3:B:626:HOH:O	2.16	0.60
1:C:391:HIS:HD2	1:C:392:ASP:OD1	1.84	0.60
1:F:156:ASN:HD21	1:F:222:ASN:H	1.49	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:156:ASN:ND2	1:G:222:ASN:H	2.03	0.57
1:G:50:GLU:HG2	3:G:686:HOH:O	2.04	0.57
1:F:310:GLN:HE22	1:G:347:GLY:HA3	1.70	0.56
1:C:391:HIS:HE1	3:C:794:HOH:O	1.89	0.56
1:D:184[B]:GLN:HA	1:D:184[B]:GLN:HE21	1.70	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:A:310:GLN:HE22	1:E:347:GLY:HA3	1.71	0.55
1:B:53:ARG:NH2	3:B:772:HOH:O	2.39	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:G:82:GLU:HG2	3:G:755:HOH:O	2.11	0.51
1:H:235:ASN:O	1:H:239:LEU:HG	2.11	0.51
1:D:82:GLU:HG2	3:D:972:HOH:O	2.10	0.50
1:G:274:ARG:NH2	3:G:645:HOH:O	2.40	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:F:92:HIS:HE1	3:F:737:HOH:O	1.93	0.50
1:G:134:ASN:HD22	1:G:135:ASN:N	2.09	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:358[B]:ASN:OD1	3:D:914:HOH:O	2.20	0.49
1:E:333:LEU:HD13	1:E:333:LEU:O	2.11	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:B:60:LEU:HD21	1:B:76:ALA:HA	1.93	0.49
1:C:160[A]:HIS:HE1	3:C:772:HOH:O	1.96	0.49
1:D:391:HIS:HE1	3:D:832:HOH:O	1.95	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:C:113:GLU:O	1:C:117:TYR:HB2	2.14	0.48
1:F:106:LEU:HB3	1:F:247:GLY:HA2	1.95	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:A:159:LYS:HD3	3:A:677:HOH:O	2.15	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:72:ASP:HA	2:A:500:GOL:H11	1.96	0.47
1:C:181:SER:HB2	1:C:182:PRO:HD2	1.97	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:B:208:ILE:HD13	1:E:379:ASP:HB3	1.97	0.46
1:H:196:ARG:HG3	1:H:197:ALA:N	2.30	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:142:LYS:HG3	1:B:160:HIS:CE1	2.51	0.45
1:B:274:ARG:HD3	3:B:942:HOH:O	2.16	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:A:233:ALA:HB2	1:H:29:GLU:HG3	1.97	0.45
1:E:318[B]:GLU:HG2	3:E:615:HOH:O	2.16	0.45
1:H:317:ARG:HD3	3:H:648:HOH:O	2.15	0.45
1:A:354[A]:ASN:ND2	3:A:964:HOH:O	2.50	0.44
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:C:30:LYS:NZ	3:C:969:HOH:O	2.51	0.44
1:F:274[B]:ARG:HD2	3:F:767:HOH:O	2.17	0.44
1:H:131:SER:HB2	1:H:174:PHE:CD2	2.52	0.44
1:D:242:ILE:HG22	1:E:274[A]:ARG:HH12	1.82	0.44
1:B:21:VAL:HG12	1:B:25:ARG:NH1	2.33	0.44
1:H:251:ALA:HB3	1:H:252:PRO:HD3	2.00	0.44
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:D:96:TYR:HE2	1:D:163:SER:HG	1.64	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:B:101:ALA:N	1:B:102:PRO:CD	2.81	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:A:137[B]:HIS:NE2	1:A:140:ASP:OD2	2.52	0.43
1:B:53:ARG:CZ	3:B:772:HOH:O	2.66	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:B:307:ILE:HG23	1:C:351:LEU:HG	2.01	0.42
1:C:181:SER:HB2	1:C:182:PRO:CD	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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: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:D:184[B]:GLN:HE21	1:D:184[B]:GLN:CA	2.30	0.42
1:F:105:GLU:HG3	1:F:117:TYR:OH	2.19	0.41
1:G:370:ARG:HB3	3:G:637:HOH:O	2.20	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:F:265:ILE:HG21	1:F:385[B]:VAL:HG13	2.02	0.41
1:H:333:LEU:O	1:H:333:LEU:HD13	2.20	0.41
1:A:370:ARG:NH2	1:C:161:PHE:HE1	2.18	0.41
1:C:53:ARG:NH2	3:C:863:HOH:O	2.52	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:A:208:ILE:HG13	1:A:209:GLY:N	2.35	0.41
1:D:129:ASN:ND2	1:D:131:SER:CB	2.84	0.41
1:D:153:TYR:OH	1:D:183:GLN:NE2	2.40	0.41
1:G:156:ASN:HD21	1:G:222:ASN:H	1.66	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:E:101:ALA:HB1	1:E:117:TYR:HE1	1.87	0.40
1:G:134:ASN:HD22	1:G:135:ASN:H	1.68	0.40
1:G:138:VAL:HG11	1:G:185:GLY:O	2.20	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%)	41	25
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%)	69	59
1	B	316/340 (93%)	308 (98%)	8 (2%)	47	31
1	C	312/340 (92%)	310 (99%)	2 (1%)	86	82
1	D	314/340 (92%)	307 (98%)	7 (2%)	52	36
1	E	314/340 (92%)	309 (98%)	5 (2%)	62	51
1	F	314/340 (92%)	308 (98%)	6 (2%)	57	43
1	G	313/340 (92%)	303 (97%)	10 (3%)	39	22
1	H	310/340 (91%)	301 (97%)	9 (3%)	42	25

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	2508/2720 (92%)	2457 (98%)	51 (2%)	55 40

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	B	502	-	5,5,5	0.87	0	5,5,5	1.39	1 (20%)
2	GOL	C	500	-	5,5,5	0.34	0	5,5,5	0.26	0
2	GOL	A	500	-	5,5,5	0.32	0	5,5,5	0.31	0
2	GOL	F	500	-	5,5,5	0.33	0	5,5,5	0.27	0
2	GOL	E	502	-	5,5,5	0.94	0	5,5,5	1.21	0
2	GOL	G	502	-	5,5,5	0.95	0	5,5,5	1.36	1 (20%)
2	GOL	G	501	-	5,5,5	0.35	0	5,5,5	0.23	0
2	GOL	H	501	-	5,5,5	0.37	0	5,5,5	0.33	0
2	GOL	E	501	-	5,5,5	0.34	0	5,5,5	0.26	0
2	GOL	H	502	-	5,5,5	0.93	0	5,5,5	1.36	0
2	GOL	B	501	-	5,5,5	0.33	0	5,5,5	0.34	0
2	GOL	D	502	-	3,4,5	0.84	0	1,4,5	1.05	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	B	502	-	-	2/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GOL	C	500	-	-	0/4/4/4	-
2	GOL	A	500	-	-	1/4/4/4	-
2	GOL	F	500	-	-	0/4/4/4	-
2	GOL	E	502	-	-	0/4/4/4	-
2	GOL	G	502	-	-	2/4/4/4	-
2	GOL	G	501	-	-	2/4/4/4	-
2	GOL	H	501	-	-	2/4/4/4	-
2	GOL	E	501	-	-	0/4/4/4	-
2	GOL	H	502	-	-	1/4/4/4	-
2	GOL	B	501	-	-	0/4/4/4	-
2	GOL	D	502	-	-	0/2/2/4	-

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	502	GOL	O3-C3-C2	2.09	120.22	110.20
2	B	502	GOL	O1-C1-C2	2.06	120.10	110.20

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	G	502	GOL	O1-C1-C2-C3
2	G	501	GOL	C1-C2-C3-O3
2	H	501	GOL	O1-C1-C2-C3
2	G	502	GOL	O1-C1-C2-O2
2	G	501	GOL	O2-C2-C3-O3
2	B	502	GOL	C1-C2-C3-O3
2	B	502	GOL	O2-C2-C3-O3
2	H	501	GOL	O1-C1-C2-O2
2	H	502	GOL	O1-C1-C2-C3
2	A	500	GOL	O1-C1-C2-C3

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, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	400/437 (91%)	-0.27	1 (0%) 94 93	14, 22, 35, 57	0
1	B	400/437 (91%)	-0.09	16 (4%) 38 36	13, 23, 48, 88	0
1	C	400/437 (91%)	-0.31	3 (0%) 86 86	13, 22, 36, 60	0
1	D	400/437 (91%)	-0.31	2 (0%) 91 91	14, 23, 40, 66	0
1	E	400/437 (91%)	-0.11	10 (2%) 57 56	14, 28, 49, 83	0
1	F	400/437 (91%)	-0.16	8 (2%) 65 65	16, 28, 43, 69	0
1	G	400/437 (91%)	0.40	41 (10%) 6 6	16, 36, 77, 99	0
1	H	400/437 (91%)	0.49	50 (12%) 3 3	15, 36, 83, 103	0
All	All	3200/3496 (91%)	-0.04	131 (4%) 37 35	13, 25, 60, 103	0

All (131) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	139	LEU	11.1
1	G	139	LEU	9.7
1	B	135	ASN	8.2
1	G	135	ASN	8.1
1	H	135	ASN	7.7
1	H	137	HIS	6.4
1	H	151	GLY	6.3
1	H	182	PRO	6.3
1	H	180	ASP	6.3
1	G	180	ASP	6.0
1	G	18	ASN	5.8
1	G	137	HIS	5.8
1	B	139	LEU	5.8
1	H	149	GLU	5.7
1	H	140	ASP	5.6
1	H	181	SER	5.6

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Mol	Chain	Res	Type	RSRZ
1	F	18	ASN	5.6
1	H	177	VAL	5.4
1	G	182	PRO	5.3
1	G	184	GLN	5.0
1	G	149	GLU	4.9
1	G	140	ASP	4.9
1	G	181	SER	4.8
1	B	180	ASP	4.8
1	G	138	VAL	4.7
1	H	227	PRO	4.7
1	B	137	HIS	4.6
1	H	154	VAL	4.5
1	H	134	ASN	4.5
1	G	290[A]	GLN	4.5
1	H	185	GLY	4.4
1	D	18	ASN	4.4
1	H	150	ASP	4.4
1	E	137[A]	HIS	4.3
1	G	151	GLY	4.2
1	E	180	ASP	4.1
1	B	290	GLN	4.1
1	E	135	ASN	4.1
1	G	178	GLN	4.0
1	H	176	VAL	4.0
1	H	179	ASP	4.0
1	G	179	ASP	4.0
1	H	183	GLN	3.9
1	H	178	GLN	3.9
1	H	152	GLY	3.8
1	H	18	ASN	3.8
1	H	221	HIS	3.7
1	G	183	GLN	3.7
1	G	134	ASN	3.6
1	G	221	HIS	3.5
1	G	146	THR	3.5
1	H	138	VAL	3.5
1	H	242	ILE	3.5
1	H	237	PHE	3.5
1	B	134	ASN	3.4
1	B	140	ASP	3.4
1	G	237	PHE	3.4
1	G	65	GLU	3.3

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Mol	Chain	Res	Type	RSRZ
1	H	290	GLN	3.3
1	G	222	ASN	3.3
1	G	185	GLY	3.3
1	G	147	PRO	3.3
1	H	136	SER	3.3
1	C	18	ASN	3.3
1	H	144	SER	3.2
1	E	134	ASN	3.2
1	B	138	VAL	3.2
1	F	65	GLU	3.2
1	H	122	GLN	3.2
1	H	147	PRO	3.2
1	G	148	THR	3.2
1	B	179	ASP	3.1
1	H	184	GLN	3.1
1	B	281	ALA	3.1
1	H	153	TYR	3.1
1	G	289	ILE	3.1
1	E	18	ASN	3.1
1	H	223	VAL	3.1
1	B	18	ASN	3.0
1	G	242	ILE	3.0
1	G	239	LEU	3.0
1	H	291	GLN	3.0
1	H	222	ASN	3.0
1	C	180	ASP	3.0
1	E	139	LEU	2.9
1	B	181	SER	2.9
1	H	286	PRO	2.9
1	H	148	THR	2.9
1	H	155	LEU	2.8
1	G	136	SER	2.8
1	C	149	GLU	2.6
1	F	161[A]	PHE	2.6
1	E	290	GLN	2.6
1	H	146	THR	2.6
1	F	66	TYR	2.6
1	G	243	GLN	2.6
1	G	198	GLY	2.5
1	H	243	GLN	2.5
1	H	239	LEU	2.5
1	H	64	ARG	2.5

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Mol	Chain	Res	Type	RSRZ
1	H	111	GLU	2.5
1	H	125	TRP	2.4
1	G	66	TYR	2.4
1	D	96	TYR	2.4
1	G	150	ASP	2.3
1	B	284	TRP	2.3
1	H	224	LYS	2.3
1	F	149	GLU	2.3
1	H	141	TRP	2.3
1	B	182	PRO	2.3
1	F	122	GLN	2.3
1	B	64[A]	ARG	2.2
1	H	145	ALA	2.2
1	E	243	GLN	2.2
1	F	160	HIS	2.2
1	G	145	ALA	2.2
1	B	132	SER	2.1
1	E	242	ILE	2.1
1	G	64	ARG	2.1
1	G	177	VAL	2.1
1	G	286	PRO	2.1
1	A	180	ASP	2.1
1	H	66	TYR	2.1
1	E	287	ALA	2.1
1	H	187	ILE	2.1
1	F	154	VAL	2.0
1	G	238	VAL	2.0
1	G	122	GLN	2.0
1	G	144	SER	2.0
1	H	225	VAL	2.0
1	G	228	ASP	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. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	GOL	H	502	6/6	0.70	0.18	40,44,46,52	0
2	GOL	G	502	6/6	0.72	0.20	38,42,45,53	0
2	GOL	E	502	6/6	0.73	0.15	36,40,45,47	0
2	GOL	B	502	6/6	0.79	0.15	30,34,41,43	0
2	GOL	G	501	6/6	0.81	0.14	40,45,48,54	0
2	GOL	F	500	6/6	0.85	0.10	35,38,43,44	0
2	GOL	H	501	6/6	0.90	0.12	40,47,53,55	0
2	GOL	D	502	5/6	0.91	0.12	34,38,41,42	0
2	GOL	D	501	1/6	0.93	0.18	31,31,31,31	0
2	GOL	C	500	6/6	0.93	0.16	27,32,35,38	0
2	GOL	B	501	6/6	0.94	0.07	26,31,33,33	0
2	GOL	E	501	6/6	0.94	0.08	33,39,43,48	0
2	GOL	A	500	6/6	0.95	0.12	25,27,33,36	0

6.5 Other polymers [i](#)

There are no such residues in this entry.