



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 01:36 AM UTC

PDB ID : 4DOY / pdb_00004doy
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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

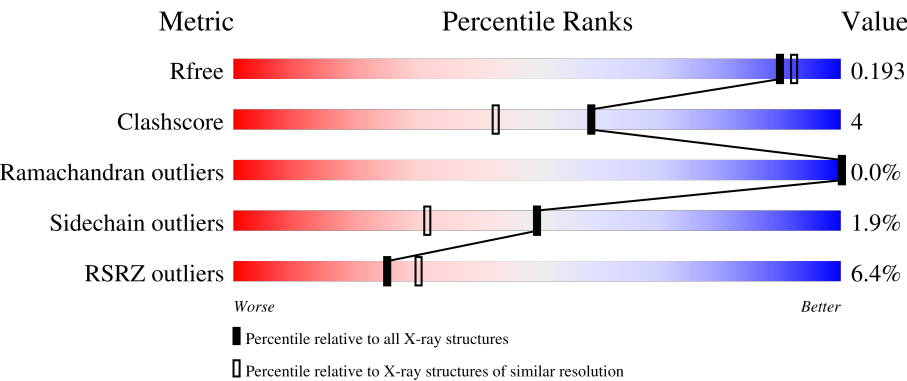
MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

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}	180053	1365 (1.78-1.78)
Clashscore	190562	1395 (1.78-1.78)
Ramachandran outliers	187476	1382 (1.78-1.78)
Sidechain outliers	187428	1382 (1.78-1.78)
RSRZ outliers	180081	1365 (1.78-1.78)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 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	0% 85%6%8%
1	B	437	4% 80%11%8%
1	C	437	84%7%8%
1	D	437	2% 85%6%8%
1	E	437	5% 85%6%8%

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Mol	Chain	Length	Quality of chain
1	F	437	% 84% 7% 8%
1	G	437	15% 82% 8% 8%
1	H	437	19% 79% 12% 8%

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
C	-12	HIS	-	expression tag	UNP Q6WNP1
C	-11	HIS	-	expression tag	UNP Q6WNP1
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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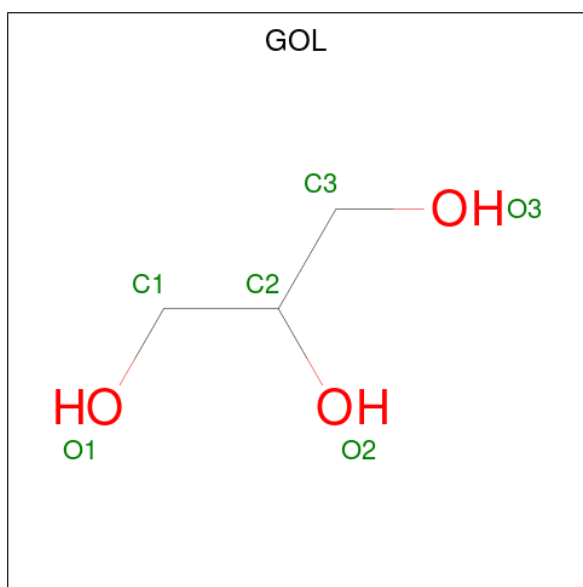
Chain	Residue	Modelled	Actual	Comment	Reference
E	-2	GLY	-	expression tag	UNP Q6WNP1
E	-1	SER	-	expression tag	UNP Q6WNP1
E	0	HIS	-	expression tag	UNP Q6WNP1
F	-19	MET	-	expression tag	UNP Q6WNP1
F	-18	GLY	-	expression tag	UNP Q6WNP1
F	-17	SER	-	expression tag	UNP Q6WNP1
F	-16	SER	-	expression tag	UNP 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 (CCD ID: 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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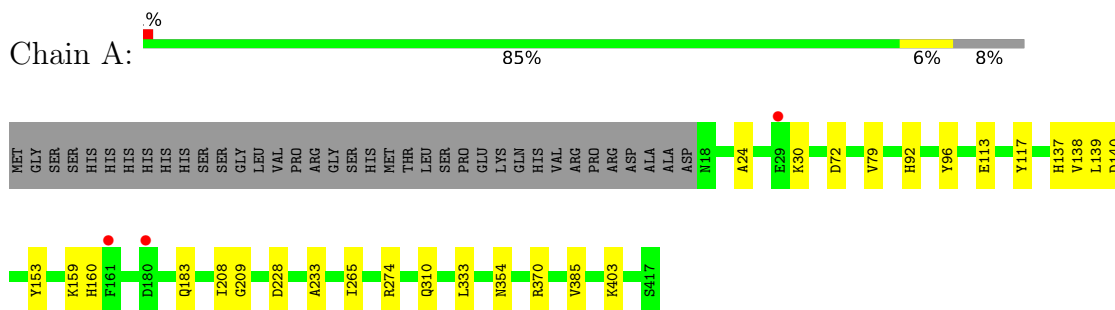
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	H	224	Total 224	O 224	0	0

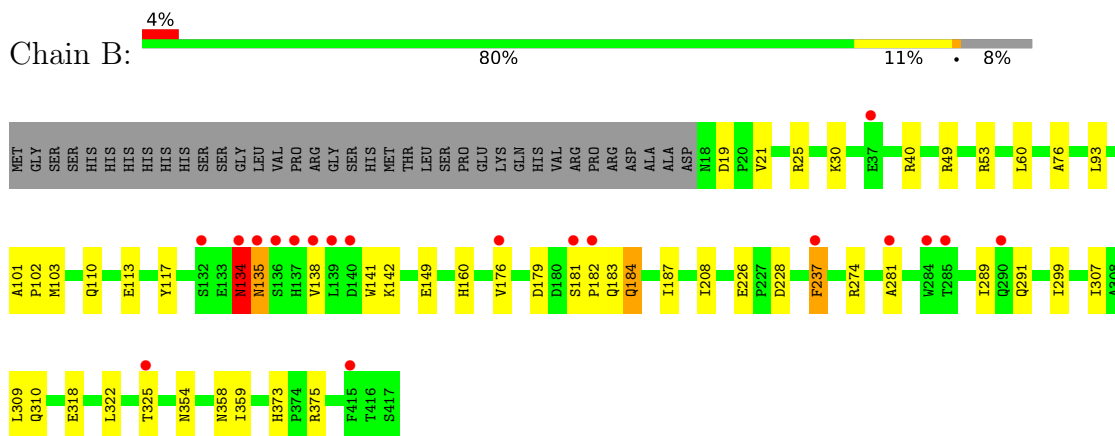
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

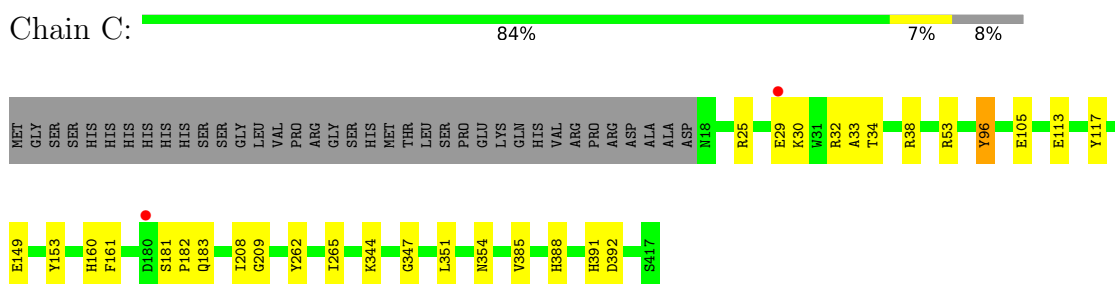
- Molecule 1: Dibenzothiophene desulfurization enzyme C



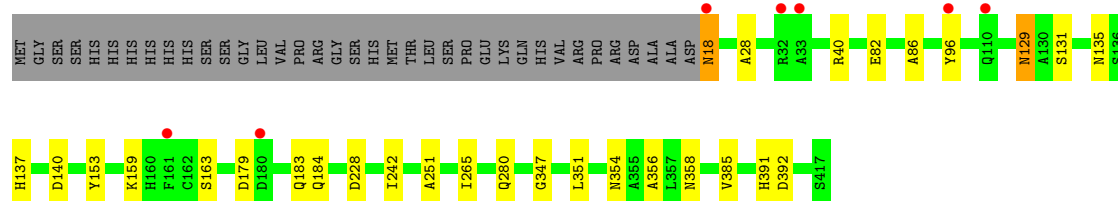
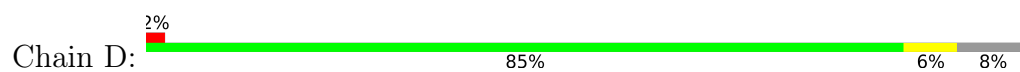
- Molecule 1: Dibenzothiophene desulfurization enzyme C



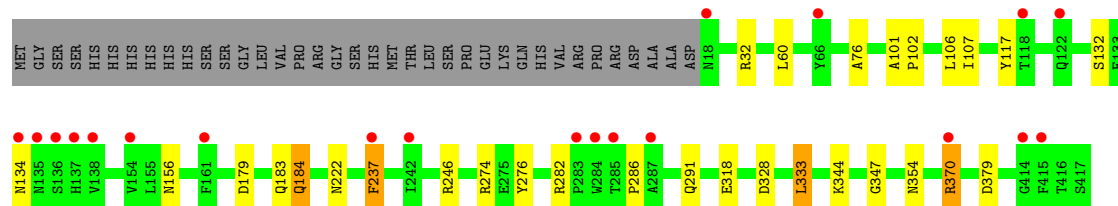
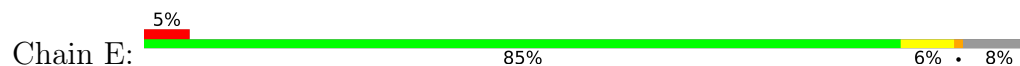
- Molecule 1: Dibenzothiophene desulfurization enzyme C



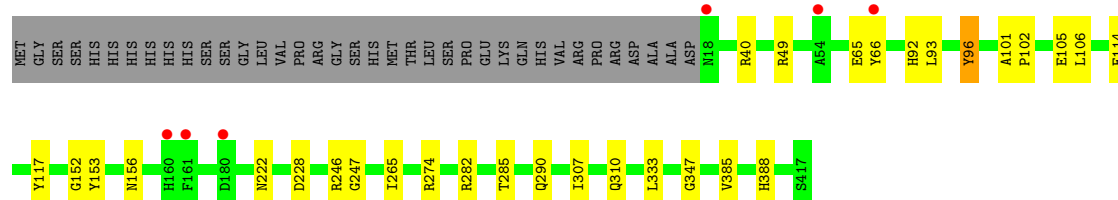
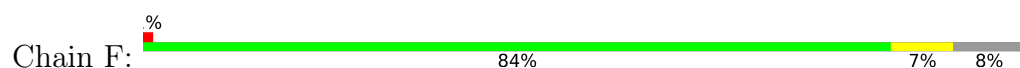
- Molecule 1: Dibenzothiophene desulfurization enzyme C



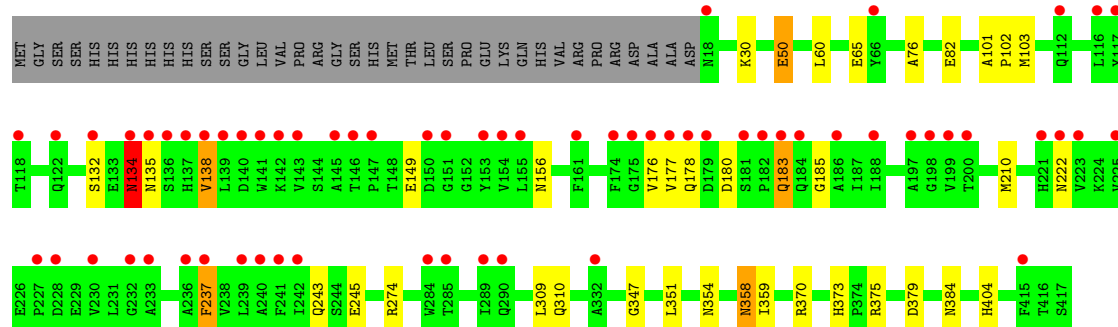
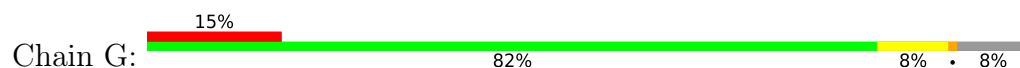
- Molecule 1: Dibenzothiophene desulfurization enzyme C



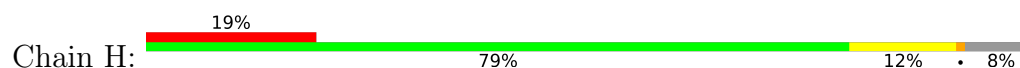
- Molecule 1: Dibenzothiophene desulfurization enzyme C

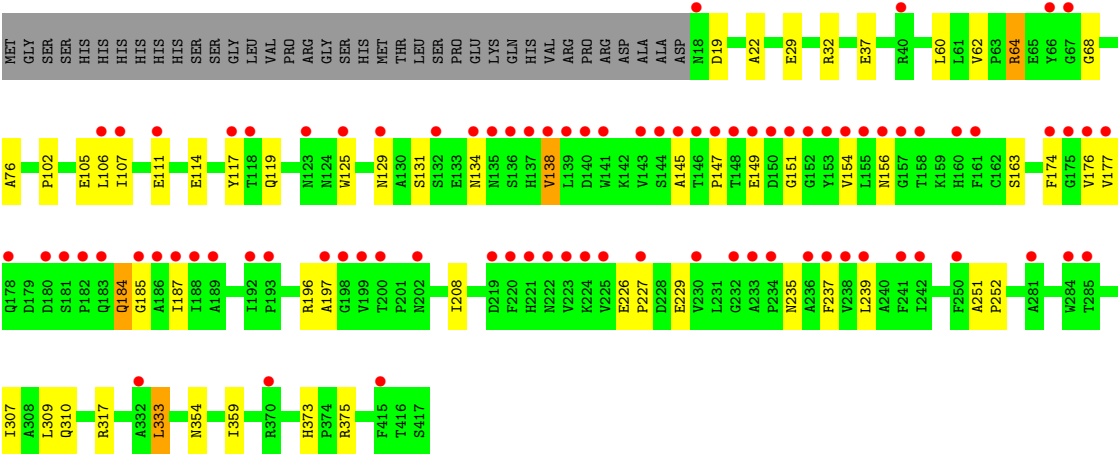


- Molecule 1: Dibenzothiophene desulfurization enzyme C



- Molecule 1: Dibenzothiophene desulfurization enzyme C





4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, 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 0.168 , 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 , 39.1	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

5.1 Standard geometry

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	0.46	0/3186	0.71	1/4344 (0.0%)
1	B	0.47	0/3195	0.73	1/4355 (0.0%)
1	C	0.48	0/3159	0.73	0/4307
1	D	0.46	0/3174	0.73	2/4326 (0.0%)
1	E	0.41	0/3182	0.67	0/4336
1	F	0.42	0/3179	0.69	0/4333
1	G	0.39	0/3167	0.68	1/4317 (0.0%)
1	H	0.38	0/3141	0.69	0/4282
All	All	0.44	0/25383	0.70	5/34600 (0.0%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	134	ASN	N-CA-C	10.63	123.87	111.11
1	D	251	ALA	CA-C-N	-5.79	112.62	119.05
1	D	251	ALA	C-N-CA	-5.79	112.62	119.05
1	A	138	VAL	N-CA-C	5.73	116.46	110.62
1	G	134	ASN	CB-CA-C	-5.27	110.48	116.54

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3109	0	2962	22	0
1	B	3119	0	2978	35	0
1	C	3083	0	2948	25	0
1	D	3099	0	2965	28	0
1	E	3105	0	2966	32	0
1	F	3103	0	2965	23	0
1	G	3092	0	2956	32	0
1	H	3066	0	2939	33	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	214	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 (214) 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:HH11	1:E:274[B]:ARG:HG3	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:274[B]:ARG:HH11	1:E:274[B]:ARG:CG	2.00	0.74
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:A:403:LYS:NZ	3:A:903:HOH:O	2.18	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:113[A]:GLU:CD	3:B:901:HOH:O	2.39	0.66
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: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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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:C:391:HIS:HE1	3:C:794:HOH:O	1.89	0.56
1:D:184[B]:GLN:HE21	1:D:184[B]:GLN:HA	1.70	0.56
1:G:358[A]:ASN:HD22	1:G:358[A]:ASN:N	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: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:HD22	1:D:18:ASN:N	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:OD1	1:B:135:ASN:N	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:134:ASN:HD22	1:G:135:ASN:N	2.09	0.50
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:CG	1:E:274[B]:ARG:NH1	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:D:358[B]:ASN:CG	3:D:914:HOH:O	2.57	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:HH11	1:H:64:ARG:HB3	1.81	0.46
1:F:92:HIS:CD2	1:F:388:HIS:NE2	2.83	0.46
1:G:183:GLN:HA	1:G:183:GLN:HE21	1.80	0.46
1:B:208:ILE:HD13	1:E:379:ASP:HB3	1.97	0.46
1:A:24:ALA:HB2	1:A:79:VAL:HG13	1.99	0.45
1:A:137[B]:HIS:CD2	1:A:140:ASP:OD2	2.69	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:274:ARG:HD3	3:B:942:HOH:O	2.16	0.45
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:H:196:ARG:HG3	1:H:197:ALA:N	2.30	0.45
1:D:351:LEU:HG	1:H:307:ILE:HG23	1.99	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: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:G:210:MET:HE2	1:G:384:ASN:HB3	1.99	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:E:274[B]:ARG:NH1	1:E:274[B]:ARG:CB	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:E:183:GLN:O	1:E:184:GLN:C	2.61	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: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:H:149:GLU:C	1:H:151:GLY:H	2.27	0.42
1:C:96:TYR:HD1	1:C:96:TYR:HA	1.66	0.42
1:E:282:ARG:NH2	1:F:290:GLN:HE22	2.16	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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:B:183:GLN:O	1:B:184:GLN:C	2.63	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:D:179:ASP:OD1	1:D:179:ASP:C	2.63	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:HH11	1:E:274[B]:ARG:CB	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:F:307:ILE:HG23	1:G:351:LEU:HG	2.03	0.41
1:G:156:ASN:HD21	1:G:222:ASN:H	1.66	0.41
1:B:138:VAL:HA	1:B:141:TRP:CD1	2.56	0.41
1:B:299:ILE:HD13	1:B:299:ILE:HA	1.91	0.41
1:E:344:LYS:NZ	3:E:740:HOH:O	2.54	0.41
1:A:137[B]:HIS:ND1	1:A:137[B]:HIS:C	2.78	0.40
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:F:152:GLY:C	1:F:153:TYR:CD1	2.99	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:B:103:MET:HE3	1:B:237:PHE:HE2	1.85	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

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	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%)	36	22
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%)	312 (99%)	3 (1%)	68	55
1	B	316/340 (93%)	308 (98%)	8 (2%)	42	22
1	C	312/340 (92%)	310 (99%)	2 (1%)	78	70
1	D	314/340 (92%)	308 (98%)	6 (2%)	50	32

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	E	314/340 (92%)	309 (98%)	5 (2%)	55	38
1	F	314/340 (92%)	309 (98%)	5 (2%)	55	38
1	G	313/340 (92%)	303 (97%)	10 (3%)	34	13
1	H	310/340 (91%)	302 (97%)	8 (3%)	40	20
All	All	2508/2720 (92%)	2461 (98%)	47 (2%)	50	32

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	96	TYR
1	A	228	ASP
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	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	228	ASP
1	F	333	LEU
1	G	50	GLU
1	G	65	GLU
1	G	134	ASN

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Mol	Chain	Res	Type
1	G	138	VAL
1	G	149	GLU
1	G	180	ASP
1	G	183	GLN
1	G	237	PHE
1	G	358[A]	ASN
1	G	358[B]	ASN
1	H	37	GLU
1	H	64	ARG
1	H	111	GLU
1	H	114	GLU
1	H	134	ASN
1	H	138	VAL
1	H	184	GLN
1	H	333	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (66) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	92	HIS
1	A	124	ASN
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	134	ASN
1	B	160	HIS
1	B	183	GLN
1	B	184	GLN
1	B	243	GLN
1	B	290	GLN
1	B	310	GLN
1	B	373	HIS
1	C	115	HIS
1	C	122	GLN
1	C	134	ASN
1	C	183	GLN
1	C	388	HIS
1	C	391	HIS
1	D	129	ASN

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Mol	Chain	Res	Type
1	D	134	ASN
1	D	135	ASN
1	D	137	HIS
1	D	183	GLN
1	D	221	HIS
1	D	290	GLN
1	D	391	HIS
1	E	156	ASN
1	E	184	GLN
1	E	202	ASN
1	E	243	GLN
1	E	290	GLN
1	F	18	ASN
1	F	92	HIS
1	F	110	GLN
1	F	122	GLN
1	F	134	ASN
1	F	156	ASN
1	F	202	ASN
1	F	221	HIS
1	F	243	GLN
1	F	280	GLN
1	F	290	GLN
1	F	310	GLN
1	G	100	ASN
1	G	115	HIS
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	310	GLN
1	G	373	HIS
1	G	404	HIS
1	H	100	ASN
1	H	122	GLN
1	H	134	ASN
1	H	221	HIS
1	H	235	ASN
1	H	243	GLN
1	H	310	GLN

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Mol	Chain	Res	Type
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 ⓘ

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

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	H	501	-	5,5,5	0.37	0	5,5,5	0.35	0
2	GOL	A	500	-	5,5,5	0.32	0	5,5,5	0.36	0
2	GOL	H	502	-	5,5,5	0.95	0	5,5,5	1.42	0
2	GOL	D	502	-	4,4,5	0.79	0	4,4,5	0.63	0
2	GOL	B	502	-	5,5,5	0.89	0	5,5,5	1.45	2 (40%)
2	GOL	E	501	-	5,5,5	0.35	0	5,5,5	0.29	0
2	GOL	F	500	-	5,5,5	0.33	0	5,5,5	0.30	0
2	GOL	E	502	-	5,5,5	0.96	0	5,5,5	1.26	0
2	GOL	G	502	-	5,5,5	0.97	0	5,5,5	1.42	1 (20%)
2	GOL	C	500	-	5,5,5	0.35	0	5,5,5	0.28	0
2	GOL	G	501	-	5,5,5	0.35	0	5,5,5	0.25	0
2	GOL	B	501	-	5,5,5	0.33	0	5,5,5	0.37	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	H	501	-	-	2/4/4/4	-
2	GOL	A	500	-	-	0/4/4/4	-
2	GOL	H	502	-	-	1/4/4/4	-
2	GOL	D	502	-	-	0/2/2/4	-
2	GOL	B	502	-	-	2/4/4/4	-
2	GOL	E	501	-	-	0/4/4/4	-
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	C	500	-	-	0/4/4/4	-
2	GOL	G	501	-	-	2/4/4/4	-
2	GOL	B	501	-	-	0/4/4/4	-

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	502	GOL	O3-C3-C2	2.19	120.22	110.38
2	B	502	GOL	O1-C1-C2	2.16	120.10	110.38
2	B	502	GOL	O3-C3-C2	2.08	119.73	110.38

There are no chirality outliers.

All (9) torsion outliers are listed below:

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

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%)	-0.12	3 (0%) 82 87	8, 22, 35, 57	7 (1%)
1	B	400/437 (91%)	0.12	19 (4%) 35 41	8, 22, 48, 88	8 (2%)
1	C	400/437 (91%)	-0.11	2 (0%) 87 91	7, 22, 36, 60	4 (1%)
1	D	400/437 (91%)	-0.02	7 (1%) 67 75	9, 23, 40, 66	6 (1%)
1	E	400/437 (91%)	0.34	20 (5%) 34 40	7, 28, 49, 83	6 (1%)
1	F	400/437 (91%)	0.24	6 (1%) 72 78	9, 28, 43, 69	6 (1%)
1	G	400/437 (91%)	0.80	64 (16%) 5 4	9, 36, 77, 99	5 (1%)
1	H	400/437 (91%)	0.96	85 (21%) 2 2	9, 36, 83, 103	2 (0%)
All	All	3200/3496 (91%)	0.28	206 (6%) 25 30	7, 25, 60, 103	44 (1%)

All (206) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	155	LEU	5.2
1	H	157	GLY	5.1
1	H	223	VAL	5.0
1	H	143	VAL	4.9
1	H	237	PHE	4.8
1	H	181	SER	4.7
1	H	177	VAL	4.6
1	H	144	SER	4.5
1	F	161[A]	PHE	4.4
1	H	154	VAL	4.4
1	H	145	ALA	4.3
1	H	151	GLY	4.3
1	H	138	VAL	4.3
1	E	138	VAL	4.2
1	G	139	LEU	4.2
1	G	237	PHE	4.2

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Mol	Chain	Res	Type	RSRZ
1	B	137	HIS	4.1
1	H	161	PHE	4.1
1	G	182	PRO	4.0
1	G	138	VAL	4.0
1	B	135	ASN	3.9
1	H	225	VAL	3.9
1	H	148	THR	3.9
1	H	227	PRO	3.9
1	H	182	PRO	3.8
1	H	139	LEU	3.8
1	B	132	SER	3.7
1	H	176	VAL	3.7
1	H	152	GLY	3.6
1	H	153	TYR	3.6
1	G	135	ASN	3.6
1	H	137	HIS	3.6
1	H	136	SER	3.5
1	H	146	THR	3.5
1	G	284	TRP	3.5
1	G	154	VAL	3.5
1	G	137	HIS	3.5
1	A	161	PHE	3.5
1	B	284	TRP	3.5
1	H	147	PRO	3.4
1	E	134	ASN	3.4
1	H	221	HIS	3.4
1	H	239	LEU	3.4
1	E	284	TRP	3.4
1	A	180	ASP	3.3
1	H	135	ASN	3.3
1	G	239	LEU	3.3
1	G	177	VAL	3.3
1	H	222	ASN	3.3
1	H	158	THR	3.3
1	H	106	LEU	3.3
1	E	415	PHE	3.3
1	B	281	ALA	3.3
1	E	18	ASN	3.3
1	H	156	ASN	3.2
1	B	138	VAL	3.2
1	H	220	PHE	3.2
1	G	198	GLY	3.2

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Mol	Chain	Res	Type	RSRZ
1	H	238	VAL	3.2
1	G	227	PRO	3.2
1	G	175	GLY	3.2
1	E	137[A]	HIS	3.2
1	H	200	THR	3.2
1	G	136	SER	3.1
1	H	186	ALA	3.1
1	B	182	PRO	3.1
1	H	187	ILE	3.1
1	G	188	ILE	3.0
1	H	192	ILE	3.0
1	D	33	ALA	3.0
1	B	285	THR	2.9
1	E	287	ALA	2.9
1	H	134	ASN	2.9
1	B	415	PHE	2.9
1	G	228	ASP	2.9
1	G	181	SER	2.9
1	B	134	ASN	2.9
1	G	116	LEU	2.9
1	G	151	GLY	2.9
1	H	125	TRP	2.9
1	G	184	GLN	2.9
1	G	161	PHE	2.9
1	H	66	TYR	2.9
1	H	107	ILE	2.9
1	G	155	LEU	2.8
1	G	233	ALA	2.8
1	B	140	ASP	2.8
1	G	141	TRP	2.8
1	H	141	TRP	2.8
1	E	161	PHE	2.8
1	H	149	GLU	2.8
1	H	224	LYS	2.8
1	H	67	GLY	2.8
1	G	176	VAL	2.7
1	G	118	THR	2.7
1	E	135	ASN	2.7
1	H	236	ALA	2.7
1	D	180	ASP	2.7
1	E	285	THR	2.7
1	H	198	GLY	2.7

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Mol	Chain	Res	Type	RSRZ
1	D	18	ASN	2.7
1	G	18	ASN	2.7
1	C	180	ASP	2.7
1	H	132	SER	2.7
1	G	290[A]	GLN	2.6
1	F	18	ASN	2.6
1	F	160	HIS	2.6
1	H	234	PRO	2.6
1	G	132	SER	2.6
1	E	237	PHE	2.6
1	G	415	PHE	2.6
1	H	415	PHE	2.6
1	C	29	GLU	2.6
1	H	199	VAL	2.6
1	G	178	GLN	2.6
1	G	183	GLN	2.6
1	G	289	ILE	2.6
1	H	160	HIS	2.6
1	G	197	ALA	2.6
1	H	250	PHE	2.6
1	B	176	VAL	2.6
1	G	225	VAL	2.6
1	E	136	SER	2.6
1	H	242	ILE	2.6
1	F	66	TYR	2.5
1	H	281	ALA	2.5
1	B	136	SER	2.5
1	G	66	TYR	2.5
1	G	153	TYR	2.5
1	G	332	ALA	2.5
1	E	414	GLY	2.5
1	G	285	THR	2.5
1	B	237	PHE	2.5
1	E	66	TYR	2.5
1	G	223	VAL	2.5
1	G	179	ASP	2.4
1	H	233	ALA	2.4
1	H	188	ILE	2.4
1	H	178	GLN	2.4
1	H	183	GLN	2.4
1	H	150	ASP	2.4
1	B	325[A]	THR	2.4

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Mol	Chain	Res	Type	RSRZ
1	G	134	ASN	2.4
1	H	193	PRO	2.4
1	H	118	THR	2.4
1	H	284	TRP	2.4
1	G	150	ASP	2.4
1	H	111	GLU	2.3
1	F	180	ASP	2.3
1	H	332	ALA	2.3
1	H	174	PHE	2.3
1	H	117	TYR	2.3
1	H	40	ARG	2.3
1	G	240	ALA	2.3
1	G	117	TYR	2.3
1	B	181	SER	2.3
1	G	230	VAL	2.3
1	B	290	GLN	2.3
1	D	110	GLN	2.3
1	G	112	GLN	2.3
1	H	129	ASN	2.3
1	D	32	ARG	2.3
1	G	232	GLY	2.3
1	H	219	ASP	2.3
1	G	145	ALA	2.2
1	H	189	ALA	2.2
1	H	197	ALA	2.2
1	E	370	ARG	2.2
1	G	241	PHE	2.2
1	H	285	THR	2.2
1	H	202	ASN	2.2
1	G	142	LYS	2.2
1	H	232	GLY	2.2
1	A	29	GLU	2.2
1	G	174	PHE	2.2
1	G	221	HIS	2.2
1	G	143	VAL	2.2
1	H	180	ASP	2.2
1	H	185	GLY	2.2
1	H	140	ASP	2.1
1	E	154	VAL	2.1
1	H	230	VAL	2.1
1	G	122	GLN	2.1
1	H	370	ARG	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	139	LEU	2.1
1	E	122[A]	GLN	2.1
1	F	54	ALA	2.1
1	G	186	ALA	2.1
1	H	123	ASN	2.1
1	G	140	ASP	2.1
1	E	118	THR	2.1
1	G	200	THR	2.1
1	E	283	PRO	2.1
1	D	161	PHE	2.1
1	H	241	PHE	2.1
1	G	242	ILE	2.1
1	G	146	THR	2.1
1	G	147	PRO	2.1
1	H	175	GLY	2.0
1	G	222	ASN	2.0
1	G	236	ALA	2.0
1	H	18	ASN	2.0
1	B	37	GLU	2.0
1	D	96	TYR	2.0
1	E	242	ILE	2.0
1	G	199	VAL	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides 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($\text{\AA}^2$)	Q<0.9
2	GOL	G	502	6/6	0.65	0.23	38,42,45,53	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	GOL	E	502	6/6	0.74	0.16	36,40,45,47	0
2	GOL	B	502	6/6	0.74	0.17	30,34,41,43	0
2	GOL	D	501	1/6	0.77	0.16	31,31,31,31	0
2	GOL	H	502	6/6	0.79	0.15	40,44,46,52	0
2	GOL	H	501	6/6	0.82	0.14	40,47,53,55	0
2	GOL	G	501	6/6	0.84	0.12	40,45,48,54	0
2	GOL	D	502	5/6	0.87	0.12	34,38,41,42	0
2	GOL	F	500	6/6	0.90	0.11	35,38,43,44	0
2	GOL	E	501	6/6	0.92	0.10	33,39,43,48	0
2	GOL	C	500	6/6	0.93	0.12	27,32,35,38	0
2	GOL	B	501	6/6	0.93	0.09	26,31,33,33	0
2	GOL	A	500	6/6	0.94	0.10	25,27,33,36	0

6.5 Other polymers [i](#)

There are no such residues in this entry.