



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 12:57 PM UTC

PDB ID : 5HUD / pdb_00005hud
Title : Non-covalent complex of and DAHP synthase and chorismate mutase from
Corynebacterium glutamicum with bound transition state analog
Authors : Burschowsky, D.; Heim, J.B.; Thorbjørnsrud, H.V.; Kregel, U.
Deposited on : 2016-01-27
Resolution : 2.15 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

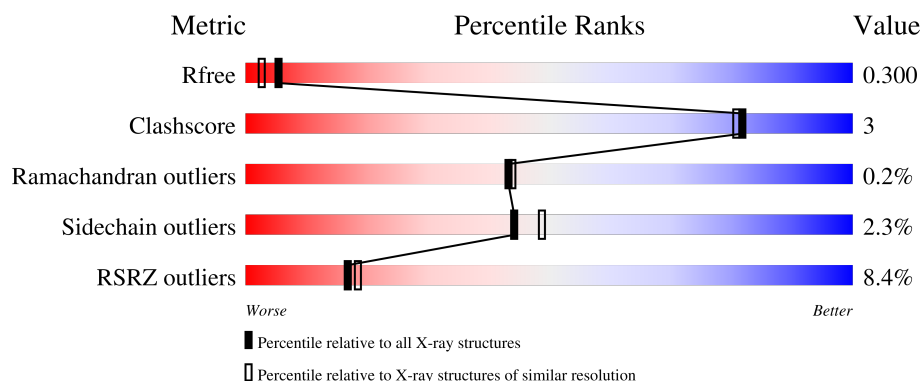
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

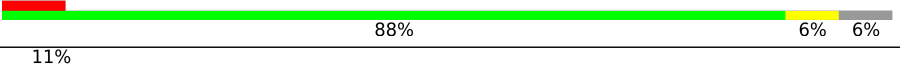
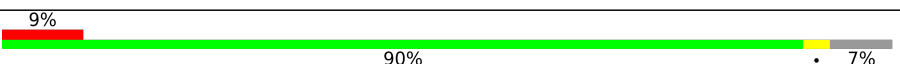
The reported resolution of this entry is 2.15 Å.

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	2057 (2.16-2.16)
Clashscore	190562	2159 (2.16-2.16)
Ramachandran outliers	187476	2134 (2.16-2.16)
Sidechain outliers	187428	2133 (2.16-2.16)
RSRZ outliers	180081	2059 (2.16-2.16)

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	472	 6% 89% 5% 6%
1	B	472	 7% 88% 6% 6%
1	C	472	 11% 88% 9% .
1	D	472	 8% 87% 9% .
2	E	90	 9% 90% . 7%

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Mol	Chain	Length	Quality of chain
2	F	90	12% 87% 6% 8%
2	G	90	11% 89% • 8%
2	H	90	4% 83% 9% 8%

2 Entry composition

There are 11 unique types of molecules in this entry. The entry contains 18279 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 3-Deoxy-D-arabino-heptulosonate 7-phosphate (DAHP) synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	444	Total	C	N	O	S	0	1	0
			3445	2141	615	672	17			
1	B	444	Total	C	N	O	S	0	1	0
			3445	2141	615	672	17			
1	C	456	Total	C	N	O	S	0	1	0
			3537	2199	630	690	18			
1	D	457	Total	C	N	O	S	0	2	0
			3553	2209	634	692	18			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	initiating methionine	UNP Q8NNL5
A	2	HIS	-	expression tag	UNP Q8NNL5
A	3	HIS	-	expression tag	UNP Q8NNL5
A	4	HIS	-	expression tag	UNP Q8NNL5
A	5	HIS	-	expression tag	UNP Q8NNL5
A	6	HIS	-	expression tag	UNP Q8NNL5
A	7	HIS	-	expression tag	UNP Q8NNL5
A	8	SER	-	expression tag	UNP Q8NNL5
A	9	SER	-	expression tag	UNP Q8NNL5
A	10	GLY	-	expression tag	UNP Q8NNL5
B	1	MET	-	initiating methionine	UNP Q8NNL5
B	2	HIS	-	expression tag	UNP Q8NNL5
B	3	HIS	-	expression tag	UNP Q8NNL5
B	4	HIS	-	expression tag	UNP Q8NNL5
B	5	HIS	-	expression tag	UNP Q8NNL5
B	6	HIS	-	expression tag	UNP Q8NNL5
B	7	HIS	-	expression tag	UNP Q8NNL5
B	8	SER	-	expression tag	UNP Q8NNL5
B	9	SER	-	expression tag	UNP Q8NNL5
B	10	GLY	-	expression tag	UNP Q8NNL5

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Chain	Residue	Modelled	Actual	Comment	Reference
C	1	MET	-	initiating methionine	UNP Q8NNL5
C	2	HIS	-	expression tag	UNP Q8NNL5
C	3	HIS	-	expression tag	UNP Q8NNL5
C	4	HIS	-	expression tag	UNP Q8NNL5
C	5	HIS	-	expression tag	UNP Q8NNL5
C	6	HIS	-	expression tag	UNP Q8NNL5
C	7	HIS	-	expression tag	UNP Q8NNL5
C	8	SER	-	expression tag	UNP Q8NNL5
C	9	SER	-	expression tag	UNP Q8NNL5
C	10	GLY	-	expression tag	UNP Q8NNL5
D	1	MET	-	initiating methionine	UNP Q8NNL5
D	2	HIS	-	expression tag	UNP Q8NNL5
D	3	HIS	-	expression tag	UNP Q8NNL5
D	4	HIS	-	expression tag	UNP Q8NNL5
D	5	HIS	-	expression tag	UNP Q8NNL5
D	6	HIS	-	expression tag	UNP Q8NNL5
D	7	HIS	-	expression tag	UNP Q8NNL5
D	8	SER	-	expression tag	UNP Q8NNL5
D	9	SER	-	expression tag	UNP Q8NNL5
D	10	GLY	-	expression tag	UNP Q8NNL5

- Molecule 2 is a protein called Chorismate mutase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	84	Total	C	N	O	S	0	1	0
			672	411	130	129	2			
2	F	83	Total	C	N	O	S	0	1	0
			666	408	132	124	2			
2	G	83	Total	C	N	O	S	0	0	0
			655	402	128	123	2			
2	H	83	Total	C	N	O	S	0	0	0
			655	402	128	123	2			

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



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

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

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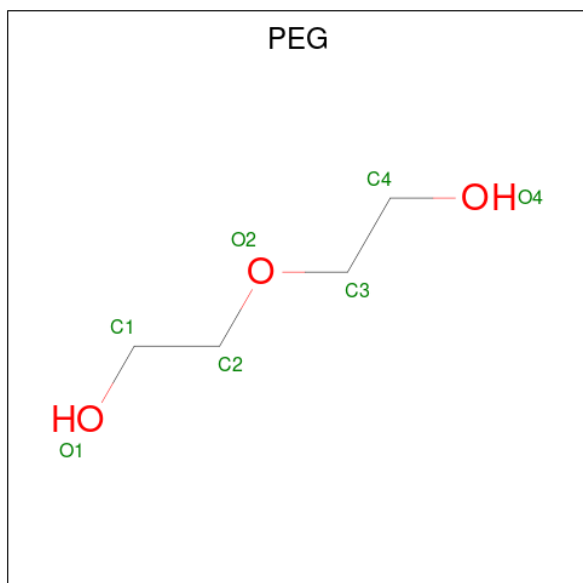
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	D	1	Total	C	O	0	0
			6	3	3		
3	D	1	Total	C	O	0	0
			6	3	3		
3	D	1	Total	C	O	0	0
			6	3	3		
3	D	1	Total	C	O	0	0
			6	3	3		
3	D	1	Total	C	O	0	0
			6	3	3		
3	D	1	Total	C	O	0	0
			6	3	3		
3	D	1	Total	C	O	0	0
			6	3	3		
3	D	1	Total	C	O	0	0
			6	3	3		
3	E	1	Total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

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	G	1	Total	C	O	0	0
			6	3	3		
3	G	1	Total	C	O	0	0
			6	3	3		
3	G	1	Total	C	O	0	0
			6	3	3		
3	G	1	Total	C	O	0	0
			6	3	3		
3	G	1	Total	C	O	0	0
			6	3	3		
3	H	1	Total	C	O	0	0
			6	3	3		
3	H	1	Total	C	O	0	0
			6	3	3		
3	H	1	Total	C	O	0	0
			6	3	3		
3	H	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



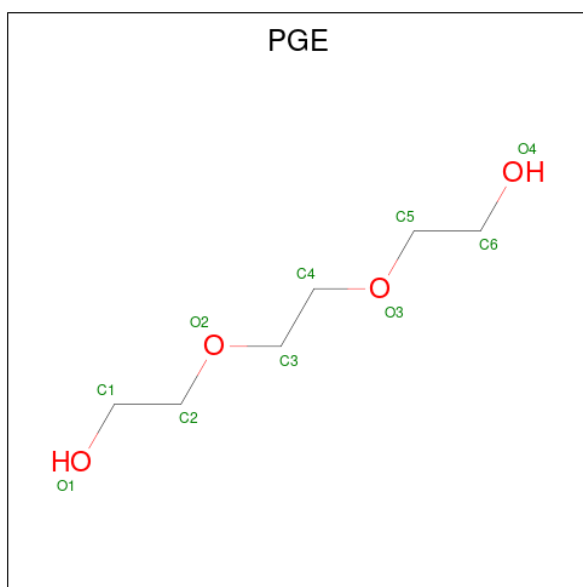
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			7	4	3		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			7	4	3		
4	B	1	Total	C	O	0	0
			7	4	3		
4	C	1	Total	C	O	0	0
			7	4	3		
4	C	1	Total	C	O	0	0
			7	4	3		
4	C	1	Total	C	O	0	0
			7	4	3		
4	D	1	Total	C	O	0	0
			7	4	3		
4	D	1	Total	C	O	0	0
			7	4	3		
4	D	1	Total	C	O	0	0
			7	4	3		
4	D	1	Total	C	O	0	0
			7	4	3		
4	D	1	Total	C	O	0	0
			7	4	3		
4	F	1	Total	C	O	0	0
			7	4	3		
4	G	1	Total	C	O	0	0
			7	4	3		
4	G	1	Total	C	O	0	0
			7	4	3		
4	G	1	Total	C	O	0	0
			7	4	3		

- Molecule 5 is TRIETHYLENE GLYCOL (CCD ID: PGE) (formula: C₆H₁₄O₄).

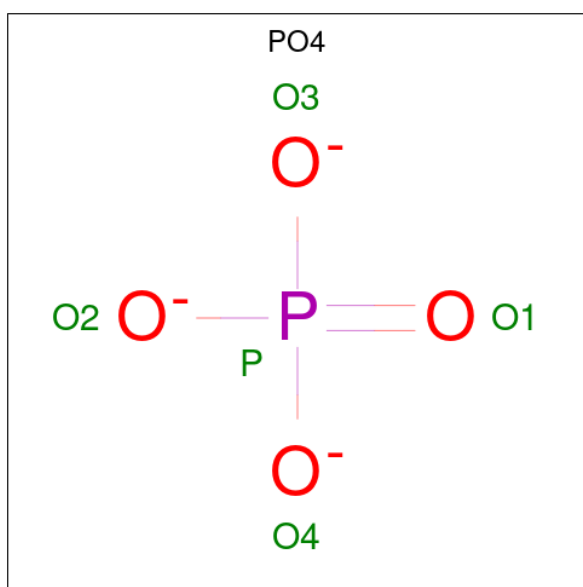


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			10	6	4		
5	C	1	Total	C	O	0	0
			10	6	4		
5	C	1	Total	C	O	0	0
			10	6	4		
5	C	1	Total	C	O	0	0
			10	6	4		
5	C	1	Total	C	O	0	0
			10	6	4		
5	C	1	Total	C	O	0	0
			10	6	4		
5	D	1	Total	C	O	0	0
			10	6	4		
5	D	1	Total	C	O	0	0
			10	6	4		
5	E	1	Total	C	O	0	0
			10	6	4		
5	E	1	Total	C	O	0	0
			10	6	4		
5	E	1	Total	C	O	0	0
			10	6	4		
5	F	1	Total	C	O	0	0
			10	6	4		
5	F	1	Total	C	O	0	0
			10	6	4		
5	G	1	Total	C	O	0	0
			10	6	4		

- Molecule 6 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

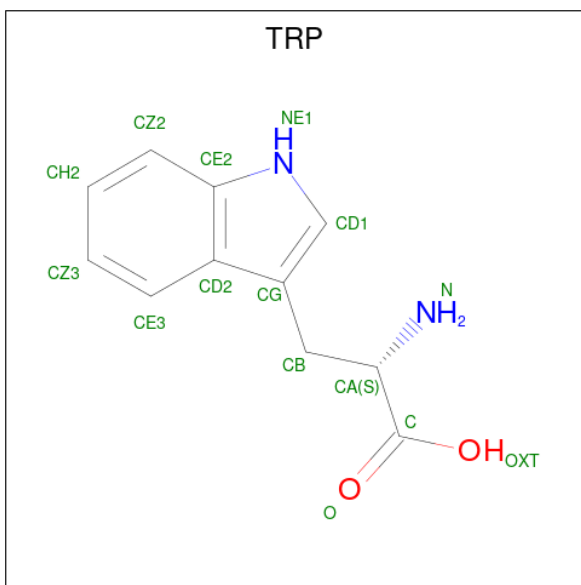
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	1	Total	Mn	0	0
			1	1		
6	B	1	Total	Mn	0	0
			1	1		
6	C	1	Total	Mn	0	0
			1	1		
6	D	1	Total	Mn	0	0
			1	1		

- Molecule 7 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



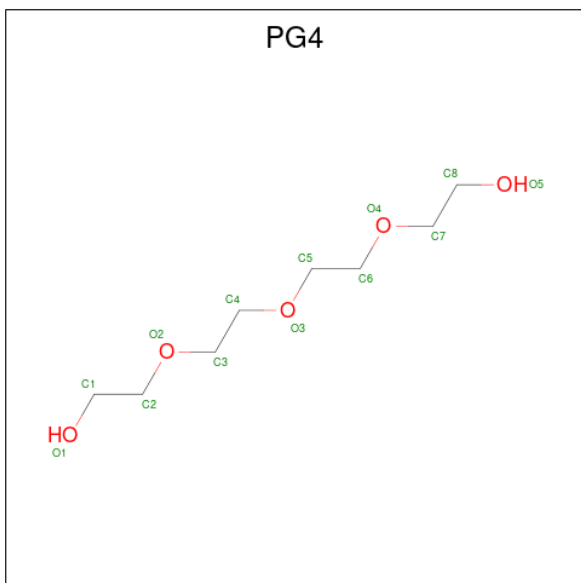
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	O	P	0	0
			5	4	1		
7	A	1	Total	O	P	0	0
			5	4	1		
7	B	1	Total	O	P	0	0
			5	4	1		
7	C	1	Total	O	P	0	0
			5	4	1		
7	D	1	Total	O	P	0	0
			5	4	1		
7	D	1	Total	O	P	0	0
			5	4	1		

- Molecule 8 is TRYPTOPHAN (CCD ID: TRP) (formula: C₁₁H₁₂N₂O₂).



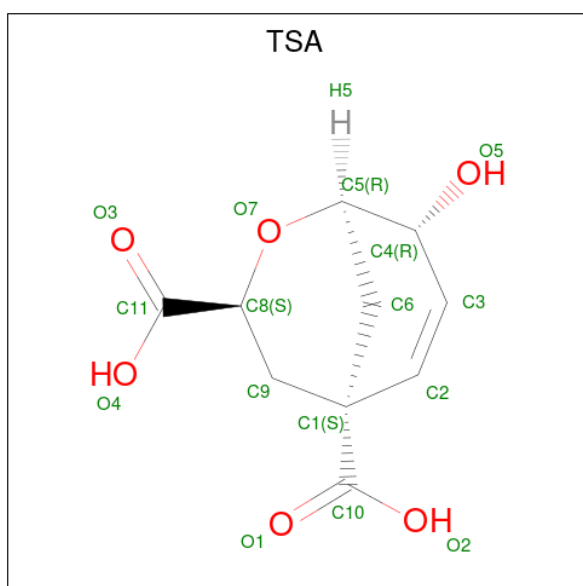
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
8	A	1	Total	C	N	O	0	0
			15	11	2	2		
8	B	1	Total	C	N	O	0	0
			15	11	2	2		
8	C	1	Total	C	N	O	0	0
			15	11	2	2		
8	D	1	Total	C	N	O	0	0
			15	11	2	2		

- Molecule 9 is TETRAETHYLENE GLYCOL (CCD ID: PG4) (formula: $C_8H_{18}O_5$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	B	1	Total	C	O	0	0
			13	8	5		
9	C	1	Total	C	O	0	0
			13	8	5		
9	D	1	Total	C	O	0	0
			13	8	5		
9	D	1	Total	C	O	0	0
			13	8	5		
9	D	1	Total	C	O	0	0
			13	8	5		

- Molecule 10 is 8-HYDROXY-2-OXA-BICYCLO[3.3.1]NON-6-ENE-3,5-DICARBOXYLIC ACID (CCD ID: TSA) (formula: C₁₀H₁₂O₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
10	E	1	Total	C	O	0	0
			16	10	6		
10	F	1	Total	C	O	0	0
			16	10	6		
10	G	1	Total	C	O	0	0
			16	10	6		
10	H	1	Total	C	O	0	0
			16	10	6		

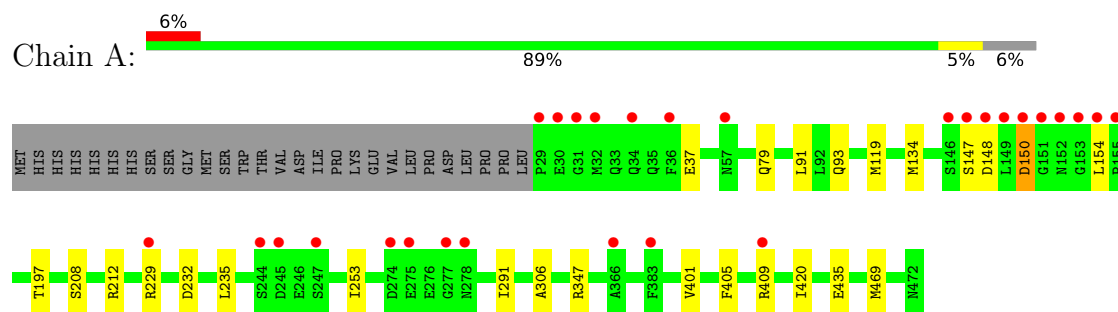
- Molecule 11 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	A	120	Total 120	O 120	0	0
11	B	117	Total 118	O 118	0	1
11	C	183	Total 183	O 183	0	0
11	D	181	Total 181	O 181	0	1
11	E	36	Total 36	O 36	0	0
11	F	49	Total 49	O 49	0	0
11	G	42	Total 42	O 42	0	0
11	H	51	Total 51	O 51	0	0

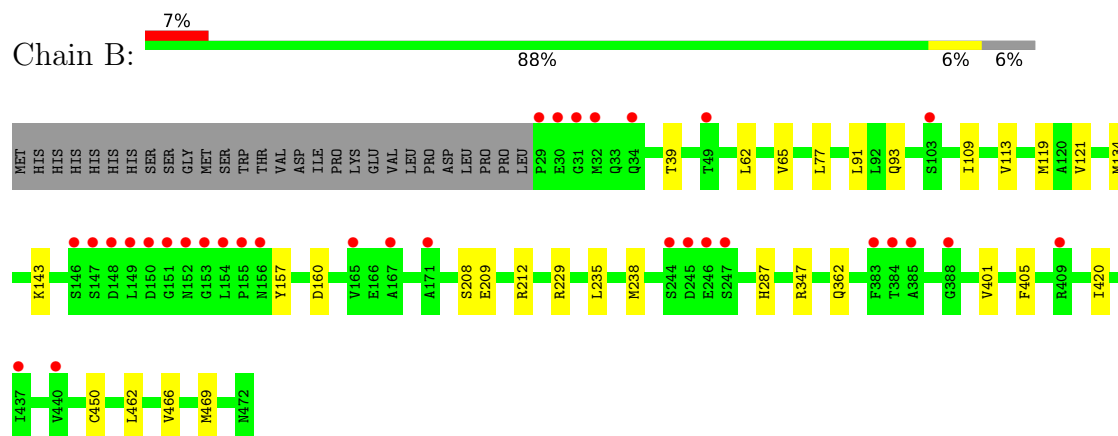
3 Residue-property plots [i](#)

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.

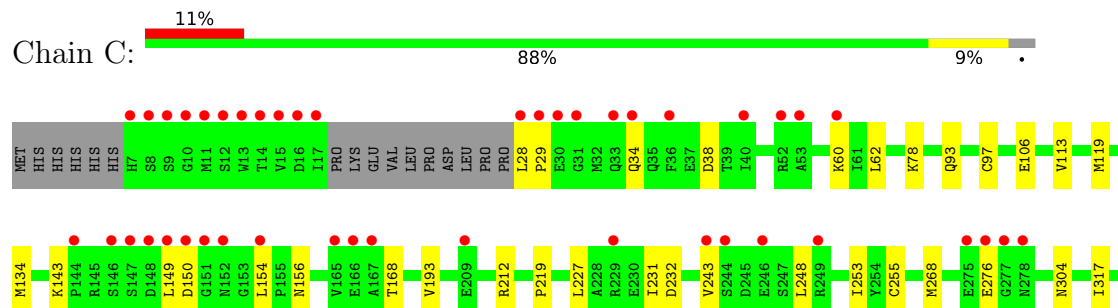
- Molecule 1: 3-Deoxy-D-arabino-heptulosonate 7-phosphate (DAHP) synthase



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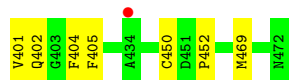
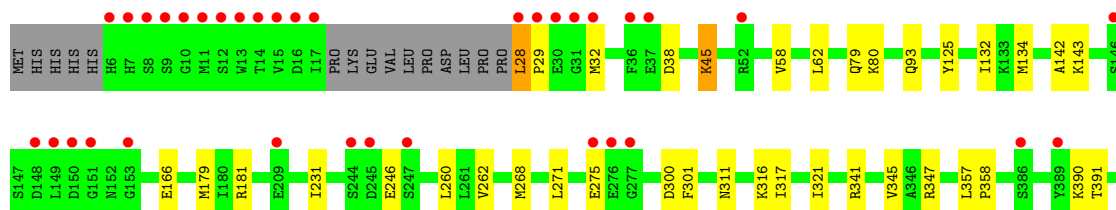
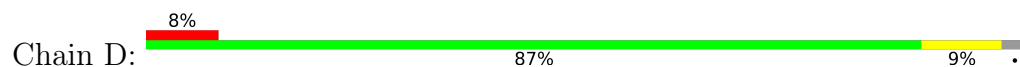


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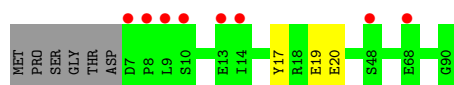
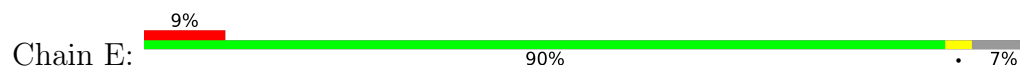




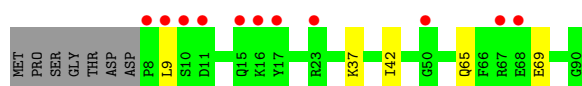
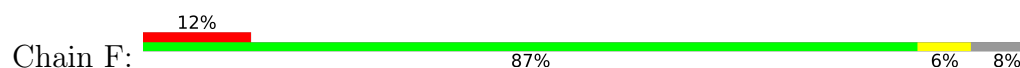
- Molecule 1: 3-Deoxy-D-arabino-heptulosonate 7-phosphate (DAHP) synthase



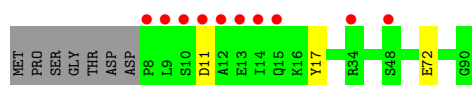
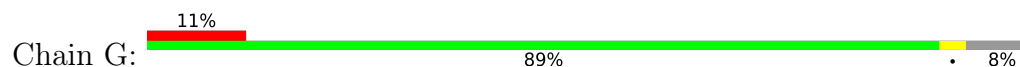
- Molecule 2: Chorismate mutase



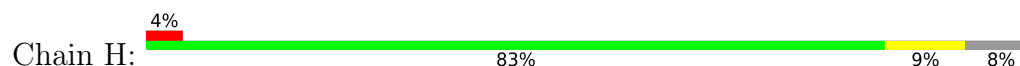
- Molecule 2: Chorismate mutase



- Molecule 2: Chorismate mutase



- Molecule 2: Chorismate mutase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	117.62Å 110.48Å 134.65Å 90.00° 101.41° 90.00°	Depositor
Resolution (Å)	40.00 – 2.15 40.00 – 2.15	Depositor EDS
% Data completeness (in resolution range)	98.0 (40.00-2.15) 98.0 (40.00-2.15)	Depositor EDS
R_{merge}	0.20	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.39 (at 2.16Å)	Xtrriage
Refinement program	REFMAC 5.8.0135	Depositor
R, R_{free}	0.250 , 0.298 0.253 , 0.300	Depositor DCC
R_{free} test set	8850 reflections (4.83%)	wwPDB-VP
Wilson B-factor (Å ²)	24.1	Xtrriage
Anisotropy	0.202	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 42.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	18279	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 22.63 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 5.4860e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: PGE, TSA, PG4, PO4, GOL, PEG, MN

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	0.71	0/3512	0.92	0/4759
1	B	0.69	0/3512	0.90	0/4759
1	C	0.87	1/3606 (0.0%)	1.02	1/4887 (0.0%)
1	D	0.87	0/3626	1.01	0/4914
2	E	0.80	0/676	0.98	0/900
2	F	0.87	0/670	1.12	1/890 (0.1%)
2	G	0.81	0/659	1.10	1/876 (0.1%)
2	H	0.91	0/659	1.14	3/876 (0.3%)
All	All	0.80	1/16920 (0.0%)	0.98	6/22861 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	193	VAL	CA-C	5.06	1.58	1.52

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	72	GLU	N-CA-C	5.47	117.68	111.11
1	C	255	CYS	N-CA-C	5.32	118.78	110.32
2	H	87	GLY	N-CA-C	-5.25	106.41	112.29
2	F	69	GLU	N-CA-C	5.20	117.74	111.40
2	H	85	GLY	N-CA-C	5.16	118.73	112.49
2	H	69	GLU	N-CA-C	5.08	116.63	111.14

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	3445	0	3348	16	0
1	B	3445	0	3348	19	0
1	C	3537	0	3432	19	0
1	D	3553	0	3447	27	0
2	E	672	0	696	3	0
2	F	666	0	700	1	0
2	G	655	0	688	1	0
2	H	655	0	688	3	0
3	A	12	0	16	0	0
3	B	60	0	80	0	0
3	C	120	0	160	5	0
3	D	78	0	104	2	0
3	E	24	0	32	0	0
3	F	36	0	48	0	0
3	G	36	0	48	0	0
3	H	30	0	40	0	0
4	A	14	0	20	1	0
4	B	7	0	10	0	0
4	C	21	0	30	0	0
4	D	42	0	60	0	0
4	F	7	0	10	0	0
4	G	21	0	30	0	0
5	A	10	0	14	0	0
5	C	50	0	70	0	0
5	D	20	0	28	0	0
5	E	30	0	42	0	0
5	F	20	0	28	0	0
5	G	10	0	14	0	0
6	A	1	0	0	0	0
6	B	1	0	0	0	0
6	C	1	0	0	0	0
6	D	1	0	0	0	0
7	A	10	0	0	0	0
7	B	5	0	0	0	0
7	C	5	0	0	0	0
7	D	10	0	0	1	0
8	A	15	0	9	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
8	B	15	0	9	0	0
8	C	15	0	9	0	0
8	D	15	0	9	0	0
9	B	13	0	18	0	0
9	C	13	0	18	0	0
9	D	39	0	54	0	0
10	E	16	0	10	0	0
10	F	16	0	10	0	0
10	G	16	0	10	0	0
10	H	16	0	10	0	0
11	A	120	0	0	1	0
11	B	118	0	0	0	0
11	C	183	0	0	0	0
11	D	181	0	0	1	0
11	E	36	0	0	2	0
11	F	49	0	0	0	0
11	G	42	0	0	0	0
11	H	51	0	0	1	0
All	All	18279	0	17397	88	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 3.

All (88) 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:401:VAL:HG12	1:D:469:MET:HE1	1.39	1.01
1:D:32:MET:HE1	1:D:271:LEU:HD21	1.48	0.95
1:A:401:VAL:HG12	1:A:469:MET:HE1	1.57	0.85
1:B:119:MET:HE1	1:B:420:ILE:HD11	1.59	0.84
1:C:119:MET:HE1	1:C:420:ILE:HD11	1.60	0.83
2:E:20:GLU:CD	11:E:201:HOH:O	2.20	0.82
1:B:401:VAL:HG12	1:B:469:MET:HE1	1.64	0.79
1:A:119:MET:HE1	1:A:420:ILE:HD11	1.65	0.78
1:A:93:GLN:CD	1:A:134:MET:HE3	2.13	0.74
1:D:401:VAL:HG12	1:D:469:MET:CE	2.20	0.70
1:D:405:PHE:CD1	1:D:469:MET:HE2	2.26	0.69
1:A:91:LEU:HD11	1:A:134:MET:HE2	1.75	0.69
1:B:93:GLN:CD	1:B:134:MET:HE3	2.23	0.64
1:C:119:MET:HE1	1:C:420:ILE:CD1	2.28	0.63
2:E:20:GLU:OE2	11:E:201:HOH:O	2.16	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:143:LYS:NZ	1:D:450:CYS:SG	2.69	0.63
1:D:405:PHE:CD1	1:D:469:MET:CE	2.82	0.62
1:B:119:MET:HE1	1:B:420:ILE:CD1	2.29	0.61
1:B:143:LYS:NZ	1:B:450:CYS:SG	2.67	0.61
1:C:304:ASN:OD1	1:C:332:LYS:NZ	2.35	0.60
1:D:181:ARG:NH1	7:D:527:PO4:O1	2.35	0.59
1:B:238:MET:HE3	1:C:243:VAL:HG21	1.85	0.57
2:E:17:TYR:HB3	2:F:42:ILE:HD11	1.88	0.55
1:D:317:ILE:HD12	1:D:321:ILE:HG21	1.89	0.55
1:B:401:VAL:CG1	1:B:469:MET:HE1	2.36	0.54
1:A:119:MET:HE1	1:A:420:ILE:CD1	2.34	0.54
1:A:405:PHE:CD1	1:A:469:MET:CE	2.90	0.54
1:A:93:GLN:HA	1:A:134:MET:O	2.07	0.54
1:D:80:LYS:HE3	3:D:511:GOL:H31	1.90	0.53
2:G:17:TYR:HB3	2:H:42:ILE:HD11	1.91	0.53
1:C:357:LEU:O	1:C:358:PRO:C	2.52	0.53
1:D:132:ILE:HD12	1:D:132:ILE:N	2.23	0.53
1:D:28:LEU:N	1:D:29:PRO:HD2	2.24	0.52
1:C:143:LYS:NZ	1:C:450:CYS:SG	2.73	0.52
1:B:91:LEU:HD11	1:B:134:MET:HE2	1.91	0.51
1:B:93:GLN:HA	1:B:134:MET:O	2.11	0.51
1:C:328:ALA:HB3	3:C:502:GOL:C3	2.43	0.49
1:B:401:VAL:HG12	1:B:469:MET:CE	2.37	0.49
1:A:291:ILE:HD11	1:A:306:ALA:CB	2.43	0.48
1:D:62:LEU:HB3	1:D:268:MET:CE	2.44	0.48
1:A:291:ILE:HD11	1:A:306:ALA:HB2	1.96	0.48
1:C:409[B]:ARG:HH11	1:C:409[B]:ARG:HG3	1.79	0.48
1:C:227:LEU:CD1	1:C:231:ILE:HD11	2.45	0.47
2:H:65:GLN:NE2	11:H:204:HOH:O	2.48	0.47
1:D:93:GLN:HA	1:D:134:MET:O	2.15	0.47
2:H:47:MET:HE1	2:H:53:ARG:NH2	2.29	0.47
1:C:28:LEU:N	1:C:29:PRO:HD3	2.30	0.46
1:C:62:LEU:HB3	1:C:268:MET:HE2	1.97	0.46
3:D:513:GOL:H32	11:D:606:HOH:O	2.15	0.46
1:D:402:GLN:HA	1:D:469:MET:HE3	1.97	0.46
1:D:311:ASN:O	1:D:341:ARG:NH2	2.49	0.45
1:D:316:LYS:HD3	1:D:345:VAL:HB	1.97	0.45
1:D:357:LEU:O	1:D:358:PRO:C	2.59	0.45
1:C:328:ALA:HB1	3:C:502:GOL:H11	1.97	0.44
1:C:113:VAL:HG13	1:C:253:ILE:CD1	2.47	0.44
1:C:212:ARG:O	3:C:508:GOL:O2	2.36	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:142:ALA:HB1	1:D:260:LEU:HD23	2.00	0.44
1:A:401:VAL:CG1	1:A:469:MET:HE1	2.38	0.44
1:B:405:PHE:CD1	1:B:469:MET:CE	3.01	0.44
1:C:93:GLN:HA	1:C:134:MET:O	2.17	0.43
1:A:150:ASP:O	1:A:154:LEU:HD13	2.18	0.43
1:B:62:LEU:O	1:B:65:VAL:HG22	2.19	0.43
1:D:401:VAL:O	1:D:404:PHE:HB3	2.18	0.43
1:A:405:PHE:CD1	1:A:469:MET:HE2	2.54	0.42
1:B:208:SER:OG	1:B:235:LEU:HD11	2.20	0.42
1:C:219:PRO:HG2	1:C:434:ALA:HB1	2.01	0.42
1:C:328:ALA:HB3	3:C:502:GOL:H32	2.01	0.42
1:B:157:TYR:OH	1:B:160:ASP:OD1	2.35	0.42
4:A:503:PEG:H12	11:A:682:HOH:O	2.19	0.42
1:A:208:SER:OG	1:A:235:LEU:HD11	2.21	0.41
1:A:401:VAL:HG12	1:A:469:MET:CE	2.39	0.41
1:D:262:VAL:HG12	1:D:301:PHE:CE2	2.56	0.41
1:B:209:GLU:OE2	1:B:212:ARG:NH1	2.46	0.41
1:D:143:LYS:CE	1:D:450:CYS:SG	3.09	0.41
1:A:212:ARG:NH1	1:A:232:ASP:OD1	2.51	0.41
1:B:462:LEU:O	1:B:466:VAL:HG23	2.21	0.41
1:D:58:VAL:HG11	1:D:179:MET:HE3	2.01	0.41
1:C:97:CYS:SG	1:C:450:CYS:HB3	2.60	0.41
1:D:45:LYS:NZ	1:D:300:ASP:OD2	2.48	0.41
1:D:391:THR:HA	1:D:452:PRO:HG3	2.02	0.41
1:B:77:LEU:HD22	1:B:287:HIS:HB3	2.03	0.41
1:B:109:ILE:O	1:B:113:VAL:HG23	2.21	0.40
1:D:80:LYS:HB3	1:D:132:ILE:HG12	2.03	0.40
1:A:197:THR:CG2	1:A:253:ILE:HD12	2.51	0.40
1:B:93:GLN:NE2	1:B:134:MET:HE3	2.36	0.40
1:D:125:TYR:CG	1:D:231:ILE:HG12	2.57	0.40
1:D:405:PHE:CE1	1:D:469:MET:HE2	2.55	0.40
1:C:231:ILE:HG21	3:C:512:GOL:C3	2.51	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	443/472 (94%)	429 (97%)	12 (3%)	2 (0%)	24	20
1	B	443/472 (94%)	431 (97%)	12 (3%)	0	100	100
1	C	453/472 (96%)	435 (96%)	17 (4%)	1 (0%)	43	44
1	D	455/472 (96%)	439 (96%)	15 (3%)	1 (0%)	43	44
2	E	83/90 (92%)	81 (98%)	2 (2%)	0	100	100
2	F	82/90 (91%)	81 (99%)	1 (1%)	0	100	100
2	G	81/90 (90%)	80 (99%)	1 (1%)	0	100	100
2	H	81/90 (90%)	80 (99%)	1 (1%)	0	100	100
All	All	2121/2248 (94%)	2056 (97%)	61 (3%)	4 (0%)	43	44

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	148	ASP
1	C	150	ASP
1	A	147	SER
1	D	246	GLU

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	363/389 (93%)	355 (98%)	8 (2%)	45	51
1	B	363/389 (93%)	358 (99%)	5 (1%)	59	66
1	C	374/389 (96%)	360 (96%)	14 (4%)	30	30
1	D	376/389 (97%)	368 (98%)	8 (2%)	47	52
2	E	71/75 (95%)	69 (97%)	2 (3%)	38	40
2	F	70/75 (93%)	67 (96%)	3 (4%)	26	24

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	G	69/75 (92%)	68 (99%)	1 (1%)	59	66
2	H	69/75 (92%)	68 (99%)	1 (1%)	59	66
All	All	1755/1856 (95%)	1713 (98%)	42 (2%)	44	47

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	37	GLU
1	A	79	GLN
1	A	150	ASP
1	A	229	ARG
1	A	347	ARG
1	A	409[A]	ARG
1	A	409[B]	ARG
1	A	435	GLU
1	B	39	THR
1	B	121	VAL
1	B	229	ARG
1	B	347	ARG
1	B	362	GLN
1	C	34	GLN
1	C	38	ASP
1	C	60	LYS
1	C	78	LYS
1	C	106	GLU
1	C	149	LEU
1	C	154	LEU
1	C	156	ASN
1	C	168	THR
1	C	232	ASP
1	C	248	LEU
1	C	276	GLU
1	C	317	ILE
1	C	347	ARG
1	D	28	LEU
1	D	38	ASP
1	D	45	LYS
1	D	79	GLN
1	D	166	GLU
1	D	275	GLU
1	D	347	ARG

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Mol	Chain	Res	Type
1	D	390	LYS
2	E	19[A]	GLU
2	E	19[B]	GLU
2	F	9	LEU
2	F	37	LYS
2	F	65	GLN
2	G	11	ASP
2	H	37	LYS

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

Mol	Chain	Res	Type
1	A	118	GLN
1	A	156	ASN
1	A	387	ASN
1	B	118	GLN
1	B	362	GLN
1	B	472	ASN
1	C	34	GLN
1	C	156	ASN
1	C	457	GLN
1	C	472	ASN
1	D	79	GLN
1	D	118	GLN
1	D	336	ASN
2	F	15	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 119 ligands modelled in this entry, 4 are monoatomic - leaving 115 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	$\# Z > 2$
3	GOL	C	508	-	5,5,5	0.63	0	5,5,5	0.38	0
7	PO4	A	507	-	4,4,4	1.12	0	6,6,6	0.50	0
3	GOL	C	516	-	5,5,5	0.41	0	5,5,5	0.18	0
3	GOL	H	101	-	5,5,5	0.48	0	5,5,5	0.49	0
3	GOL	C	509	-	5,5,5	0.27	0	5,5,5	0.36	0
4	PEG	C	521	-	6,6,6	0.79	0	5,5,5	0.88	0
5	PGE	A	505	-	9,9,9	0.64	0	8,8,8	0.47	0
3	GOL	D	507	-	5,5,5	0.69	0	5,5,5	0.64	0
3	GOL	B	502	-	5,5,5	0.42	0	5,5,5	0.30	0
3	GOL	E	103	-	5,5,5	0.41	0	5,5,5	0.39	0
5	PGE	E	105	-	9,9,9	0.54	0	8,8,8	0.35	0
5	PGE	F	108	-	9,9,9	0.65	0	8,8,8	0.42	0
5	PGE	C	528	-	9,9,9	0.64	0	8,8,8	0.80	0
3	GOL	F	103	-	5,5,5	0.44	0	5,5,5	1.24	0
3	GOL	B	508	-	5,5,5	0.41	0	5,5,5	0.15	0
3	GOL	B	506	-	5,5,5	0.42	0	5,5,5	0.57	0
8	TRP	B	515	-	15,16,16	0.76	0	18,22,22	0.71	1 (5%)
3	GOL	E	104	-	5,5,5	0.44	0	5,5,5	0.19	0
4	PEG	A	503	-	6,6,6	0.43	0	5,5,5	0.30	0
7	PO4	B	514	-	4,4,4	0.94	0	6,6,6	0.54	0
4	PEG	D	518	-	6,6,6	0.49	0	5,5,5	0.37	0
3	GOL	D	503	-	5,5,5	0.32	0	5,5,5	0.27	0
3	GOL	C	511	-	5,5,5	0.62	0	5,5,5	0.58	0
9	PG4	C	529	-	12,12,12	0.65	0	11,11,11	0.54	0
4	PEG	G	109	-	6,6,6	0.50	0	5,5,5	0.43	0
3	GOL	B	507	-	5,5,5	0.29	0	5,5,5	0.49	0
3	GOL	F	101	-	5,5,5	0.50	0	5,5,5	0.31	0
10	TSA	H	106	-	16,17,17	1.49	4 (25%)	16,26,26	1.76	4 (25%)
5	PGE	C	527	-	9,9,9	0.64	0	8,8,8	0.68	0
3	GOL	D	505	-	5,5,5	0.65	0	5,5,5	0.72	0
3	GOL	C	504	-	5,5,5	0.44	0	5,5,5	0.51	0
9	PG4	D	524	-	12,12,12	0.62	0	11,11,11	0.92	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	GOL	C	514	-	5,5,5	0.46	0	5,5,5	0.62	0
3	GOL	C	515	-	5,5,5	0.56	0	5,5,5	0.54	0
4	PEG	B	511	-	6,6,6	0.45	0	5,5,5	0.33	0
3	GOL	A	502	-	5,5,5	0.46	0	5,5,5	0.26	0
8	TRP	A	509	-	15,16,16	0.83	1 (6%)	18,22,22	0.64	1 (5%)
3	GOL	C	513	-	5,5,5	0.60	0	5,5,5	0.36	0
3	GOL	G	105	-	5,5,5	0.46	0	5,5,5	0.52	0
4	PEG	A	504	-	6,6,6	0.50	0	5,5,5	0.43	0
9	PG4	B	512	-	12,12,12	0.66	0	11,11,11	0.37	0
9	PG4	D	523	-	12,12,12	0.57	0	11,11,11	0.30	0
4	PEG	D	519	-	6,6,6	0.61	0	5,5,5	0.25	0
7	PO4	C	531	-	4,4,4	1.08	0	6,6,6	0.73	0
3	GOL	B	509	-	5,5,5	0.34	0	5,5,5	0.17	0
7	PO4	D	527	-	4,4,4	0.82	0	6,6,6	0.82	0
3	GOL	H	105	-	5,5,5	0.53	0	5,5,5	0.35	0
3	GOL	C	506	-	5,5,5	0.31	0	5,5,5	0.30	0
3	GOL	C	510	-	5,5,5	0.33	0	5,5,5	0.26	0
3	GOL	C	501	-	5,5,5	0.45	0	5,5,5	0.39	0
3	GOL	F	106	-	5,5,5	0.39	0	5,5,5	0.35	0
3	GOL	C	517	-	5,5,5	0.43	0	5,5,5	0.21	0
3	GOL	F	105	-	5,5,5	0.36	0	5,5,5	0.29	0
3	GOL	C	520	-	5,5,5	0.42	0	5,5,5	0.41	0
3	GOL	C	503	-	5,5,5	0.26	0	5,5,5	0.37	0
4	PEG	D	514	-	6,6,6	0.60	0	5,5,5	1.10	0
4	PEG	F	107	-	6,6,6	0.56	0	5,5,5	0.52	0
7	PO4	D	526	-	4,4,4	0.82	0	6,6,6	0.66	0
4	PEG	C	522	-	6,6,6	0.57	0	5,5,5	0.58	0
3	GOL	D	511	-	5,5,5	0.71	0	5,5,5	0.61	0
3	GOL	D	504	-	5,5,5	0.50	0	5,5,5	0.54	0
9	PG4	D	522	-	12,12,12	0.69	0	11,11,11	1.02	1 (9%)
5	PGE	E	107	-	9,9,9	0.63	0	8,8,8	0.42	0
5	PGE	D	521	-	9,9,9	0.64	0	8,8,8	0.38	0
5	PGE	C	526	-	9,9,9	0.65	0	8,8,8	0.52	0
4	PEG	D	516	-	6,6,6	0.55	0	5,5,5	0.27	0
4	PEG	G	108	-	6,6,6	0.62	0	5,5,5	0.30	0
8	TRP	C	532	-	15,16,16	0.82	1 (6%)	18,22,22	0.73	1 (5%)
3	GOL	G	106	-	5,5,5	0.35	0	5,5,5	0.22	0
3	GOL	H	102	-	5,5,5	0.34	0	5,5,5	0.59	0
3	GOL	G	102	-	5,5,5	0.73	0	5,5,5	0.62	0
3	GOL	C	502	-	5,5,5	0.57	0	5,5,5	0.92	0
4	PEG	D	515	-	6,6,6	0.59	0	5,5,5	0.49	0
5	PGE	G	110	-	9,9,9	0.56	0	8,8,8	0.34	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	GOL	G	104	-	5,5,5	0.42	0	5,5,5	0.58	0
3	GOL	G	103	-	5,5,5	0.37	0	5,5,5	0.25	0
3	GOL	C	505	-	5,5,5	0.58	0	5,5,5	0.48	0
3	GOL	H	104	-	5,5,5	0.70	0	5,5,5	0.64	0
5	PGE	F	109	-	9,9,9	0.62	0	8,8,8	0.39	0
5	PGE	C	525	-	9,9,9	0.53	0	8,8,8	0.51	0
3	GOL	F	102	-	5,5,5	0.46	0	5,5,5	0.30	0
10	TSA	F	110	-	16,17,17	1.53	4 (25%)	16,26,26	1.41	2 (12%)
3	GOL	B	504	-	5,5,5	0.40	0	5,5,5	0.33	0
3	GOL	C	507	-	5,5,5	0.41	0	5,5,5	0.98	0
3	GOL	E	101	-	5,5,5	0.40	0	5,5,5	0.19	0
3	GOL	C	512	-	5,5,5	0.24	0	5,5,5	0.36	0
3	GOL	G	101	-	5,5,5	0.52	0	5,5,5	0.63	0
3	GOL	A	501	-	5,5,5	0.37	0	5,5,5	0.39	0
7	PO4	A	508	-	4,4,4	0.97	0	6,6,6	0.57	0
3	GOL	D	502	-	5,5,5	0.57	0	5,5,5	0.83	0
3	GOL	F	104	-	5,5,5	0.41	0	5,5,5	0.65	0
8	TRP	D	528	-	15,16,16	0.85	0	18,22,22	0.68	0
3	GOL	B	505	-	5,5,5	0.49	0	5,5,5	0.49	0
3	GOL	D	508	-	5,5,5	0.78	0	5,5,5	0.61	0
3	GOL	E	102	-	5,5,5	0.50	0	5,5,5	0.24	0
3	GOL	C	519	-	5,5,5	0.47	0	5,5,5	0.44	0
3	GOL	D	509	-	5,5,5	0.35	0	5,5,5	0.38	0
3	GOL	D	513	-	5,5,5	0.53	0	5,5,5	0.62	0
10	TSA	G	111	-	16,17,17	1.34	1 (6%)	16,26,26	1.39	4 (25%)
5	PGE	E	106	-	9,9,9	0.70	0	8,8,8	0.47	0
3	GOL	B	503	-	5,5,5	0.42	0	5,5,5	0.67	0
3	GOL	B	510	-	5,5,5	0.25	0	5,5,5	0.36	0
3	GOL	B	501	-	5,5,5	0.28	0	5,5,5	0.52	0
4	PEG	G	107	-	6,6,6	0.72	0	5,5,5	0.92	0
3	GOL	C	518	-	5,5,5	0.57	0	5,5,5	0.65	0
3	GOL	D	506	-	5,5,5	0.29	0	5,5,5	0.32	0
3	GOL	D	510	-	5,5,5	0.30	0	5,5,5	0.28	0
5	PGE	D	520	-	9,9,9	0.56	0	8,8,8	0.39	0
5	PGE	C	524	-	9,9,9	0.55	0	8,8,8	0.76	0
3	GOL	D	501	-	5,5,5	0.51	0	5,5,5	0.24	0
3	GOL	D	512	-	5,5,5	0.40	0	5,5,5	0.42	0
3	GOL	H	103	-	5,5,5	0.38	0	5,5,5	0.67	0
4	PEG	D	517	-	6,6,6	0.55	0	5,5,5	0.18	0
10	TSA	E	108	-	16,17,17	1.12	0	16,26,26	1.67	4 (25%)
4	PEG	C	523	-	6,6,6	0.53	0	5,5,5	0.44	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
3	GOL	C	508	-	-	4/4/4/4	-
3	GOL	C	516	-	-	2/4/4/4	-
3	GOL	H	101	-	-	2/4/4/4	-
3	GOL	C	509	-	-	0/4/4/4	-
4	PEG	C	521	-	-	2/4/4/4	-
5	PGE	A	505	-	-	5/7/7/7	-
3	GOL	D	507	-	-	2/4/4/4	-
3	GOL	B	502	-	-	2/4/4/4	-
3	GOL	E	103	-	-	1/4/4/4	-
5	PGE	E	105	-	-	2/7/7/7	-
5	PGE	F	108	-	-	4/7/7/7	-
5	PGE	C	528	-	-	5/7/7/7	-
3	GOL	F	103	-	-	2/4/4/4	-
3	GOL	B	508	-	-	2/4/4/4	-
3	GOL	B	506	-	-	0/4/4/4	-
8	TRP	B	515	-	-	0/8/8/8	0/2/2/2
3	GOL	E	104	-	-	4/4/4/4	-
4	PEG	A	503	-	-	2/4/4/4	-
4	PEG	D	518	-	-	2/4/4/4	-
3	GOL	D	503	-	-	2/4/4/4	-
3	GOL	C	511	-	-	2/4/4/4	-
9	PG4	C	529	-	-	7/10/10/10	-
4	PEG	G	109	-	-	4/4/4/4	-
3	GOL	B	507	-	-	2/4/4/4	-
3	GOL	F	101	-	-	4/4/4/4	-
10	TSA	H	106	-	-	0/10/34/34	1/3/2/2
5	PGE	C	527	-	-	3/7/7/7	-
3	GOL	D	505	-	-	4/4/4/4	-
3	GOL	C	504	-	-	2/4/4/4	-
9	PG4	D	524	-	-	4/10/10/10	-
3	GOL	C	514	-	-	2/4/4/4	-
3	GOL	C	515	-	-	2/4/4/4	-
4	PEG	B	511	-	-	0/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GOL	A	502	-	-	3/4/4/4	-
8	TRP	A	509	-	-	1/8/8/8	0/2/2/2
3	GOL	C	513	-	-	3/4/4/4	-
3	GOL	G	105	-	-	0/4/4/4	-
4	PEG	A	504	-	-	3/4/4/4	-
9	PG4	B	512	-	-	1/10/10/10	-
9	PG4	D	523	-	-	5/10/10/10	-
4	PEG	D	519	-	-	1/4/4/4	-
3	GOL	B	509	-	-	2/4/4/4	-
3	GOL	H	105	-	-	4/4/4/4	-
3	GOL	C	506	-	-	2/4/4/4	-
3	GOL	C	510	-	-	0/4/4/4	-
3	GOL	C	501	-	-	2/4/4/4	-
3	GOL	F	106	-	-	2/4/4/4	-
3	GOL	C	517	-	-	2/4/4/4	-
3	GOL	F	105	-	-	2/4/4/4	-
3	GOL	C	520	-	-	0/4/4/4	-
3	GOL	C	503	-	-	0/4/4/4	-
4	PEG	D	514	-	-	3/4/4/4	-
4	PEG	F	107	-	-	3/4/4/4	-
4	PEG	C	522	-	-	1/4/4/4	-
3	GOL	D	511	-	-	4/4/4/4	-
3	GOL	D	504	-	-	4/4/4/4	-
9	PG4	D	522	-	-	6/10/10/10	-
5	PGE	E	107	-	-	2/7/7/7	-
5	PGE	D	521	-	-	3/7/7/7	-
5	PGE	C	526	-	-	3/7/7/7	-
4	PEG	D	516	-	-	3/4/4/4	-
4	PEG	G	108	-	-	2/4/4/4	-
8	TRP	C	532	-	-	0/8/8/8	0/2/2/2
3	GOL	G	106	-	-	3/4/4/4	-
3	GOL	H	102	-	-	2/4/4/4	-
3	GOL	G	102	-	-	2/4/4/4	-
3	GOL	C	502	-	-	3/4/4/4	-
4	PEG	D	515	-	-	1/4/4/4	-
5	PGE	G	110	-	-	3/7/7/7	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GOL	G	104	-	-	2/4/4/4	-
3	GOL	G	103	-	-	2/4/4/4	-
3	GOL	C	505	-	-	2/4/4/4	-
3	GOL	H	104	-	-	2/4/4/4	-
5	PGE	F	109	-	-	2/7/7/7	-
5	PGE	C	525	-	-	4/7/7/7	-
3	GOL	F	102	-	-	4/4/4/4	-
10	TSA	F	110	-	-	1/10/34/34	1/3/2/2
3	GOL	B	504	-	-	3/4/4/4	-
3	GOL	C	507	-	-	2/4/4/4	-
3	GOL	E	101	-	-	0/4/4/4	-
3	GOL	C	512	-	-	2/4/4/4	-
3	GOL	G	101	-	-	0/4/4/4	-
3	GOL	A	501	-	-	2/4/4/4	-
3	GOL	D	502	-	-	1/4/4/4	-
3	GOL	F	104	-	-	4/4/4/4	-
8	TRP	D	528	-	-	0/8/8/8	0/2/2/2
3	GOL	B	505	-	-	0/4/4/4	-
3	GOL	D	508	-	-	4/4/4/4	-
3	GOL	E	102	-	-	0/4/4/4	-
3	GOL	C	519	-	-	4/4/4/4	-
3	GOL	D	509	-	-	2/4/4/4	-
3	GOL	D	513	-	-	2/4/4/4	-
10	TSA	G	111	-	-	1/10/34/34	1/3/2/2
5	PGE	E	106	-	-	3/7/7/7	-
3	GOL	B	503	-	-	2/4/4/4	-
3	GOL	B	510	-	-	3/4/4/4	-
3	GOL	B	501	-	-	3/4/4/4	-
4	PEG	G	107	-	-	2/4/4/4	-
3	GOL	C	518	-	-	2/4/4/4	-
3	GOL	D	506	-	-	0/4/4/4	-
3	GOL	D	510	-	-	0/4/4/4	-
5	PGE	D	520	-	-	5/7/7/7	-
5	PGE	C	524	-	-	5/7/7/7	-
3	GOL	D	501	-	-	4/4/4/4	-
3	GOL	D	512	-	-	2/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GOL	H	103	-	-	1/4/4/4	-
4	PEG	D	517	-	-	2/4/4/4	-
10	TSA	E	108	-	-	1/10/34/34	1/3/2/2
4	PEG	C	523	-	-	1/4/4/4	-

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
10	H	106	TSA	C3-C2	2.79	1.36	1.32
10	F	110	TSA	C3-C2	2.71	1.36	1.32
10	F	110	TSA	C9-C1	-2.65	1.51	1.55
10	G	111	TSA	C6-C1	-2.57	1.51	1.55
10	H	106	TSA	C6-C1	-2.51	1.51	1.55
10	H	106	TSA	C9-C1	-2.47	1.51	1.55
10	F	110	TSA	C6-C1	-2.31	1.51	1.55
10	F	110	TSA	O4-C11	-2.28	1.23	1.30
8	A	509	TRP	OXT-C	-2.20	1.23	1.30
10	H	106	TSA	C8-C11	-2.16	1.49	1.52
8	C	532	TRP	OXT-C	-2.02	1.24	1.30

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
10	H	106	TSA	O7-C8-C9	3.94	115.54	109.95
10	F	110	TSA	O7-C8-C9	3.50	114.92	109.95
10	E	108	TSA	O7-C8-C9	3.50	114.91	109.95
10	E	108	TSA	O4-C11-C8	3.39	121.52	112.71
10	H	106	TSA	O4-C11-C8	3.10	120.78	112.71
10	G	111	TSA	O7-C8-C11	2.76	112.93	107.72
10	G	111	TSA	O4-C11-O3	-2.70	117.97	124.08
8	B	515	TRP	OXT-C-O	-2.56	118.27	124.08
10	G	111	TSA	O4-C11-C8	2.55	119.34	112.71
10	E	108	TSA	O3-C11-C8	-2.55	117.35	122.85
10	H	106	TSA	O3-C11-C8	-2.46	117.54	122.85
10	H	106	TSA	C9-C1-C10	-2.31	104.40	109.84
10	F	110	TSA	C9-C1-C10	-2.27	104.50	109.84
10	E	108	TSA	C9-C1-C10	-2.25	104.54	109.84
8	A	509	TRP	OXT-C-O	-2.23	119.02	124.08
9	D	522	PG4	C7-O4-C6	2.15	122.67	113.26
8	C	532	TRP	OXT-C-O	-2.08	119.36	124.08
10	G	111	TSA	C9-C1-C10	-2.06	104.98	109.84

There are no chirality outliers.

All (244) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	501	GOL	O1-C1-C2-O2
3	A	501	GOL	O1-C1-C2-C3
3	A	502	GOL	O1-C1-C2-C3
3	B	502	GOL	O1-C1-C2-O2
3	B	503	GOL	O1-C1-C2-C3
3	B	508	GOL	O1-C1-C2-O2
3	B	508	GOL	O1-C1-C2-C3
3	C	502	GOL	O1-C1-C2-C3
3	C	506	GOL	O2-C2-C3-O3
3	C	507	GOL	C1-C2-C3-O3
3	C	508	GOL	O1-C1-C2-C3
3	C	511	GOL	O1-C1-C2-C3
3	C	513	GOL	C1-C2-C3-O3
3	C	513	GOL	O2-C2-C3-O3
3	C	514	GOL	O1-C1-C2-O2
3	C	518	GOL	O1-C1-C2-C3
3	C	519	GOL	O1-C1-C2-O2
3	C	519	GOL	O1-C1-C2-C3
3	D	501	GOL	O1-C1-C2-C3
3	D	501	GOL	C1-C2-C3-O3
3	D	503	GOL	C1-C2-C3-O3
3	D	504	GOL	O1-C1-C2-C3
3	D	505	GOL	O1-C1-C2-O2
3	D	505	GOL	C1-C2-C3-O3
3	D	507	GOL	C1-C2-C3-O3
3	D	508	GOL	O1-C1-C2-C3
3	D	511	GOL	O1-C1-C2-C3
3	D	511	GOL	C1-C2-C3-O3
3	D	512	GOL	C1-C2-C3-O3
3	D	513	GOL	C1-C2-C3-O3
3	E	104	GOL	C1-C2-C3-O3
3	F	101	GOL	O1-C1-C2-C3
3	F	102	GOL	O1-C1-C2-C3
3	F	102	GOL	C1-C2-C3-O3
3	F	103	GOL	O1-C1-C2-C3
3	F	104	GOL	C1-C2-C3-O3
3	F	105	GOL	O1-C1-C2-O2
3	F	105	GOL	O1-C1-C2-C3
3	F	106	GOL	C1-C2-C3-O3
3	G	103	GOL	C1-C2-C3-O3

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Mol	Chain	Res	Type	Atoms
3	G	103	GOL	O2-C2-C3-O3
3	G	104	GOL	O1-C1-C2-O2
3	G	104	GOL	O1-C1-C2-C3
3	G	106	GOL	O1-C1-C2-O2
3	G	106	GOL	O1-C1-C2-C3
3	H	101	GOL	O1-C1-C2-O2
3	H	101	GOL	O1-C1-C2-C3
3	H	104	GOL	C1-C2-C3-O3
3	H	105	GOL	C1-C2-C3-O3
9	D	524	PG4	C6-C5-O3-C4
5	F	108	PGE	O2-C3-C4-O3
9	B	512	PG4	O2-C3-C4-O3
9	C	529	PG4	O2-C3-C4-O3
9	D	522	PG4	O3-C5-C6-O4
9	D	522	PG4	O2-C3-C4-O3
4	G	108	PEG	O2-C3-C4-O4
5	D	520	PGE	O2-C3-C4-O3
5	E	105	PGE	O2-C3-C4-O3
5	C	528	PGE	O2-C3-C4-O3
5	D	521	PGE	O2-C3-C4-O3
3	D	501	GOL	O1-C1-C2-O2
3	D	508	GOL	O1-C1-C2-O2
3	F	101	GOL	O2-C2-C3-O3
3	F	103	GOL	O1-C1-C2-O2
5	C	528	PGE	C6-C5-O3-C4
4	F	107	PEG	O1-C1-C2-O2
5	A	505	PGE	O3-C5-C6-O4
5	F	108	PGE	O1-C1-C2-O2
4	D	514	PEG	C4-C3-O2-C2
5	C	528	PGE	C3-C4-O3-C5
5	G	110	PGE	O2-C3-C4-O3
4	A	503	PEG	O2-C3-C4-O4
4	A	504	PEG	O1-C1-C2-O2
4	A	504	PEG	O2-C3-C4-O4
4	C	522	PEG	O2-C3-C4-O4
4	D	514	PEG	O1-C1-C2-O2
4	D	515	PEG	O2-C3-C4-O4
4	G	109	PEG	O2-C3-C4-O4
5	G	110	PGE	O3-C5-C6-O4
5	E	107	PGE	O2-C3-C4-O3
4	D	514	PEG	C1-C2-O2-C3
9	D	523	PG4	O3-C5-C6-O4

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Mol	Chain	Res	Type	Atoms
3	B	501	GOL	O1-C1-C2-C3
3	B	502	GOL	O1-C1-C2-C3
3	B	504	GOL	C1-C2-C3-O3
3	B	507	GOL	C1-C2-C3-O3
3	B	509	GOL	C1-C2-C3-O3
3	B	510	GOL	O1-C1-C2-C3
3	C	501	GOL	O1-C1-C2-C3
3	C	505	GOL	C1-C2-C3-O3
3	C	506	GOL	C1-C2-C3-O3
3	C	508	GOL	C1-C2-C3-O3
3	C	512	GOL	C1-C2-C3-O3
3	C	514	GOL	O1-C1-C2-C3
3	C	515	GOL	O1-C1-C2-C3
3	C	516	GOL	C1-C2-C3-O3
3	C	517	GOL	C1-C2-C3-O3
3	C	519	GOL	C1-C2-C3-O3
3	D	504	GOL	C1-C2-C3-O3
3	D	505	GOL	O1-C1-C2-C3
3	D	508	GOL	C1-C2-C3-O3
3	E	103	GOL	O1-C1-C2-C3
3	E	104	GOL	O1-C1-C2-C3
3	F	101	GOL	C1-C2-C3-O3
3	F	104	GOL	O1-C1-C2-C3
3	G	102	GOL	O1-C1-C2-C3
3	H	102	GOL	C1-C2-C3-O3
3	H	103	GOL	C1-C2-C3-O3
3	H	105	GOL	O1-C1-C2-C3
4	D	517	PEG	O1-C1-C2-O2
4	D	519	PEG	O1-C1-C2-O2
5	D	520	PGE	O1-C1-C2-O2
5	F	108	PGE	O3-C5-C6-O4
9	D	523	PG4	O4-C7-C8-O5
9	D	524	PG4	C3-C4-O3-C5
5	C	524	PGE	C4-C3-O2-C2
3	A	502	GOL	O1-C1-C2-O2
3	B	501	GOL	O1-C1-C2-O2
3	C	502	GOL	O1-C1-C2-O2
3	C	505	GOL	O2-C2-C3-O3
3	C	508	GOL	O1-C1-C2-O2
3	C	511	GOL	O1-C1-C2-O2
3	C	512	GOL	O2-C2-C3-O3
3	C	516	GOL	O2-C2-C3-O3

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Mol	Chain	Res	Type	Atoms
3	D	503	GOL	O2-C2-C3-O3
3	D	504	GOL	O1-C1-C2-O2
3	D	505	GOL	O2-C2-C3-O3
3	D	508	GOL	O2-C2-C3-O3
3	D	511	GOL	O1-C1-C2-O2
3	D	511	GOL	O2-C2-C3-O3
3	D	513	GOL	O2-C2-C3-O3
3	E	104	GOL	O2-C2-C3-O3
3	F	101	GOL	O1-C1-C2-O2
3	F	102	GOL	O2-C2-C3-O3
3	F	104	GOL	O2-C2-C3-O3
3	F	106	GOL	O2-C2-C3-O3
3	H	104	GOL	O2-C2-C3-O3
3	H	105	GOL	O2-C2-C3-O3
4	G	107	PEG	C4-C3-O2-C2
4	C	521	PEG	O2-C3-C4-O4
4	D	518	PEG	O2-C3-C4-O4
9	D	524	PG4	O1-C1-C2-O2
9	D	524	PG4	O2-C3-C4-O3
9	D	522	PG4	C8-C7-O4-C6
4	D	517	PEG	O2-C3-C4-O4
5	C	526	PGE	O3-C5-C6-O4
9	C	529	PG4	O3-C5-C6-O4
4	D	516	PEG	O1-C1-C2-O2
4	F	107	PEG	O2-C3-C4-O4
5	C	524	PGE	O1-C1-C2-O2
5	C	525	PGE	O3-C5-C6-O4
5	C	527	PGE	O3-C5-C6-O4
5	C	528	PGE	O1-C1-C2-O2
5	C	525	PGE	O2-C3-C4-O3
3	B	503	GOL	O1-C1-C2-O2
3	C	507	GOL	O2-C2-C3-O3
3	C	518	GOL	O1-C1-C2-O2
3	F	102	GOL	O1-C1-C2-O2
3	F	104	GOL	O1-C1-C2-O2
3	H	105	GOL	O1-C1-C2-O2
5	E	106	PGE	O3-C5-C6-O4
3	B	509	GOL	O2-C2-C3-O3
3	C	501	GOL	O1-C1-C2-O2
3	C	519	GOL	O2-C2-C3-O3
3	D	501	GOL	O2-C2-C3-O3
3	D	512	GOL	O2-C2-C3-O3

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Mol	Chain	Res	Type	Atoms
3	E	104	GOL	O1-C1-C2-O2
3	G	102	GOL	O1-C1-C2-O2
5	E	105	PGE	O1-C1-C2-O2
5	C	524	PGE	O3-C5-C6-O4
3	B	504	GOL	O1-C1-C2-O2
3	B	510	GOL	O1-C1-C2-O2
3	C	508	GOL	O2-C2-C3-O3
3	D	509	GOL	O1-C1-C2-O2
5	F	109	PGE	C1-C2-O2-C3
4	D	516	PEG	O2-C3-C4-O4
5	D	521	PGE	C6-C5-O3-C4
4	F	107	PEG	C1-C2-O2-C3
5	C	526	PGE	C3-C4-O3-C5
9	C	529	PG4	C3-C4-O3-C5
9	D	522	PG4	C5-C6-O4-C7
5	A	505	PGE	C4-C3-O2-C2
9	C	529	PG4	C8-C7-O4-C6
9	D	523	PG4	C8-C7-O4-C6
4	G	108	PEG	O1-C1-C2-O2
5	A	505	PGE	O1-C1-C2-O2
5	C	527	PGE	O1-C1-C2-O2
5	D	520	PGE	O3-C5-C6-O4
5	E	106	PGE	C3-C4-O3-C5
5	F	108	PGE	C6-C5-O3-C4
5	G	110	PGE	C3-C4-O3-C5
5	C	525	PGE	C3-C4-O3-C5
5	D	520	PGE	C3-C4-O3-C5
3	C	502	GOL	O2-C2-C3-O3
3	C	517	GOL	O2-C2-C3-O3
3	D	507	GOL	O2-C2-C3-O3
5	E	107	PGE	O3-C5-C6-O4
9	C	529	PG4	C4-C3-O2-C2
5	C	527	PGE	O2-C3-C4-O3
4	D	516	PEG	C1-C2-O2-C3
4	A	503	PEG	O1-C1-C2-O2
3	D	502	GOL	O2-C2-C3-O3
3	C	504	GOL	O1-C1-C2-C3
5	C	524	PGE	O2-C3-C4-O3
4	G	109	PEG	O1-C1-C2-O2
9	D	522	PG4	C3-C4-O3-C5
3	B	501	GOL	O2-C2-C3-O3
3	B	507	GOL	O2-C2-C3-O3

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Mol	Chain	Res	Type	Atoms
3	C	515	GOL	O1-C1-C2-O2
3	D	504	GOL	O2-C2-C3-O3
4	C	521	PEG	C4-C3-O2-C2
4	A	504	PEG	C1-C2-O2-C3
4	D	518	PEG	C4-C3-O2-C2
9	D	522	PG4	C1-C2-O2-C3
3	A	502	GOL	O2-C2-C3-O3
3	B	504	GOL	O2-C2-C3-O3
3	B	510	GOL	O2-C2-C3-O3
3	C	504	GOL	O1-C1-C2-O2
4	G	107	PEG	O2-C3-C4-O4
5	C	528	PGE	O3-C5-C6-O4
9	C	529	PG4	C1-C2-O2-C3
9	D	523	PG4	C4-C3-O2-C2
3	C	513	GOL	O1-C1-C2-C3
3	D	509	GOL	O1-C1-C2-C3
5	D	520	PGE	C6-C5-O3-C4
5	E	106	PGE	C4-C3-O2-C2
5	D	521	PGE	C4-C3-O2-C2
10	F	110	TSA	C2-C1-C10-O1
10	G	111	TSA	C2-C1-C10-O1
5	C	525	PGE	C6-C5-O3-C4
4	C	523	PEG	O2-C3-C4-O4
5	F	109	PGE	O3-C5-C6-O4
9	C	529	PG4	O1-C1-C2-O2
10	E	108	TSA	C9-C1-C10-O2
5	C	524	PGE	C1-C2-O2-C3
5	C	526	PGE	O2-C3-C4-O3
5	A	505	PGE	C6-C5-O3-C4
3	G	106	GOL	O2-C2-C3-O3
3	H	102	GOL	O2-C2-C3-O3
5	A	505	PGE	C1-C2-O2-C3
4	G	109	PEG	C1-C2-O2-C3
4	G	109	PEG	C4-C3-O2-C2
8	A	509	TRP	OXT-C-CA-N
9	D	523	PG4	O2-C3-C4-O3

All (4) ring outliers are listed below:

Mol	Chain	Res	Type	Atoms
10	F	110	TSA	C1-C2-C3-C4-C5-C8-C9-O7
10	H	106	TSA	C1-C2-C3-C4-C5-C8-C9-O7

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Mol	Chain	Res	Type	Atoms
10	E	108	TSA	C1-C2-C3-C4-C5-C8-C9-O7
10	G	111	TSA	C1-C2-C3-C4-C5-C8-C9-O7

7 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	508	GOL	1	0
4	A	503	PEG	1	0
7	D	527	PO4	1	0
3	D	511	GOL	1	0
3	C	502	GOL	3	0
3	C	512	GOL	1	0
3	D	513	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	444/472 (94%)	0.37	28 (6%)	26	29	14, 30, 63, 99	1 (0%)
1	B	444/472 (94%)	0.50	32 (7%)	21	25	16, 34, 68, 93	1 (0%)
1	C	456/472 (96%)	0.67	50 (10%)	10	12	12, 25, 71, 111	1 (0%)
1	D	457/472 (96%)	0.50	36 (7%)	18	20	12, 24, 63, 100	2 (0%)
2	E	84/90 (93%)	0.79	8 (9%)	14	15	12, 33, 57, 81	1 (1%)
2	F	83/90 (92%)	0.94	11 (13%)	7	8	16, 28, 76, 102	1 (1%)
2	G	83/90 (92%)	0.59	10 (12%)	9	9	19, 28, 54, 71	0
2	H	83/90 (92%)	0.67	4 (4%)	35	40	18, 25, 63, 98	0
All	All	2134/2248 (94%)	0.55	179 (8%)	17	18	12, 29, 67, 111	7 (0%)

All (179) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	13	TRP	10.8
1	C	17	ILE	10.4
1	C	28	LEU	9.3
1	A	149	LEU	8.6
1	D	17	ILE	8.3
1	A	29	PRO	8.1
1	A	152	ASN	8.0
2	H	9	LEU	8.0
2	F	8	PRO	7.7
1	D	28	LEU	7.3
2	H	8	PRO	7.1
1	D	13	TRP	6.9
1	C	15	VAL	6.9
1	C	7	HIS	6.7
1	A	153	GLY	6.6
1	D	15	VAL	6.5

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Mol	Chain	Res	Type	RSRZ
1	C	149	LEU	6.5
1	D	9	SER	6.3
1	D	7	HIS	6.0
2	F	9	LEU	5.9
2	G	9	LEU	5.9
1	A	31	GLY	5.8
1	C	151	GLY	5.7
1	B	146	SER	5.7
2	G	8	PRO	5.7
2	E	7	ASP	5.5
2	E	8	PRO	5.4
1	C	9	SER	5.4
1	B	29	PRO	5.3
1	D	31	GLY	5.3
1	B	154	LEU	5.2
1	B	149	LEU	5.1
1	D	8	SER	5.0
1	C	31	GLY	4.9
1	D	10	GLY	4.8
1	C	8	SER	4.8
1	C	29	PRO	4.8
1	C	14	THR	4.7
1	A	154	LEU	4.6
1	C	30	GLU	4.6
1	A	146	SER	4.5
1	A	151	GLY	4.5
1	C	275	GLU	4.5
1	D	275	GLU	4.4
2	E	9	LEU	4.3
1	C	389	TYR	4.3
1	D	386	SER	4.2
1	A	148	ASP	4.1
1	B	151	GLY	4.1
1	C	277	GLY	4.1
1	D	14	THR	3.9
2	H	68	GLU	3.9
1	B	31	GLY	3.9
1	B	153	GLY	3.9
2	F	10	SER	3.9
2	E	10	SER	3.8
1	D	277	GLY	3.8
1	B	155	PRO	3.8

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Mol	Chain	Res	Type	RSRZ
1	C	146	SER	3.7
1	C	52	ARG	3.7
1	D	29	PRO	3.7
2	G	10	SER	3.7
1	A	274	ASP	3.7
1	B	246	GLU	3.6
1	C	11	MET	3.6
1	C	148	ASP	3.6
1	D	6	HIS	3.5
1	C	10	GLY	3.5
1	B	152	ASN	3.5
1	D	12	SER	3.4
1	A	32	MET	3.4
1	D	52	ARG	3.4
1	A	247	SER	3.4
1	C	150	ASP	3.4
1	B	384	THR	3.3
1	C	16	ASP	3.3
1	C	436	ASP	3.3
1	B	148	ASP	3.3
1	D	16	ASP	3.3
1	A	147	SER	3.2
1	D	244	SER	3.2
2	F	50	GLY	3.2
1	D	389	TYR	3.1
1	A	30	GLU	3.1
1	C	276	GLU	3.1
1	B	30	GLU	3.0
2	F	16	LYS	3.0
2	E	14	ILE	3.0
1	B	150	ASP	3.0
1	B	244	SER	3.0
1	D	276	GLU	3.0
1	B	247	SER	2.9
1	B	385	ALA	2.9
2	F	17	TYR	2.9
1	C	12	SER	2.9
1	C	244	SER	2.9
2	E	48	SER	2.8
1	C	53	ALA	2.8
1	D	149	LEU	2.8
1	C	147	SER	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	34	GLN	2.8
1	D	150	ASP	2.8
1	D	11	MET	2.8
1	B	147	SER	2.7
1	D	32	MET	2.7
1	B	171	ALA	2.7
1	C	154	LEU	2.7
1	C	249	ARG	2.7
1	C	152	ASN	2.7
1	C	386	SER	2.7
1	D	146	SER	2.7
1	A	150	ASP	2.7
1	A	245	ASP	2.7
2	G	14	ILE	2.7
1	C	388	GLY	2.7
1	B	49	THR	2.6
1	C	167	ALA	2.6
1	D	247	SER	2.6
2	F	23	ARG	2.6
1	A	409[A]	ARG	2.6
1	D	30	GLU	2.6
1	D	434	ALA	2.6
1	A	155	PRO	2.6
1	A	57	ASN	2.5
1	C	36	PHE	2.5
1	C	40	ILE	2.5
2	F	11	ASP	2.5
1	A	244	SER	2.5
1	B	32	MET	2.5
1	A	229	ARG	2.5
1	A	275	GLU	2.4
1	C	33	GLN	2.4
2	E	13	GLU	2.4
1	B	409[A]	ARG	2.4
1	B	440	VAL	2.4
1	D	151	GLY	2.4
1	C	165	VAL	2.4
1	C	243	VAL	2.4
1	C	229	ARG	2.4
1	A	278	ASN	2.4
2	G	12	ALA	2.3
1	C	34	GLN	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	383	PHE	2.3
1	B	245	ASP	2.3
2	G	48	SER	2.3
2	F	68	GLU	2.3
1	C	144	PRO	2.3
1	C	443	PRO	2.3
1	C	60	LYS	2.3
1	C	209	GLU	2.3
1	D	148	ASP	2.3
1	B	388	GLY	2.3
1	D	153	GLY	2.3
1	C	246	GLU	2.2
1	D	209	GLU	2.2
1	B	165	VAL	2.2
1	B	167	ALA	2.2
1	A	34	GLN	2.2
1	C	328	ALA	2.2
1	B	103	SER	2.2
1	B	437	ILE	2.2
1	D	36	PHE	2.2
2	E	68	GLU	2.2
2	H	65	GLN	2.1
2	G	34	ARG	2.1
1	A	277	GLY	2.1
2	F	67	ARG	2.1
1	C	166	GLU	2.1
2	F	15	GLN	2.1
1	A	383	PHE	2.1
1	B	156	ASN	2.1
1	D	245	ASP	2.1
1	D	37	GLU	2.0
2	G	13	GLU	2.0
2	G	11	ASP	2.0
2	G	15	GLN	2.0
1	A	36	PHE	2.0
1	A	366	ALA	2.0
1	C	278	ASN	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 [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
3	GOL	C	520	6/6	0.56	0.21	46,49,51,52	0
3	GOL	C	519	6/6	0.57	0.19	54,56,58,58	0
5	PGE	E	106	10/10	0.62	0.27	63,66,68,69	0
3	GOL	D	507	6/6	0.65	0.25	57,61,63,64	0
3	GOL	C	518	6/6	0.65	0.23	53,53,55,55	0
4	PEG	D	516	7/7	0.66	0.22	46,52,58,59	0
3	GOL	G	103	6/6	0.67	0.20	57,61,62,65	0
5	PGE	E	105	10/10	0.67	0.21	58,64,65,66	0
3	GOL	H	105	6/6	0.67	0.20	54,56,58,59	0
8	TRP	D	528	15/15	0.69	0.20	55,57,61,63	0
3	GOL	B	505	6/6	0.70	0.18	54,54,56,58	0
3	GOL	D	504	6/6	0.70	0.21	55,60,61,62	0
5	PGE	F	109	10/10	0.70	0.20	56,61,64,65	0
3	GOL	C	503	6/6	0.70	0.22	55,56,57,59	0
3	GOL	C	513	6/6	0.71	0.22	54,58,58,59	0
4	PEG	F	107	7/7	0.71	0.19	42,48,50,51	0
3	GOL	D	508	6/6	0.72	0.26	49,56,57,61	0
3	GOL	E	103	6/6	0.72	0.17	56,58,60,60	0
3	GOL	G	104	6/6	0.73	0.17	49,54,56,57	0
3	GOL	D	511	6/6	0.73	0.21	50,51,55,55	0
3	GOL	D	512	6/6	0.74	0.20	51,56,59,59	0
3	GOL	B	506	6/6	0.74	0.18	60,65,66,66	0
5	PGE	E	107	10/10	0.74	0.21	56,63,65,66	0
3	GOL	F	105	6/6	0.74	0.19	50,53,53,54	0
3	GOL	C	507	6/6	0.74	0.16	39,45,49,50	0
4	PEG	G	109	7/7	0.75	0.18	49,53,57,57	0
3	GOL	C	515	6/6	0.75	0.20	57,62,63,63	0
4	PEG	C	523	7/7	0.75	0.18	46,49,50,51	0
4	PEG	D	515	7/7	0.75	0.18	41,42,45,45	0
3	GOL	C	511	6/6	0.75	0.19	60,62,63,64	0
5	PGE	G	110	10/10	0.75	0.20	45,65,69,70	0
3	GOL	F	101	6/6	0.75	0.16	48,51,52,52	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
9	PG4	B	512	13/13	0.75	0.19	50,57,60,61	0
8	TRP	A	509	15/15	0.76	0.19	58,61,64,67	0
8	TRP	C	532	15/15	0.76	0.18	47,51,54,55	0
5	PGE	A	505	10/10	0.76	0.21	53,65,67,68	0
3	GOL	B	509	6/6	0.76	0.20	59,64,66,67	0
9	PG4	C	529	13/13	0.76	0.20	48,53,61,62	0
3	GOL	C	517	6/6	0.77	0.22	58,65,68,69	0
4	PEG	C	522	7/7	0.78	0.15	39,43,46,49	0
4	PEG	G	108	7/7	0.78	0.18	49,51,53,54	0
3	GOL	F	106	6/6	0.78	0.16	45,47,48,49	0
3	GOL	C	508	6/6	0.78	0.21	46,51,53,54	0
5	PGE	C	524	10/10	0.78	0.23	53,66,70,71	0
5	PGE	C	527	10/10	0.78	0.19	44,51,58,58	0
4	PEG	A	504	7/7	0.78	0.19	52,53,59,64	0
4	PEG	D	517	7/7	0.78	0.21	52,56,58,59	0
9	PG4	D	524	13/13	0.78	0.19	51,52,60,63	0
3	GOL	H	101	6/6	0.79	0.20	50,54,56,57	0
3	GOL	D	505	6/6	0.79	0.17	38,41,43,43	0
3	GOL	C	505	6/6	0.79	0.22	43,49,50,52	0
3	GOL	C	506	6/6	0.79	0.15	57,59,64,71	0
3	GOL	C	501	6/6	0.79	0.15	35,37,39,41	0
3	GOL	B	502	6/6	0.79	0.17	51,56,57,59	0
5	PGE	D	521	10/10	0.79	0.17	51,54,57,58	0
3	GOL	E	102	6/6	0.79	0.15	48,54,57,58	0
3	GOL	G	105	6/6	0.79	0.17	53,58,59,59	0
5	PGE	F	108	10/10	0.80	0.15	43,50,56,58	0
3	GOL	H	102	6/6	0.80	0.14	38,42,43,47	0
3	GOL	D	503	6/6	0.80	0.17	42,45,46,47	0
5	PGE	C	525	10/10	0.81	0.18	42,49,55,56	0
4	PEG	D	514	7/7	0.81	0.16	47,49,50,51	0
5	PGE	D	520	10/10	0.81	0.16	43,51,53,54	0
4	PEG	G	107	7/7	0.81	0.17	41,45,49,51	0
3	GOL	B	504	6/6	0.81	0.17	57,62,64,65	0
3	GOL	B	510	6/6	0.81	0.18	55,60,63,65	0
3	GOL	B	507	6/6	0.81	0.17	50,51,52,55	0
9	PG4	D	522	13/13	0.81	0.18	28,35,56,56	0
4	PEG	D	518	7/7	0.81	0.17	47,51,53,53	0
5	PGE	C	526	10/10	0.82	0.17	39,47,50,51	0
3	GOL	B	503	6/6	0.82	0.17	44,50,55,59	0
7	PO4	D	527	5/5	0.82	0.15	52,56,59,61	0
3	GOL	B	501	6/6	0.82	0.16	45,49,50,50	0
9	PG4	D	523	13/13	0.82	0.19	57,62,67,67	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	GOL	G	101	6/6	0.82	0.18	40,45,45,47	0
8	TRP	B	515	15/15	0.83	0.18	62,64,65,65	0
3	GOL	F	104	6/6	0.83	0.11	32,37,39,45	0
3	GOL	H	103	6/6	0.83	0.16	46,53,54,55	0
3	GOL	A	502	6/6	0.83	0.14	44,48,48,49	0
5	PGE	C	528	10/10	0.83	0.15	43,47,48,48	0
3	GOL	E	104	6/6	0.83	0.15	50,52,53,53	0
4	PEG	C	521	7/7	0.83	0.21	34,41,46,47	0
3	GOL	C	502	6/6	0.83	0.17	36,39,40,42	0
3	GOL	C	509	6/6	0.84	0.22	54,56,57,58	0
3	GOL	C	510	6/6	0.84	0.20	58,60,65,66	0
3	GOL	E	101	6/6	0.84	0.12	53,55,56,57	0
4	PEG	D	519	7/7	0.84	0.13	44,47,47,48	0
3	GOL	C	514	6/6	0.85	0.11	41,43,43,46	0
3	GOL	G	106	6/6	0.85	0.18	67,68,68,68	0
3	GOL	D	510	6/6	0.85	0.14	47,49,55,58	0
3	GOL	C	504	6/6	0.85	0.15	56,57,58,60	0
3	GOL	D	506	6/6	0.86	0.18	41,45,47,48	0
7	PO4	A	508	5/5	0.86	0.09	68,68,70,72	0
3	GOL	C	512	6/6	0.86	0.18	57,61,63,65	0
3	GOL	G	102	6/6	0.87	0.14	31,34,35,36	0
3	GOL	B	508	6/6	0.87	0.17	47,51,51,52	0
3	GOL	D	501	6/6	0.87	0.13	34,37,39,41	0
3	GOL	D	502	6/6	0.87	0.13	38,42,44,46	0
3	GOL	F	103	6/6	0.88	0.14	47,48,51,53	0
3	GOL	H	104	6/6	0.88	0.10	31,37,37,39	0
3	GOL	C	516	6/6	0.88	0.11	50,55,56,57	0
3	GOL	D	509	6/6	0.88	0.12	43,47,49,50	0
3	GOL	D	513	6/6	0.88	0.12	36,41,42,44	0
3	GOL	A	501	6/6	0.88	0.14	48,49,56,59	0
3	GOL	F	102	6/6	0.88	0.12	35,36,38,38	0
4	PEG	A	503	7/7	0.89	0.17	33,36,42,42	0
4	PEG	B	511	7/7	0.89	0.17	40,40,44,44	0
10	TSA	E	108	16/16	0.91	0.09	31,33,34,35	0
10	TSA	G	111	16/16	0.92	0.08	27,28,29,29	0
10	TSA	H	106	16/16	0.95	0.07	15,17,22,23	0
10	TSA	F	110	16/16	0.96	0.06	16,18,23,24	0
6	MN	B	513	1/1	0.98	0.05	40,40,40,40	0
6	MN	A	506	1/1	0.99	0.05	40,40,40,40	0
7	PO4	B	514	5/5	0.99	0.06	33,34,35,35	0
7	PO4	C	531	5/5	0.99	0.06	28,29,30,30	0
7	PO4	D	526	5/5	0.99	0.07	26,27,28,28	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
6	MN	D	525	1/1	0.99	0.02	27,27,27,27	0
7	PO4	A	507	5/5	0.99	0.05	31,32,33,34	0
6	MN	C	530	1/1	1.00	0.01	29,29,29,29	0

6.5 Other polymers [i](#)

There are no such residues in this entry.