



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 03:43 PM UTC

PDB ID : 7KG5 / pdb_00007kg5
Title : Dihydrodipicolinate synthase (DHDPS) from C.jejuni, H56W mutant with pyruvate bound in the active site
Authors : Saran, S.; Majdi Yazdi, M.; Sanders, D.A.R.
Deposited on : 2020-10-15
Resolution : 1.95 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:

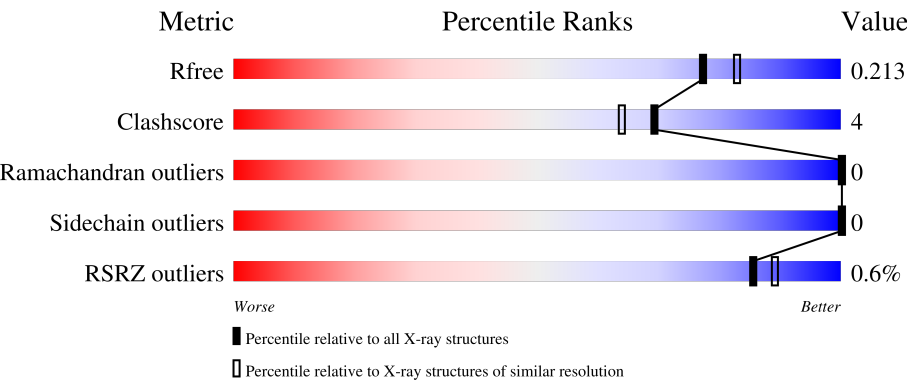
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (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 1.95 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	3494 (1.96-1.96)
Clashscore	190562	3612 (1.96-1.96)
Ramachandran outliers	187476	3587 (1.96-1.96)
Sidechain outliers	187428	3587 (1.96-1.96)
RSRZ outliers	180081	3495 (1.96-1.96)

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	310	87%9%.
1	B	310	93%. 5%
1	C	310	% 89%6%5%
1	D	310	% 90%5%5%
1	E	310	89%6%5%

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Mol	Chain	Length	Quality of chain
1	F	310	% 86% 9% .

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	EDO	C	304	-	-	X	-
4	PEG	A	310	-	-	X	-
4	PEG	F	305	-	-	X	-
5	ACT	D	307	-	-	X	-

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 15020 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 4-hydroxy-tetrahydrodipicolinate synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	297	Total	C	N	O	S	0	2	0
			2306	1469	381	443	13			
1	B	296	Total	C	N	O	S	0	1	0
			2298	1466	380	439	13			
1	C	296	Total	C	N	O	S	0	1	0
			2298	1466	380	439	13			
1	D	296	Total	C	N	O	S	0	2	0
			2286	1455	378	440	13			
1	E	296	Total	C	N	O	S	0	1	0
			2285	1455	378	439	13			
1	F	297	Total	C	N	O	S	0	3	0
			2296	1460	379	444	13			

There are 78 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-11	MET	-	expression tag	UNP Q9PPB4
A	-10	ARG	-	expression tag	UNP Q9PPB4
A	-9	GLY	-	expression tag	UNP Q9PPB4
A	-8	SER	-	expression tag	UNP Q9PPB4
A	-7	HIS	-	expression tag	UNP Q9PPB4
A	-6	HIS	-	expression tag	UNP Q9PPB4
A	-5	HIS	-	expression tag	UNP Q9PPB4
A	-4	HIS	-	expression tag	UNP Q9PPB4
A	-3	HIS	-	expression tag	UNP Q9PPB4
A	-2	HIS	-	expression tag	UNP Q9PPB4
A	-1	GLY	-	expression tag	UNP Q9PPB4
A	0	SER	-	expression tag	UNP Q9PPB4
A	56	TRP	HIS	engineered mutation	UNP Q9PPB4
B	-11	MET	-	expression tag	UNP Q9PPB4
B	-10	ARG	-	expression tag	UNP Q9PPB4
B	-9	GLY	-	expression tag	UNP Q9PPB4
B	-8	SER	-	expression tag	UNP Q9PPB4

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-7	HIS	-	expression tag	UNP Q9PPB4
B	-6	HIS	-	expression tag	UNP Q9PPB4
B	-5	HIS	-	expression tag	UNP Q9PPB4
B	-4	HIS	-	expression tag	UNP Q9PPB4
B	-3	HIS	-	expression tag	UNP Q9PPB4
B	-2	HIS	-	expression tag	UNP Q9PPB4
B	-1	GLY	-	expression tag	UNP Q9PPB4
B	0	SER	-	expression tag	UNP Q9PPB4
B	56	TRP	HIS	engineered mutation	UNP Q9PPB4
C	-11	MET	-	expression tag	UNP Q9PPB4
C	-10	ARG	-	expression tag	UNP Q9PPB4
C	-9	GLY	-	expression tag	UNP Q9PPB4
C	-8	SER	-	expression tag	UNP Q9PPB4
C	-7	HIS	-	expression tag	UNP Q9PPB4
C	-6	HIS	-	expression tag	UNP Q9PPB4
C	-5	HIS	-	expression tag	UNP Q9PPB4
C	-4	HIS	-	expression tag	UNP Q9PPB4
C	-3	HIS	-	expression tag	UNP Q9PPB4
C	-2	HIS	-	expression tag	UNP Q9PPB4
C	-1	GLY	-	expression tag	UNP Q9PPB4
C	0	SER	-	expression tag	UNP Q9PPB4
C	56	TRP	HIS	engineered mutation	UNP Q9PPB4
D	-11	MET	-	expression tag	UNP Q9PPB4
D	-10	ARG	-	expression tag	UNP Q9PPB4
D	-9	GLY	-	expression tag	UNP Q9PPB4
D	-8	SER	-	expression tag	UNP Q9PPB4
D	-7	HIS	-	expression tag	UNP Q9PPB4
D	-6	HIS	-	expression tag	UNP Q9PPB4
D	-5	HIS	-	expression tag	UNP Q9PPB4
D	-4	HIS	-	expression tag	UNP Q9PPB4
D	-3	HIS	-	expression tag	UNP Q9PPB4
D	-2	HIS	-	expression tag	UNP Q9PPB4
D	-1	GLY	-	expression tag	UNP Q9PPB4
D	0	SER	-	expression tag	UNP Q9PPB4
D	56	TRP	HIS	engineered mutation	UNP Q9PPB4
E	-11	MET	-	expression tag	UNP Q9PPB4
E	-10	ARG	-	expression tag	UNP Q9PPB4
E	-9	GLY	-	expression tag	UNP Q9PPB4
E	-8	SER	-	expression tag	UNP Q9PPB4
E	-7	HIS	-	expression tag	UNP Q9PPB4
E	-6	HIS	-	expression tag	UNP Q9PPB4
E	-5	HIS	-	expression tag	UNP Q9PPB4

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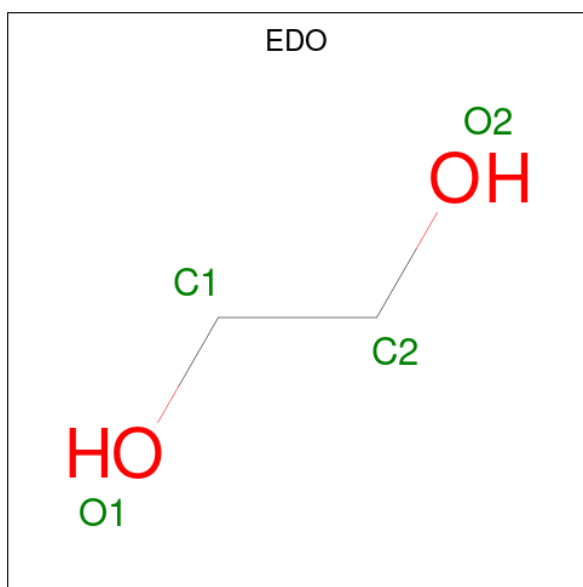
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Chain	Residue	Modelled	Actual	Comment	Reference
E	-4	HIS	-	expression tag	UNP Q9PPB4
E	-3	HIS	-	expression tag	UNP Q9PPB4
E	-2	HIS	-	expression tag	UNP Q9PPB4
E	-1	GLY	-	expression tag	UNP Q9PPB4
E	0	SER	-	expression tag	UNP Q9PPB4
E	56	TRP	HIS	engineered mutation	UNP Q9PPB4
F	-11	MET	-	expression tag	UNP Q9PPB4
F	-10	ARG	-	expression tag	UNP Q9PPB4
F	-9	GLY	-	expression tag	UNP Q9PPB4
F	-8	SER	-	expression tag	UNP Q9PPB4
F	-7	HIS	-	expression tag	UNP Q9PPB4
F	-6	HIS	-	expression tag	UNP Q9PPB4
F	-5	HIS	-	expression tag	UNP Q9PPB4
F	-4	HIS	-	expression tag	UNP Q9PPB4
F	-3	HIS	-	expression tag	UNP Q9PPB4
F	-2	HIS	-	expression tag	UNP Q9PPB4
F	-1	GLY	-	expression tag	UNP Q9PPB4
F	0	SER	-	expression tag	UNP Q9PPB4
F	56	TRP	HIS	engineered mutation	UNP Q9PPB4

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	2	Total Mg 2 2	0	0
2	B	3	Total Mg 3 3	0	0
2	C	2	Total Mg 2 2	0	0
2	D	2	Total Mg 2 2	0	0
2	E	2	Total Mg 2 2	0	0
2	F	2	Total Mg 2 2	0	0

- Molecule 3 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



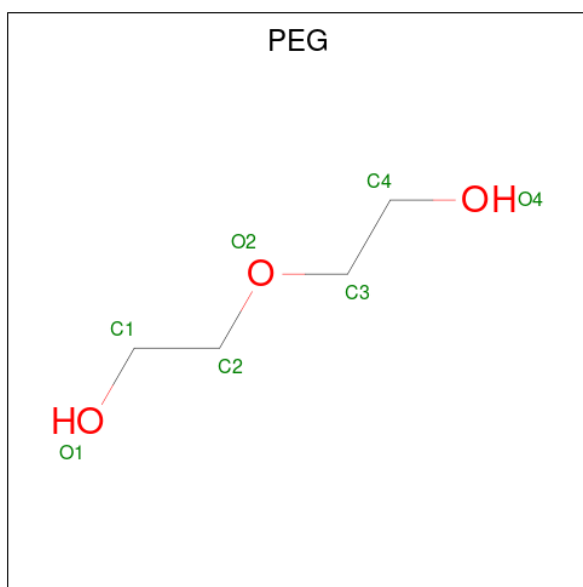
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		
3	C	1	Total	C	O	0	0
			4	2	2		

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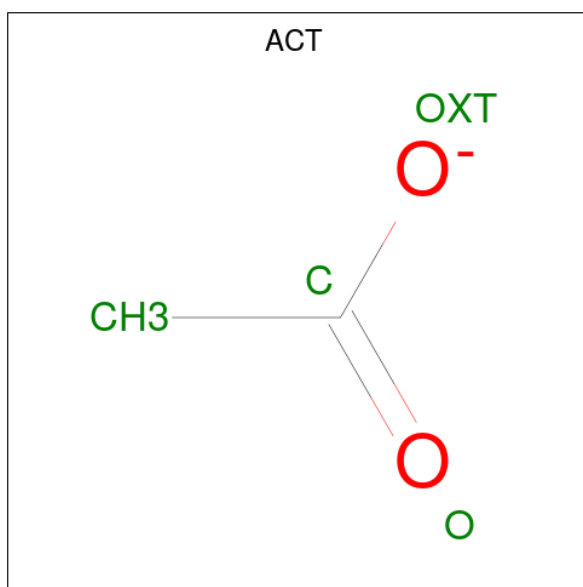
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	C	1	Total	C	O	0	0
			4	2	2		
3	C	1	Total	C	O	0	0
			4	2	2		
3	C	1	Total	C	O	0	0
			4	2	2		
3	C	1	Total	C	O	0	0
			4	2	2		
3	D	1	Total	C	O	0	0
			4	2	2		
3	D	1	Total	C	O	0	0
			4	2	2		
3	E	1	Total	C	O	0	0
			4	2	2		
3	E	1	Total	C	O	0	0
			4	2	2		
3	E	1	Total	C	O	0	0
			4	2	2		
3	E	1	Total	C	O	0	0
			4	2	2		
3	E	1	Total	C	O	0	0
			4	2	2		
3	F	1	Total	C	O	0	0
			4	2	2		
3	F	1	Total	C	O	0	0
			4	2	2		

- Molecule 4 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: C₄H₁₀O₃).



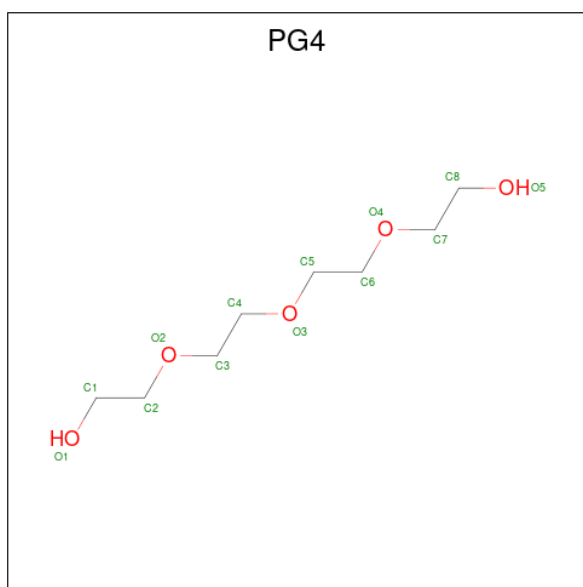
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			7	4	3		
4	A	1	Total	C	O	0	0
			7	4	3		
4	C	1	Total	C	O	0	0
			7	4	3		
4	C	1	Total	C	O	0	0
			7	4	3		
4	C	1	Total	C	O	0	0
			7	4	3		
4	D	1	Total	C	O	0	0
			7	4	3		
4	D	1	Total	C	O	0	0
			7	4	3		
4	F	1	Total	C	O	0	0
			7	4	3		
4	F	1	Total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

- Molecule 5 is ACETATE ION (CCD ID: ACT) (formula: $C_2H_3O_2$).



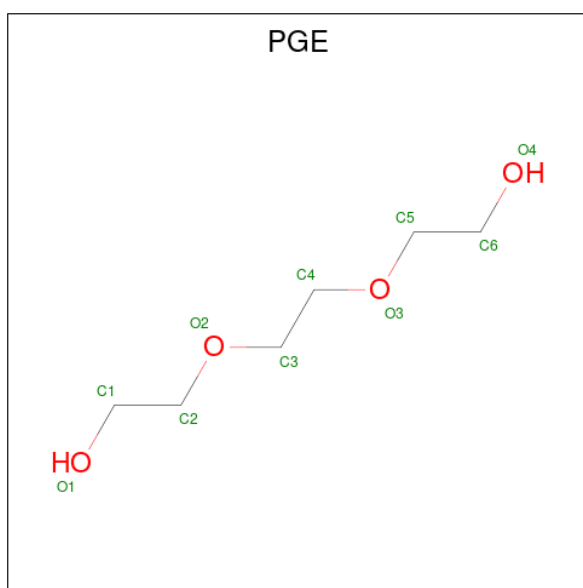
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			4	2	2		
5	C	1	Total	C	O	0	0
			4	2	2		
5	C	1	Total	C	O	0	0
			4	2	2		
5	D	1	Total	C	O	0	0
			4	2	2		
5	D	1	Total	C	O	0	0
			4	2	2		
5	E	1	Total	C	O	0	0
			4	2	2		
5	F	1	Total	C	O	0	0
			4	2	2		
5	F	1	Total	C	O	0	0
			4	2	2		

- Molecule 6 is TETRAETHYLENE GLYCOL (CCD ID: PG4) (formula: C₈H₁₈O₅).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			13	8	5		
6	D	1	Total	C	O	0	0
			13	8	5		

- Molecule 7 is TRIETHYLENE GLYCOL (CCD ID: PGE) (formula: $C_6H_{14}O_4$).



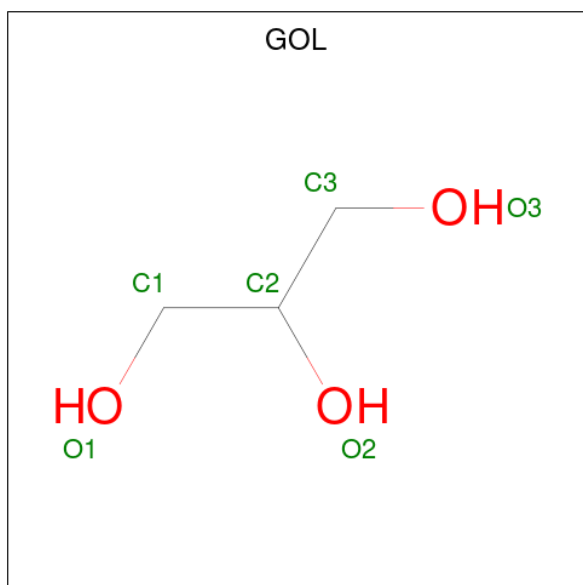
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	B	1	Total	C	O	0	0
			10	6	4		
7	B	1	Total	C	O	0	0
			10	6	4		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	C	1	Total	C	O	0	0
			10	6	4		
7	E	1	Total	C	O	0	0
			10	6	4		

- Molecule 8 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	161	Total	O	0	0
			161	161		
9	B	159	Total	O	0	0
			159	159		
9	C	155	Total	O	0	0
			155	155		

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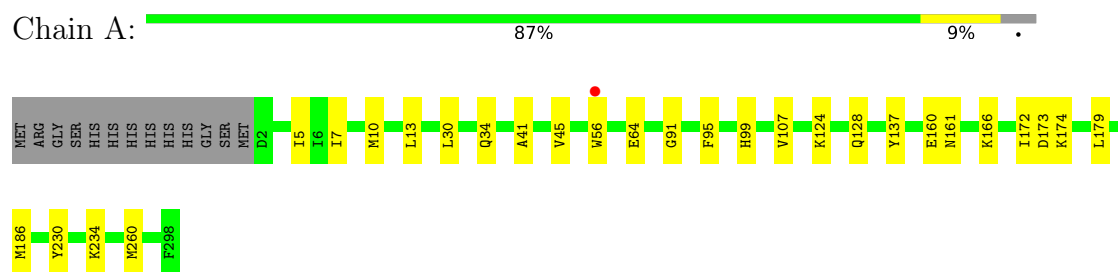
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	D	153	Total 153	O 153	0	0
9	E	153	Total 153	O 153	0	0
9	F	160	Total 160	O 160	0	0

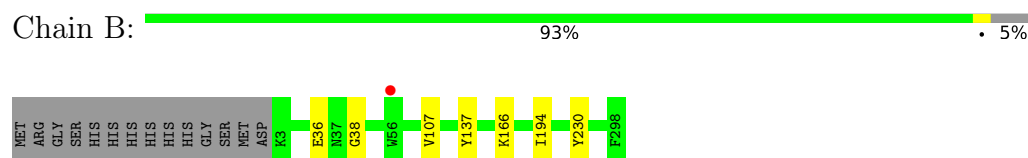
3 Residue-property plots [i](#)

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.

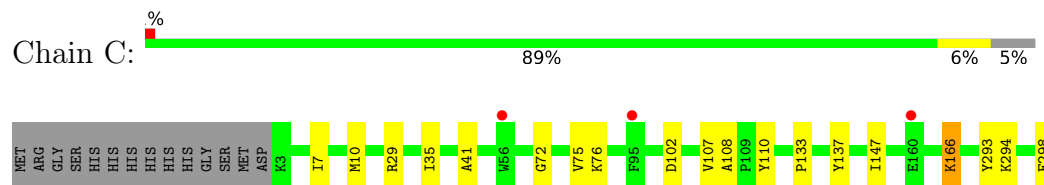
- Molecule 1: 4-hydroxy-tetrahydrodipicolinate synthase



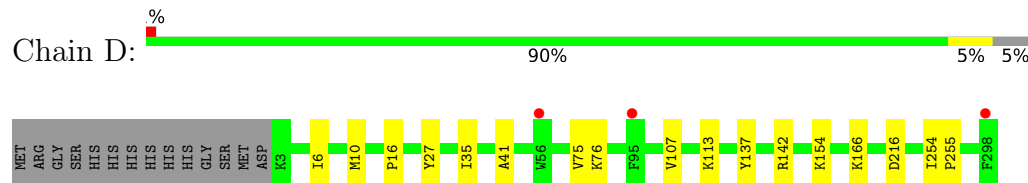
- Molecule 1: 4-hydroxy-tetrahydrodipicolinate synthase



- Molecule 1: 4-hydroxy-tetrahydrodipicolinate synthase



- Molecule 1: 4-hydroxy-tetrahydrodipicolinate synthase

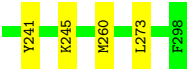
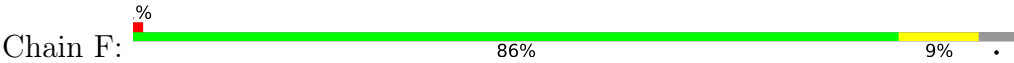


- Molecule 1: 4-hydroxy-tetrahydrodipicolinate synthase





● Molecule 1: 4-hydroxy-tetrahydronicotinate synthase



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	85.13Å 231.25Å 202.10Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.30 – 1.95 46.30 – 1.95	Depositor EDS
% Data completeness (in resolution range)	98.1 (46.30-1.95) 98.1 (46.30-1.95)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.46 (at 1.95Å)	Xtriage
Refinement program	PHENIX dev_2398	Depositor
R, R_{free}	0.174 , 0.210 0.178 , 0.213	Depositor DCC
R_{free} test set	7105 reflections (4.91%)	wwPDB-VP
Wilson B-factor (Å ²)	29.5	Xtriage
Anisotropy	0.274	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 35.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	15020	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 2.31% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: PEG, KPI, GOL, EDO, MG, PGE, ACT, PG4

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	0.37	2/2339 (0.1%)	0.49	1/3163 (0.0%)
1	B	0.30	0/2325	0.49	0/3144
1	C	0.28	0/2325	0.46	0/3144
1	D	0.35	0/2321	0.48	0/3137
1	E	0.29	0/2315	0.47	0/3129
1	F	0.29	0/2336	0.50	0/3159
All	All	0.31	2/13961 (0.0%)	0.48	1/18876 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	160	GLU	C-O	-6.34	1.16	1.24
1	A	161	ASN	C-O	-5.14	1.17	1.24

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	161	ASN	N-CA-C	5.21	117.87	111.82

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2306	0	2337	29	0
1	B	2298	0	2327	6	0
1	C	2298	0	2331	15	0
1	D	2286	0	2324	11	0
1	E	2285	0	2322	15	0
1	F	2296	0	2328	23	0
2	A	2	0	0	0	0
2	B	3	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0
2	E	2	0	0	0	0
2	F	2	0	0	0	0
3	A	24	0	36	7	0
3	B	28	0	42	4	0
3	C	20	0	30	4	0
3	D	8	0	12	0	0
3	E	20	0	30	7	0
3	F	8	0	12	1	0
4	A	14	0	20	9	0
4	C	21	0	30	1	0
4	D	14	0	20	0	0
4	F	14	0	20	6	0
5	A	4	0	3	1	0
5	C	8	0	6	0	0
5	D	8	0	6	2	0
5	E	4	0	3	0	0
5	F	12	0	9	1	0
6	A	13	0	18	0	0
6	D	13	0	18	1	0
7	B	20	0	28	2	0
7	C	10	0	14	0	0
7	E	10	0	14	0	0
8	B	6	0	8	0	0
8	C	6	0	8	0	0
8	E	12	0	16	1	0
9	A	161	0	0	0	0
9	B	159	0	0	1	0
9	C	155	0	0	2	0
9	D	153	0	0	1	0
9	E	153	0	0	1	0
9	F	160	0	0	4	0
All	All	15020	0	14372	100	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 (100) 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:F:159:CYS:HA	4:F:305:PEG:H21	1.36	1.07
1:C:166:KPI:H1A	9:C:460:HOH:O	1.58	1.04
1:A:30:LEU:HB3	4:A:310:PEG:H31	1.60	0.84
1:F:5:ILE:HD12	1:F:186:MET:HE2	1.66	0.78
1:A:7:ILE:HG21	1:A:186:MET:HE3	1.66	0.76
1:E:5:ILE:HD12	1:E:186:MET:HE2	1.68	0.74
1:C:76:LYS:HD2	3:C:304:EDO:H22	1.70	0.73
1:E:7:ILE:HG21	1:E:186:MET:HE3	1.69	0.73
1:F:166:KPI:H1A	9:F:428:HOH:O	1.90	0.72
1:A:7:ILE:HG21	1:A:186:MET:CE	2.19	0.71
1:F:7:ILE:HG21	1:F:186:MET:HE3	1.71	0.71
1:A:64:GLU:HG2	3:A:308:EDO:H12	1.73	0.70
1:A:107:VAL:HA	1:A:137:TYR:HB3	1.76	0.68
1:A:34:GLN:HG2	4:A:310:PEG:H11	1.75	0.68
1:D:107:VAL:HA	1:D:137:TYR:HB3	1.76	0.66
1:D:142:ARG:HE	5:D:307:ACT:H2	1.59	0.65
1:B:107:VAL:HA	1:B:137:TYR:HB3	1.79	0.65
1:F:142:ARG:HE	5:F:308:ACT:H2	1.61	0.65
1:E:107:VAL:HA	1:E:137:TYR:HB3	1.78	0.64
1:A:5:ILE:HD12	1:A:186:MET:HE2	1.80	0.64
1:C:107:VAL:HA	1:C:137:TYR:HB3	1.79	0.63
1:F:107:VAL:HA	1:F:137:TYR:HB3	1.79	0.63
1:D:154:LYS:HD3	6:D:309:PG4:H12	1.81	0.63
1:F:7:ILE:HG21	1:F:186:MET:CE	2.28	0.63
1:A:30:LEU:HB3	4:A:310:PEG:H12	1.80	0.62
1:A:13:LEU:HD22	4:A:310:PEG:H42	1.84	0.60
1:E:7:ILE:HG21	1:E:186:MET:CE	2.33	0.59
1:C:35:ILE:HG12	1:C:75:VAL:HG21	1.84	0.58
1:C:7:ILE:H	3:C:304:EDO:H21	1.69	0.57
1:C:7:ILE:HG13	3:E:304:EDO:H11	1.86	0.57
1:A:45:VAL:O	3:A:303:EDO:O1	2.25	0.55
7:B:302:PGE:O4	3:B:307:EDO:O2	2.26	0.54
1:A:260:MET:CE	4:A:310:PEG:H21	2.38	0.53
1:D:113:LYS:NZ	9:D:405:HOH:O	2.41	0.53
1:B:38:GLY:HA2	3:B:307:EDO:H11	1.91	0.53
1:E:24:GLU:HB2	3:E:306:EDO:H22	1.90	0.52
1:A:230:TYR:HD2	1:B:230:TYR:HD2	1.58	0.52
1:F:128:GLN:HG2	4:F:305:PEG:H32	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:174:LYS:HA	3:A:305:EDO:H11	1.91	0.51
1:A:230:TYR:HD2	1:B:230:TYR:CD2	2.28	0.51
3:E:305:EDO:H21	1:F:273:LEU:HD21	1.93	0.51
1:F:214:LEU:HB3	1:F:217:MET:HE3	1.94	0.50
1:F:97:LYS:NZ	9:F:408:HOH:O	2.44	0.50
1:F:56:TRP:CE3	1:F:57:GLU:HG2	2.46	0.49
3:B:307:EDO:H22	9:B:401:HOH:O	2.12	0.49
1:E:23:ASP:HA	3:E:306:EDO:H11	1.95	0.49
1:E:3:LYS:N	9:E:406:HOH:O	2.46	0.48
1:F:74:LYS:NZ	9:F:410:HOH:O	2.46	0.48
1:F:56:TRP:CZ2	1:F:95:PHE:CE1	3.01	0.48
1:C:7:ILE:N	3:C:304:EDO:H21	2.28	0.47
1:F:56:TRP:HE3	1:F:57:GLU:HG2	1.78	0.47
1:C:10:MET:HG2	1:C:41:ALA:HB3	1.97	0.47
1:C:166:KPI:C1	9:C:460:HOH:O	2.38	0.47
1:D:6:ILE:HG12	1:D:76:LYS:HD3	1.96	0.47
1:C:293:TYR:C	1:C:294:LYS:HD3	2.40	0.46
1:F:127:ALA:O	4:F:306:PEG:H32	2.15	0.46
1:A:7:ILE:HG21	1:A:186:MET:HE2	1.97	0.46
1:A:230:TYR:HE2	5:A:311:ACT:H1	1.81	0.46
1:F:60:ARG:HB2	1:F:95:PHE:HZ	1.81	0.46
1:A:99:HIS:O	3:A:308:EDO:H11	2.16	0.45
1:F:117:GLN:HB2	3:F:304:EDO:H11	1.98	0.45
1:D:10:MET:HG2	1:D:41:ALA:HB3	1.97	0.45
1:A:260:MET:HE2	4:A:310:PEG:H21	1.97	0.45
3:C:304:EDO:H12	3:E:304:EDO:O1	2.16	0.45
1:E:287:GLU:HG3	3:E:307:EDO:H11	1.97	0.45
4:F:306:PEG:H41	9:F:441:HOH:O	2.17	0.45
1:E:24:GLU:H	3:E:306:EDO:H11	1.82	0.44
1:C:108:ALA:HB2	1:C:147:ILE:HD11	2.00	0.44
1:B:36:GLU:O	7:B:302:PGE:H62	2.18	0.44
1:A:234:LYS:HD3	3:B:311:EDO:H11	2.00	0.43
1:A:172:ILE:HG23	1:B:194:ILE:HG21	2.01	0.43
1:C:102:ASP:O	1:C:133:PRO:HD2	2.19	0.43
1:A:91:GLY:HA3	4:A:309:PEG:H31	2.00	0.43
1:A:260:MET:HE1	4:A:310:PEG:H21	2.01	0.43
1:E:177:ASP:OD1	8:E:310:GOL:H32	2.19	0.42
1:A:174:LYS:CA	3:A:305:EDO:H11	2.49	0.42
4:F:305:PEG:H41	4:F:305:PEG:H22	1.50	0.42
1:E:102:ASP:O	1:E:133:PRO:HD2	2.19	0.42
1:A:179:LEU:HD22	3:A:307:EDO:H21	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:241:TYR:CE2	1:F:245:LYS:HG3	2.53	0.42
1:D:35:ILE:HG12	1:D:75:VAL:HG21	2.01	0.42
1:A:173:ASP:OD1	3:A:305:EDO:O1	2.34	0.41
1:E:231:LYS:HD3	1:E:231:LYS:HA	1.72	0.41
1:C:29:ARG:HG2	1:C:298:PHE:CE1	2.56	0.41
1:C:110:TYR:CZ	4:C:309:PEG:H42	2.55	0.41
1:A:124:LYS:HE2	1:A:128:GLN:NE2	2.35	0.41
1:D:142:ARG:NE	5:D:307:ACT:H2	2.32	0.41
1:E:60:ARG:HG3	1:E:95:PHE:CZ	2.56	0.41
1:F:182:GLU:HB3	1:F:185:MET:HE3	2.03	0.41
1:A:10:MET:HG2	1:A:41:ALA:HB3	2.03	0.41
1:C:72:GLY:C	1:E:295:ILE:HD12	2.46	0.41
1:F:160:GLU:H	4:F:305:PEG:C1	2.34	0.41
1:E:130:VAL:O	1:E:161:ASN:ND2	2.53	0.41
1:F:16:PRO:HD2	1:F:27:TYR:HD1	1.85	0.40
1:F:14:ILE:HD13	1:F:260:MET:HG3	2.04	0.40
1:D:16:PRO:HD2	1:D:27:TYR:HD1	1.86	0.40
1:A:56[B]:TRP:CH2	1:A:95:PHE:HB3	2.56	0.40
1:D:254:ILE:HB	1:D:255:PRO:HD3	2.03	0.40
1:A:34:GLN:HE21	4:A:310:PEG:C1	2.34	0.40
1:D:216:ASP:OD1	1:D:216:ASP:N	2.54	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	296/310 (96%)	290 (98%)	6 (2%)	0	100	100
1	B	294/310 (95%)	289 (98%)	5 (2%)	0	100	100
1	C	294/310 (95%)	288 (98%)	6 (2%)	0	100	100
1	D	295/310 (95%)	289 (98%)	6 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	E	294/310 (95%)	288 (98%)	6 (2%)	0	100	100
1	F	297/310 (96%)	292 (98%)	5 (2%)	0	100	100
All	All	1770/1860 (95%)	1736 (98%)	34 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	251/260 (96%)	251 (100%)	0	100	100
1	B	249/260 (96%)	249 (100%)	0	100	100
1	C	249/260 (96%)	249 (100%)	0	100	100
1	D	250/260 (96%)	250 (100%)	0	100	100
1	E	249/260 (96%)	249 (100%)	0	100	100
1	F	252/260 (97%)	252 (100%)	0	100	100
All	All	1500/1560 (96%)	1500 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	19	ASN
1	A	229	ASN
1	B	19	ASN
1	B	128	GLN
1	C	19	ASN
1	D	99	HIS
1	E	37	ASN
1	E	201	ASN
1	F	201	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

6 non-standard protein/DNA/RNA residues are modelled in this entry.

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
1	KPI	F	166	1	11,13,14	1.08	1 (9%)	9,15,17	2.67	4 (44%)
1	KPI	E	166	1	11,13,14	1.00	0	9,15,17	2.52	4 (44%)
1	KPI	B	166	1	11,13,14	1.75	4 (36%)	9,15,17	2.78	4 (44%)
1	KPI	A	166	1	11,13,14	1.25	1 (9%)	9,15,17	2.62	5 (55%)
1	KPI	D	166	1	11,13,14	1.13	1 (9%)	9,15,17	2.33	4 (44%)
1	KPI	C	166	1	11,13,14	1.15	2 (18%)	9,15,17	2.90	5 (55%)

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
1	KPI	F	166	1	-	0/13/14/16	-
1	KPI	E	166	1	-	2/13/14/16	-
1	KPI	B	166	1	-	5/13/14/16	-
1	KPI	A	166	1	-	1/13/14/16	-
1	KPI	D	166	1	-	0/13/14/16	-
1	KPI	C	166	1	-	0/13/14/16	-

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	166	KPI	C1-CX1	3.47	1.56	1.49
1	A	166	KPI	CX2-CX1	3.34	1.53	1.49

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	166	KPI	O2-CX2	2.49	1.28	1.22
1	C	166	KPI	CX2-CX1	2.47	1.52	1.49
1	D	166	KPI	CX2-CX1	2.27	1.52	1.49
1	C	166	KPI	CA-N	-2.22	1.41	1.48
1	B	166	KPI	CX2-CX1	2.19	1.52	1.49
1	B	166	KPI	O1-CX2	-2.11	1.24	1.30
1	F	166	KPI	O2-CX2	2.09	1.27	1.22

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	166	KPI	CE-NZ-CX1	5.76	137.42	121.58
1	F	166	KPI	CE-NZ-CX1	5.46	136.57	121.58
1	D	166	KPI	CE-NZ-CX1	5.17	135.79	121.58
1	A	166	KPI	CE-NZ-CX1	5.09	135.56	121.58
1	C	166	KPI	CE-NZ-CX1	5.03	135.40	121.58
1	B	166	KPI	CE-NZ-CX1	4.87	134.95	121.58
1	C	166	KPI	O2-CX2-CX1	-4.75	115.55	121.35
1	B	166	KPI	C1-CX1-CX2	4.73	122.57	118.11
1	B	166	KPI	O2-CX2-CX1	3.89	126.10	121.35
1	F	166	KPI	O2-CX2-CX1	-3.64	116.90	121.35
1	A	166	KPI	CX2-CX1-NZ	3.63	123.42	114.88
1	C	166	KPI	O1-CX2-CX1	3.31	123.54	116.50
1	F	166	KPI	CX2-CX1-NZ	3.15	122.29	114.88
1	D	166	KPI	CX2-CX1-NZ	2.93	121.78	114.88
1	F	166	KPI	C1-CX1-CX2	-2.76	115.51	118.11
1	A	166	KPI	O2-CX2-CX1	-2.74	118.00	121.35
1	C	166	KPI	CX2-CX1-NZ	2.66	121.13	114.88
1	E	166	KPI	CX2-CX1-NZ	2.53	120.83	114.88
1	C	166	KPI	C1-CX1-CX2	2.48	120.46	118.11
1	A	166	KPI	C1-CX1-CX2	2.44	120.41	118.11
1	E	166	KPI	O2-CX2-CX1	-2.42	118.39	121.35
1	A	166	KPI	O1-CX2-CX1	2.25	121.30	116.50
1	E	166	KPI	O1-CX2-CX1	2.15	121.08	116.50
1	D	166	KPI	O2-CX2-CX1	-2.11	118.77	121.35
1	D	166	KPI	O1-CX2-CX1	2.08	120.94	116.50
1	B	166	KPI	O1-CX2-O2	-2.00	119.14	123.90

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	B	166	KPI	NZ-CX1-CX2-O1
1	B	166	KPI	NZ-CX1-CX2-O2
1	B	166	KPI	C1-CX1-CX2-O1
1	B	166	KPI	C1-CX1-CX2-O2
1	E	166	KPI	C-CA-CB-CG
1	A	166	KPI	NZ-CX1-CX2-O2
1	E	166	KPI	N-CA-CB-CG
1	B	166	KPI	CD-CE-NZ-CX1

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	F	166	KPI	1	0
1	C	166	KPI	2	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 68 ligands modelled in this entry, 13 are monoatomic - leaving 55 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	$\# Z > 2$
5	ACT	E	309	-	3,3,3	0.82	0	3,3,3	1.41	0
5	ACT	D	308	-	3,3,3	0.83	0	3,3,3	1.33	0
7	PGE	B	301	-	9,9,9	0.30	0	8,8,8	0.31	0
4	PEG	D	305	-	6,6,6	0.49	0	5,5,5	0.29	0
3	EDO	C	307	-	3,3,3	0.49	0	2,2,2	0.31	0
7	PGE	B	302	-	9,9,9	0.31	0	8,8,8	0.32	0
4	PEG	F	305	-	6,6,6	0.60	0	5,5,5	0.56	0
4	PEG	A	309	-	6,6,6	0.49	0	5,5,5	0.31	0
3	EDO	E	308	-	3,3,3	0.43	0	2,2,2	0.42	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	EDO	F	303	-	3,3,3	0.51	0	2,2,2	0.25	0
4	PEG	C	310	-	6,6,6	0.49	0	5,5,5	0.19	0
3	EDO	B	306	-	3,3,3	0.47	0	2,2,2	0.36	0
4	PEG	A	310	-	6,6,6	0.70	0	5,5,5	0.51	0
5	ACT	D	307	-	3,3,3	0.80	0	3,3,3	1.26	0
8	GOL	E	311	-	5,5,5	0.36	0	5,5,5	0.37	0
3	EDO	D	303	-	3,3,3	0.36	0	2,2,2	0.47	0
3	EDO	A	308	-	3,3,3	0.41	0	2,2,2	0.39	0
5	ACT	C	312	-	3,3,3	0.80	0	3,3,3	1.27	0
5	ACT	F	307	-	3,3,3	0.81	0	3,3,3	1.33	0
3	EDO	B	310	-	3,3,3	0.56	0	2,2,2	0.23	0
3	EDO	A	303	-	3,3,3	0.40	0	2,2,2	0.51	0
4	PEG	C	309	2	6,6,6	0.53	0	5,5,5	0.27	0
3	EDO	C	306	-	3,3,3	0.50	0	2,2,2	0.26	0
3	EDO	C	304	-	3,3,3	0.31	0	2,2,2	0.50	0
3	EDO	A	307	-	3,3,3	0.35	0	2,2,2	0.70	0
5	ACT	A	311	-	3,3,3	0.84	0	3,3,3	1.30	0
5	ACT	F	308	-	3,3,3	0.77	0	3,3,3	1.37	0
6	PG4	D	309	-	12,12,12	0.53	0	11,11,11	0.27	0
8	GOL	B	313	-	5,5,5	0.29	0	5,5,5	0.38	0
8	GOL	E	310	-	5,5,5	0.33	0	5,5,5	0.33	0
3	EDO	B	307	-	3,3,3	0.45	0	2,2,2	0.33	0
3	EDO	F	304	-	3,3,3	0.35	0	2,2,2	0.47	0
4	PEG	D	306	-	6,6,6	0.49	0	5,5,5	0.22	0
3	EDO	E	304	-	3,3,3	0.44	0	2,2,2	0.45	0
3	EDO	B	311	-	3,3,3	0.44	0	2,2,2	0.41	0
5	ACT	C	313	-	3,3,3	0.86	0	3,3,3	1.51	0
4	PEG	F	306	-	6,6,6	0.50	0	5,5,5	0.65	0
3	EDO	E	306	-	3,3,3	0.39	0	2,2,2	0.45	0
6	PG4	A	312	-	12,12,12	0.52	0	11,11,11	0.19	0
5	ACT	F	309	-	3,3,3	0.80	0	3,3,3	1.34	0
3	EDO	B	308	-	3,3,3	0.42	0	2,2,2	0.46	0
3	EDO	B	312	-	3,3,3	0.42	0	2,2,2	0.42	0
3	EDO	E	305	-	3,3,3	0.47	0	2,2,2	0.15	0
8	GOL	C	314	-	5,5,5	0.32	0	5,5,5	0.36	0
3	EDO	A	306	-	3,3,3	0.54	0	2,2,2	0.19	0
3	EDO	C	305	-	3,3,3	0.47	0	2,2,2	0.32	0
4	PEG	C	311	-	6,6,6	0.50	0	5,5,5	0.42	0
7	PGE	E	301	-	9,9,9	0.34	0	8,8,8	0.19	0
3	EDO	D	304	-	3,3,3	0.44	0	2,2,2	0.36	0
3	EDO	B	309	-	3,3,3	0.45	0	2,2,2	0.30	0
3	EDO	C	308	-	3,3,3	0.46	0	2,2,2	0.41	0
3	EDO	A	305	-	3,3,3	0.37	0	2,2,2	0.44	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
7	PGE	C	301	-	9,9,9	0.32	0	8,8,8	0.29	0
3	EDO	E	307	-	3,3,3	0.48	0	2,2,2	0.35	0
3	EDO	A	304	-	3,3,3	0.38	0	2,2,2	0.55	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
7	PGE	B	301	-	-	1/7/7/7	-
4	PEG	D	305	-	-	0/4/4/4	-
3	EDO	C	307	-	-	1/1/1/1	-
7	PGE	B	302	-	-	6/7/7/7	-
4	PEG	F	305	-	-	2/4/4/4	-
4	PEG	A	309	-	-	1/4/4/4	-
3	EDO	E	308	-	-	1/1/1/1	-
3	EDO	F	303	-	-	0/1/1/1	-
4	PEG	C	310	-	-	2/4/4/4	-
3	EDO	B	306	-	-	0/1/1/1	-
4	PEG	A	310	-	-	0/4/4/4	-
8	GOL	E	311	-	-	2/4/4/4	-
3	EDO	D	303	-	-	0/1/1/1	-
3	EDO	A	308	-	-	1/1/1/1	-
3	EDO	B	310	-	-	0/1/1/1	-
3	EDO	A	303	-	-	0/1/1/1	-
4	PEG	C	309	2	-	2/4/4/4	-
3	EDO	C	306	-	-	1/1/1/1	-
3	EDO	C	304	-	-	1/1/1/1	-
3	EDO	A	307	-	-	1/1/1/1	-
6	PG4	D	309	-	-	2/10/10/10	-
8	GOL	B	313	-	-	2/4/4/4	-
8	GOL	E	310	-	-	2/4/4/4	-
3	EDO	B	307	-	-	1/1/1/1	-
3	EDO	F	304	-	-	0/1/1/1	-
4	PEG	D	306	-	-	1/4/4/4	-
3	EDO	E	304	-	-	1/1/1/1	-
3	EDO	B	311	-	-	0/1/1/1	-
4	PEG	F	306	-	-	3/4/4/4	-
3	EDO	E	306	-	-	1/1/1/1	-
6	PG4	A	312	-	-	0/10/10/10	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	EDO	B	308	-	-	1/1/1/1	-
3	EDO	B	312	-	-	1/1/1/1	-
3	EDO	E	305	-	-	0/1/1/1	-
8	GOL	C	314	-	-	4/4/4/4	-
3	EDO	A	306	-	-	1/1/1/1	-
3	EDO	C	305	-	-	0/1/1/1	-
4	PEG	C	311	-	-	4/4/4/4	-
7	PGE	E	301	-	-	0/7/7/7	-
3	EDO	D	304	-	-	1/1/1/1	-
3	EDO	B	309	-	-	0/1/1/1	-
3	EDO	C	308	-	-	1/1/1/1	-
3	EDO	A	305	-	-	0/1/1/1	-
7	PGE	C	301	-	-	1/7/7/7	-
3	EDO	E	307	-	-	1/1/1/1	-
3	EDO	A	304	-	-	0/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (50) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
8	B	313	GOL	O1-C1-C2-C3
8	E	311	GOL	O1-C1-C2-C3
4	F	306	PEG	C4-C3-O2-C2
4	C	309	PEG	O2-C3-C4-O4
4	C	311	PEG	O2-C3-C4-O4
4	F	306	PEG	O2-C3-C4-O4
8	C	314	GOL	O1-C1-C2-C3
8	C	314	GOL	C1-C2-C3-O3
8	E	310	GOL	O1-C1-C2-C3
8	B	313	GOL	O1-C1-C2-O2
4	C	311	PEG	O1-C1-C2-O2
4	F	306	PEG	O1-C1-C2-O2
7	B	302	PGE	O3-C5-C6-O4
3	A	306	EDO	O1-C1-C2-O2
4	F	305	PEG	C4-C3-O2-C2
4	C	310	PEG	O2-C3-C4-O4
4	C	309	PEG	C1-C2-O2-C3
7	B	302	PGE	O2-C3-C4-O3

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Mol	Chain	Res	Type	Atoms
8	E	311	GOL	O1-C1-C2-O2
4	A	309	PEG	O2-C3-C4-O4
3	E	308	EDO	O1-C1-C2-O2
8	C	314	GOL	O1-C1-C2-O2
8	C	314	GOL	O2-C2-C3-O3
8	E	310	GOL	O1-C1-C2-O2
7	B	302	PGE	C4-C3-O2-C2
7	B	302	PGE	C6-C5-O3-C4
6	D	309	PG4	C3-C4-O3-C5
7	B	302	PGE	C1-C2-O2-C3
4	C	311	PEG	C1-C2-O2-C3
7	C	301	PGE	C1-C2-O2-C3
4	F	305	PEG	O1-C1-C2-O2
3	B	312	EDO	O1-C1-C2-O2
3	C	304	EDO	O1-C1-C2-O2
3	D	304	EDO	O1-C1-C2-O2
7	B	301	PGE	O1-C1-C2-O2
3	A	308	EDO	O1-C1-C2-O2
3	B	307	EDO	O1-C1-C2-O2
3	C	307	EDO	O1-C1-C2-O2
3	C	308	EDO	O1-C1-C2-O2
3	E	306	EDO	O1-C1-C2-O2
4	D	306	PEG	O1-C1-C2-O2
4	C	311	PEG	C4-C3-O2-C2
3	C	306	EDO	O1-C1-C2-O2
3	E	304	EDO	O1-C1-C2-O2
3	E	307	EDO	O1-C1-C2-O2
6	D	309	PG4	C1-C2-O2-C3
7	B	302	PGE	O1-C1-C2-O2
3	A	307	EDO	O1-C1-C2-O2
4	C	310	PEG	C4-C3-O2-C2
3	B	308	EDO	O1-C1-C2-O2

There are no ring outliers.

23 monomers are involved in 45 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	B	302	PGE	2	0
4	F	305	PEG	4	0
4	A	309	PEG	1	0
4	A	310	PEG	8	0
5	D	307	ACT	2	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	308	EDO	2	0
3	A	303	EDO	1	0
4	C	309	PEG	1	0
3	C	304	EDO	4	0
3	A	307	EDO	1	0
5	A	311	ACT	1	0
5	F	308	ACT	1	0
6	D	309	PG4	1	0
8	E	310	GOL	1	0
3	B	307	EDO	3	0
3	F	304	EDO	1	0
3	E	304	EDO	2	0
3	B	311	EDO	1	0
4	F	306	PEG	2	0
3	E	306	EDO	3	0
3	E	305	EDO	1	0
3	A	305	EDO	3	0
3	E	307	EDO	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	296/310 (95%)	-0.29	1 (0%) 90 92	21, 27, 41, 59	2 (0%)
1	B	295/310 (95%)	-0.25	1 (0%) 90 92	21, 27, 41, 54	1 (0%)
1	C	295/310 (95%)	-0.21	3 (1%) 79 84	21, 28, 41, 53	1 (0%)
1	D	295/310 (95%)	-0.16	3 (1%) 79 84	21, 28, 44, 56	3 (1%)
1	E	295/310 (95%)	-0.18	1 (0%) 90 92	20, 29, 42, 54	1 (0%)
1	F	296/310 (95%)	-0.10	2 (0%) 84 88	19, 29, 44, 58	4 (1%)
All	All	1772/1860 (95%)	-0.20	11 (0%) 85 89	19, 28, 43, 59	12 (0%)

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	56[A]	TRP	5.2
1	A	56[A]	TRP	4.3
1	D	56	TRP	4.3
1	C	56[A]	TRP	4.2
1	F	56	TRP	4.0
1	F	95	PHE	3.2
1	D	298	PHE	2.7
1	E	56	TRP	2.6
1	C	95	PHE	2.1
1	D	95	PHE	2.1
1	C	160	GLU	2.0

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

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
1	KPI	A	166	14/15	0.75	0.16	20,20,20,20	0
1	KPI	D	166	14/15	0.75	0.17	20,20,20,20	0
1	KPI	B	166	14/15	0.76	0.15	20,20,20,20	0
1	KPI	E	166	14/15	0.76	0.15	20,20,20,20	0
1	KPI	C	166	14/15	0.78	0.15	20,20,20,20	0
1	KPI	F	166	14/15	0.78	0.15	20,20,20,20	0

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
8	GOL	E	311	6/6	0.67	0.17	51,54,62,63	0
5	ACT	C	313	4/4	0.68	0.23	37,48,49,50	0
4	PEG	C	311	7/7	0.70	0.18	42,47,52,61	0
8	GOL	B	313	6/6	0.73	0.16	42,52,60,76	0
8	GOL	C	314	6/6	0.75	0.16	40,49,50,53	0
5	ACT	D	307	4/4	0.76	0.17	39,41,48,58	0
8	GOL	E	310	6/6	0.76	0.14	48,50,58,58	0
4	PEG	C	309	7/7	0.76	0.15	38,42,56,57	0
4	PEG	C	310	7/7	0.77	0.17	35,44,56,62	0
4	PEG	F	306	7/7	0.77	0.16	32,40,45,45	0
5	ACT	F	307	4/4	0.77	0.15	47,54,55,65	0
5	ACT	F	309	4/4	0.77	0.14	40,47,50,53	0
7	PGE	B	302	10/10	0.80	0.15	39,44,52,62	0
3	EDO	C	308	4/4	0.80	0.16	46,52,54,57	0
3	EDO	E	304	4/4	0.80	0.14	33,36,43,47	0
2	MG	C	302	1/1	0.80	0.16	56,56,56,56	0
5	ACT	A	311	4/4	0.80	0.16	42,44,47,53	0
3	EDO	B	312	4/4	0.82	0.15	41,42,44,56	0
3	EDO	E	308	4/4	0.82	0.16	36,37,38,44	0
3	EDO	A	307	4/4	0.84	0.21	29,38,38,40	0
3	EDO	A	306	4/4	0.85	0.13	34,39,41,47	0
3	EDO	A	308	4/4	0.85	0.17	36,39,46,56	0
3	EDO	C	305	4/4	0.86	0.15	40,41,47,61	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	EDO	B	308	4/4	0.86	0.12	38,44,45,49	0
2	MG	F	302	1/1	0.86	0.12	55,55,55,55	0
3	EDO	A	304	4/4	0.87	0.11	37,38,41,42	0
3	EDO	B	306	4/4	0.88	0.14	35,36,37,56	0
4	PEG	D	306	7/7	0.88	0.11	42,49,53,55	0
5	ACT	D	308	4/4	0.88	0.11	31,41,43,51	0
3	EDO	E	305	4/4	0.88	0.14	31,31,44,50	0
3	EDO	E	306	4/4	0.88	0.14	33,44,48,48	0
4	PEG	A	309	7/7	0.89	0.10	38,45,55,56	0
5	ACT	C	312	4/4	0.89	0.12	39,46,47,48	0
3	EDO	C	307	4/4	0.89	0.10	43,44,44,53	0
3	EDO	A	303	4/4	0.89	0.16	34,36,37,38	0
6	PG4	D	309	13/13	0.90	0.10	36,43,59,62	0
7	PGE	B	301	10/10	0.90	0.09	32,43,47,47	0
3	EDO	B	310	4/4	0.90	0.12	28,33,36,41	0
7	PGE	C	301	10/10	0.90	0.10	34,39,47,50	0
3	EDO	F	303	4/4	0.90	0.13	39,40,44,47	0
3	EDO	F	304	4/4	0.90	0.15	27,38,41,48	0
5	ACT	F	308	4/4	0.90	0.12	33,42,44,52	0
3	EDO	C	304	4/4	0.90	0.15	35,36,37,38	0
7	PGE	E	301	10/10	0.91	0.10	41,45,49,50	0
3	EDO	B	311	4/4	0.91	0.10	36,40,42,42	0
4	PEG	D	305	7/7	0.91	0.11	42,45,49,52	0
3	EDO	B	307	4/4	0.91	0.13	32,37,43,49	0
3	EDO	D	304	4/4	0.91	0.13	38,39,41,47	0
2	MG	C	303	1/1	0.92	0.09	51,51,51,51	0
4	PEG	F	305	7/7	0.92	0.11	25,29,31,33	0
3	EDO	C	306	4/4	0.92	0.11	30,31,37,42	0
3	EDO	A	305	4/4	0.92	0.16	27,33,41,43	0
2	MG	E	303	1/1	0.92	0.09	44,44,44,44	0
2	MG	D	301	1/1	0.93	0.14	58,58,58,58	0
3	EDO	D	303	4/4	0.93	0.10	29,36,50,54	0
4	PEG	A	310	7/7	0.93	0.18	23,30,34,34	0
2	MG	F	301	1/1	0.94	0.07	36,36,36,36	0
3	EDO	E	307	4/4	0.94	0.09	31,33,34,36	0
6	PG4	A	312	13/13	0.94	0.08	30,35,46,47	0
2	MG	B	303	1/1	0.95	0.05	31,31,31,31	0
2	MG	A	302	1/1	0.95	0.07	46,46,46,46	0
2	MG	E	302	1/1	0.95	0.10	42,42,42,42	0
3	EDO	B	309	4/4	0.95	0.08	33,35,38,43	0
2	MG	B	304	1/1	0.96	0.14	48,48,48,48	0
2	MG	B	305	1/1	0.97	0.06	39,39,39,39	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	MG	A	301	1/1	0.97	0.03	30,30,30,30	0
5	ACT	E	309	4/4	0.98	0.06	30,33,34,35	0
2	MG	D	302	1/1	0.98	0.04	36,36,36,36	0

6.5 Other polymers [i](#)

There are no such residues in this entry.