



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 15, 2017 – 03:54 am GMT

PDB ID : 4MLR
Title : dihydrodipicolinate synthase from *C. jejuni*, Y110F mutation with pyruvate and Lysine
Authors : Conly, C.J.T.
Deposited on : 2013-09-06
Resolution : 2.20 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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : recalc28949

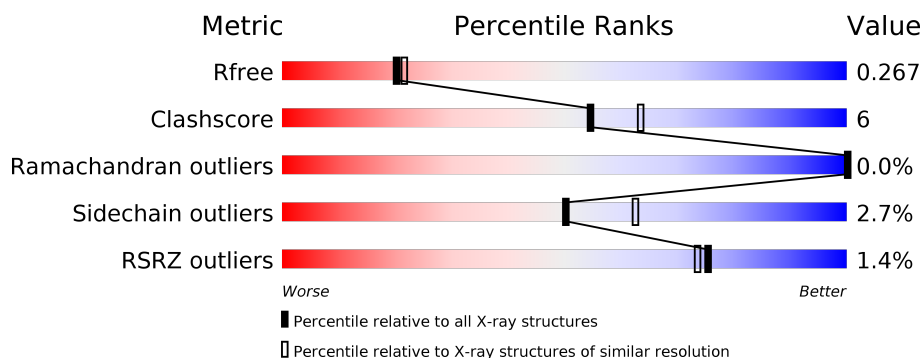
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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.



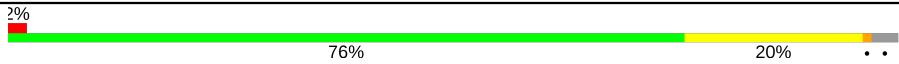

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	100719	4002 (2.20-2.20)
Clashscore	112137	4730 (2.20-2.20)
Ramachandran outliers	110173	4656 (2.20-2.20)
Sidechain outliers	110143	4657 (2.20-2.20)
RSRZ outliers	101464	4033 (2.20-2.20)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	306	<div> <div>84%</div> <div>12%</div> <div>••</div> </div>
1	B	306	<div> <div>%</div> <div>82%</div> <div>14%</div> <div>••</div> </div>
1	C	306	<div> <div>82%</div> <div>14%</div> <div>•</div> </div>
1	D	306	<div> <div>%</div> <div>78%</div> <div>17%</div> <div>••</div> </div>
1	E	306	<div> <div>%</div> <div>83%</div> <div>13%</div> <div>•</div> </div>
1	F	306	<div> <div>%</div> <div>81%</div> <div>14%</div> <div>••</div> </div>

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Mol	Chain	Length	Quality of chain
1	G	306	
1	H	306	

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	LYS	B	301	-	-	-	X
2	LYS	C	301	-	-	-	X
2	LYS	G	301	-	-	-	X
3	EDO	D	301	-	-	-	X
3	EDO	G	302	-	-	-	X
5	PG4	C	302	-	-	-	X
6	PGE	B	303	-	-	-	X

2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 18601 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 dihydrodipicolinate synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	296	Total	C	N	O	S	0	1	0
			2282	1454	379	436	13			
1	B	296	Total	C	N	O	S	1	0	0
			2278	1450	379	436	13			
1	C	295	Total	C	N	O	S	0	0	0
			2269	1444	377	435	13			
1	D	295	Total	C	N	O	S	0	1	0
			2278	1447	382	436	13			
1	E	294	Total	C	N	O	S	0	0	0
			2261	1440	375	433	13			
1	F	296	Total	C	N	O	S	1	0	0
			2278	1450	379	436	13			
1	G	296	Total	C	N	O	S	0	0	0
			2278	1450	379	436	13			
1	H	295	Total	C	N	O	S	0	0	0
			2269	1444	377	435	13			

There are 72 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
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	ALA	-	EXPRESSION TAG	UNP Q9PPB4
A	0	SER	-	EXPRESSION TAG	UNP Q9PPB4
A	110	PHE	TYR	ENGINEERED MUTATION	UNP Q9PPB4
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

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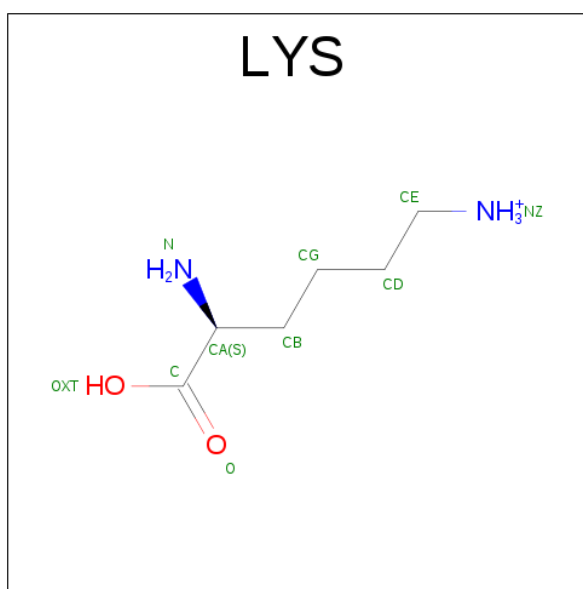
Chain	Residue	Modelled	Actual	Comment	Reference
B	-3	HIS	-	EXPRESSION TAG	UNP Q9PPB4
B	-2	HIS	-	EXPRESSION TAG	UNP Q9PPB4
B	-1	ALA	-	EXPRESSION TAG	UNP Q9PPB4
B	0	SER	-	EXPRESSION TAG	UNP Q9PPB4
B	110	PHE	TYR	ENGINEERED MUTATION	UNP Q9PPB4
C	-7	HIS	-	EXPRESSION TAG	UNP Q9PPB4
C	-6	HIS	-	EXPRESSION TAG	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	ALA	-	EXPRESSION TAG	UNP Q9PPB4
C	0	SER	-	EXPRESSION TAG	UNP Q9PPB4
C	110	PHE	TYR	ENGINEERED MUTATION	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	ALA	-	EXPRESSION TAG	UNP Q9PPB4
D	0	SER	-	EXPRESSION TAG	UNP Q9PPB4
D	110	PHE	TYR	ENGINEERED MUTATION	UNP Q9PPB4
E	-7	HIS	-	EXPRESSION TAG	UNP Q9PPB4
E	-6	HIS	-	EXPRESSION TAG	UNP Q9PPB4
E	-5	HIS	-	EXPRESSION TAG	UNP Q9PPB4
E	-4	HIS	-	EXPRESSION TAG	UNP Q9PPB4
E	-3	HIS	-	EXPRESSION TAG	UNP Q9PPB4
E	-2	HIS	-	EXPRESSION TAG	UNP Q9PPB4
E	-1	ALA	-	EXPRESSION TAG	UNP Q9PPB4
E	0	SER	-	EXPRESSION TAG	UNP Q9PPB4
E	110	PHE	TYR	ENGINEERED MUTATION	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	ALA	-	EXPRESSION TAG	UNP Q9PPB4
F	0	SER	-	EXPRESSION TAG	UNP Q9PPB4
F	110	PHE	TYR	ENGINEERED MUTATION	UNP Q9PPB4
G	-7	HIS	-	EXPRESSION TAG	UNP Q9PPB4

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Chain	Residue	Modelled	Actual	Comment	Reference
G	-6	HIS	-	EXPRESSION TAG	UNP Q9PPB4
G	-5	HIS	-	EXPRESSION TAG	UNP Q9PPB4
G	-4	HIS	-	EXPRESSION TAG	UNP Q9PPB4
G	-3	HIS	-	EXPRESSION TAG	UNP Q9PPB4
G	-2	HIS	-	EXPRESSION TAG	UNP Q9PPB4
G	-1	ALA	-	EXPRESSION TAG	UNP Q9PPB4
G	0	SER	-	EXPRESSION TAG	UNP Q9PPB4
G	110	PHE	TYR	ENGINEERED MUTATION	UNP Q9PPB4
H	-7	HIS	-	EXPRESSION TAG	UNP Q9PPB4
H	-6	HIS	-	EXPRESSION TAG	UNP Q9PPB4
H	-5	HIS	-	EXPRESSION TAG	UNP Q9PPB4
H	-4	HIS	-	EXPRESSION TAG	UNP Q9PPB4
H	-3	HIS	-	EXPRESSION TAG	UNP Q9PPB4
H	-2	HIS	-	EXPRESSION TAG	UNP Q9PPB4
H	-1	ALA	-	EXPRESSION TAG	UNP Q9PPB4
H	0	SER	-	EXPRESSION TAG	UNP Q9PPB4
H	110	PHE	TYR	ENGINEERED MUTATION	UNP Q9PPB4

- Molecule 2 is LYSINE (three-letter code: LYS) (formula: $C_6H_{15}N_2O_2$).



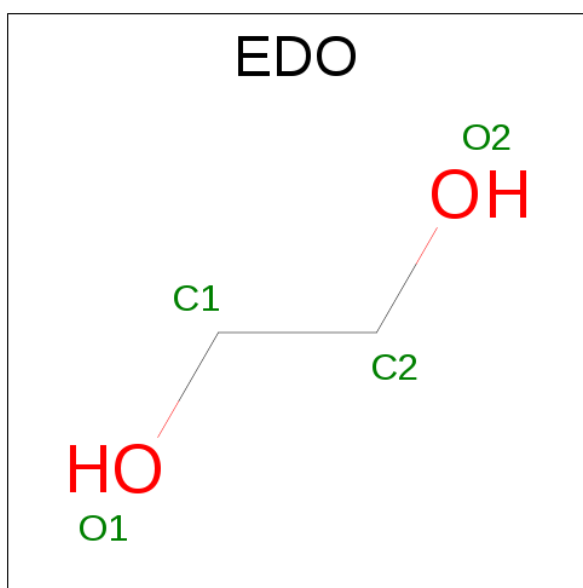
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			10	6	2	2		
2	A	1	Total	C	N	O	0	0
			10	6	2	2		
2	B	1	Total	C	N	O	0	0
			10	6	2	2		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	C	1	Total	C	N	O	0	0
			10	6	2	2		
2	E	1	Total	C	N	O	0	0
			10	6	2	2		
2	F	1	Total	C	N	O	0	0
			10	6	2	2		
2	G	1	Total	C	N	O	0	0
			10	6	2	2		
2	H	1	Total	C	N	O	0	0
			10	6	2	2		

- Molecule 3 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



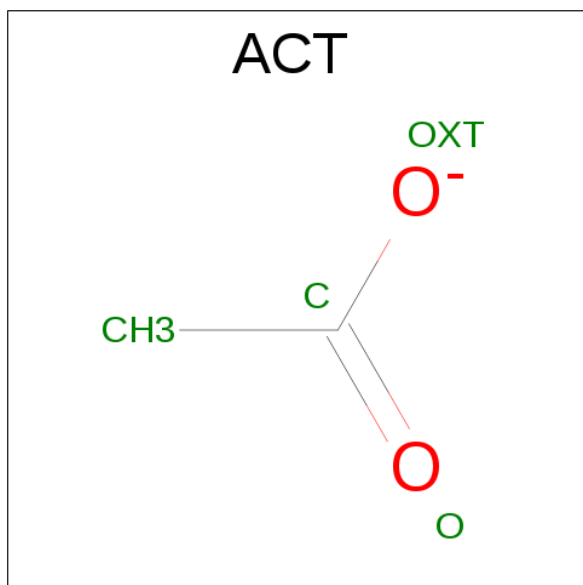
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	D	1	Total	C	O	0	0
			4	2	2		
3	E	1	Total	C	O	0	0
			4	2	2		
3	E	1	Total	C	O	0	0
			4	2	2		

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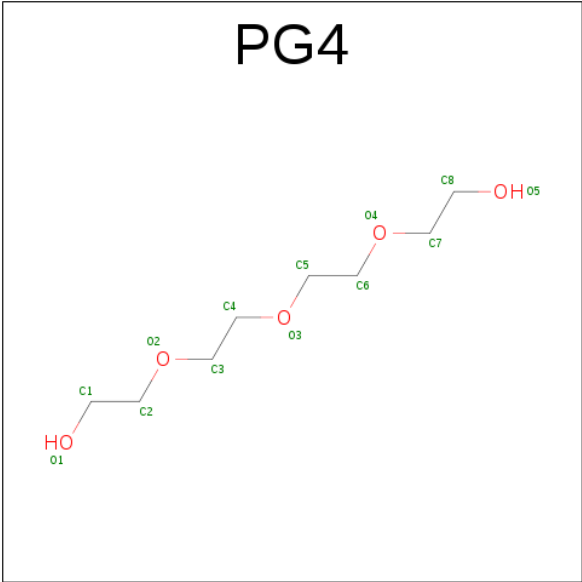
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	G	1	Total	C	O	0	0
			4	2	2		
3	G	1	Total	C	O	0	0
			4	2	2		
3	G	1	Total	C	O	0	0
			4	2	2		

- Molecule 4 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).



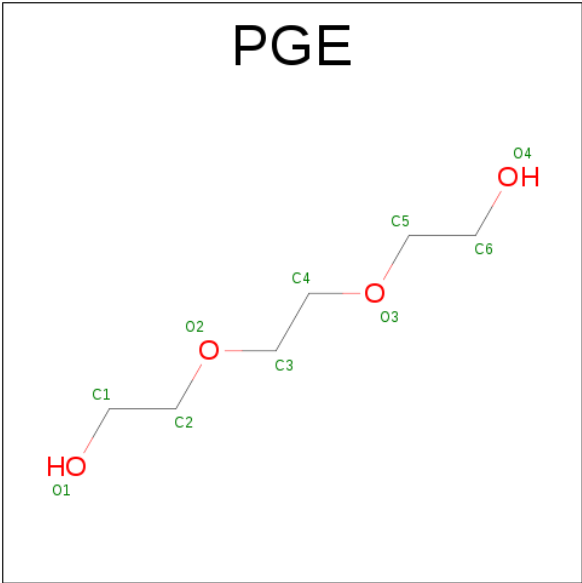
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			4	2	2		
4	F	1	Total	C	O	0	0
			4	2	2		

- Molecule 5 is TETRAETHYLENE GLYCOL (three-letter code: PG4) (formula: $C_8H_{18}O_5$).



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

- Molecule 6 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula: C₆H₁₄O₄).



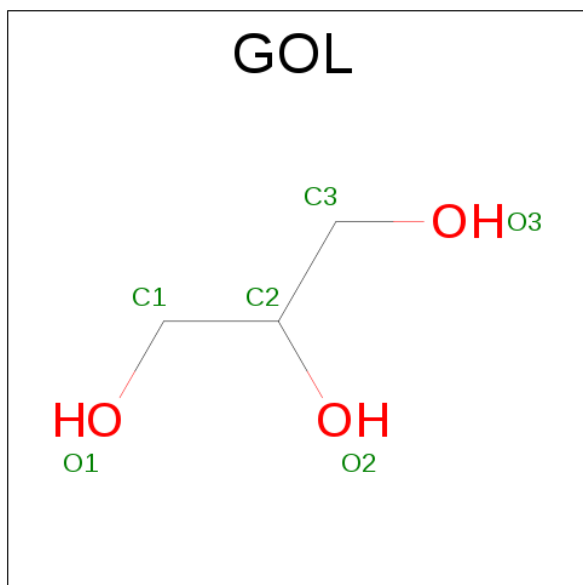
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	B	1	Total	C	O	0	0
			10	6	4		
6	C	1	Total	C	O	0	0
			10	6	4		

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

- Molecule 7 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



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

- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	42	Total	O	0	0
			42	42		
8	B	29	Total	O	0	0
			29	29		
8	C	23	Total	O	0	0
			23	23		
8	D	23	Total	O	0	0
			23	23		
8	E	32	Total	O	0	0
			32	32		

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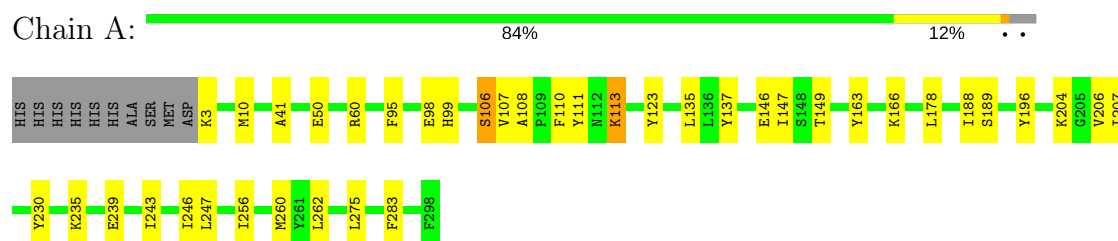
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	F	23	Total 23	O 23	0	0
8	G	17	Total 17	O 17	0	0
8	H	13	Total 13	O 13	0	0

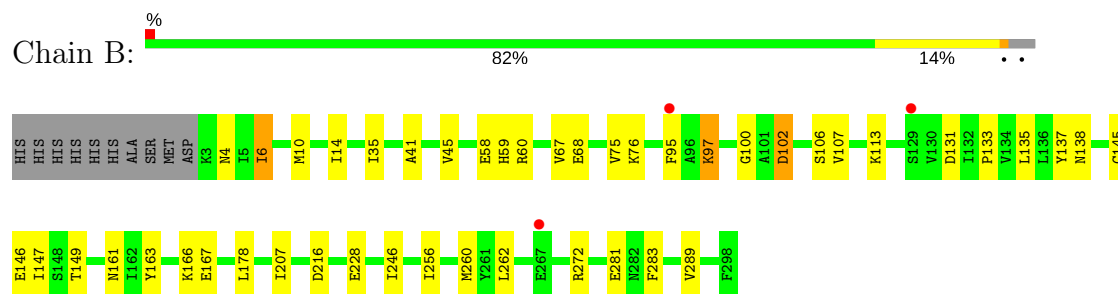
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

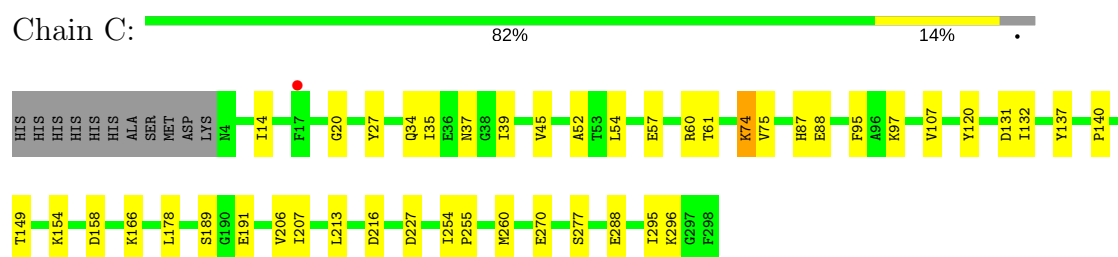
• Molecule 1: dihydrodipicolinate synthase



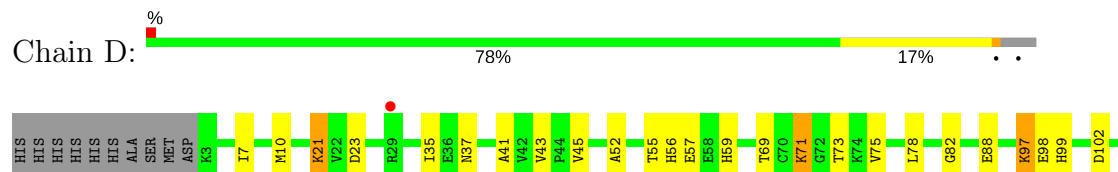
• Molecule 1: dihydrodipicolinate synthase

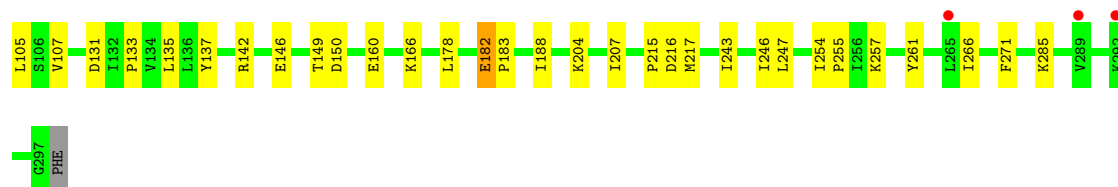


• Molecule 1: dihydrodipicolinate synthase

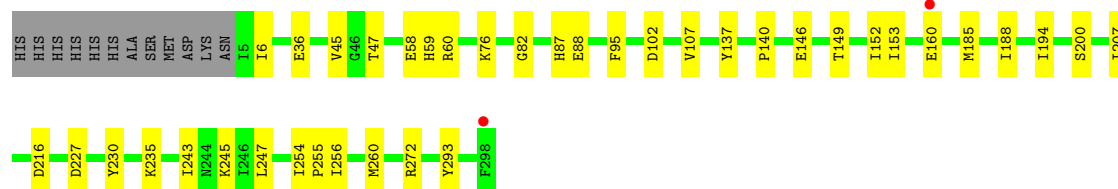
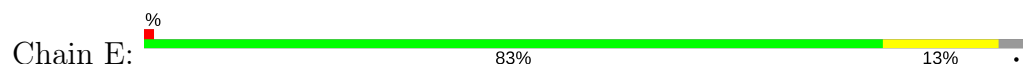


• Molecule 1: dihydrodipicolinate synthase

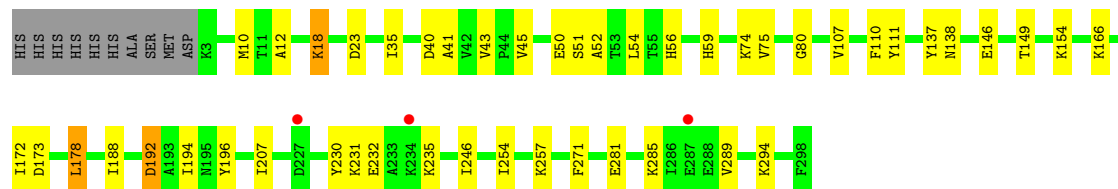
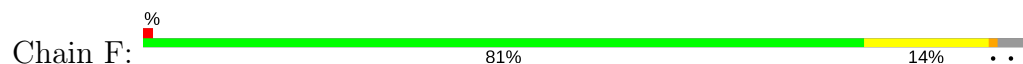




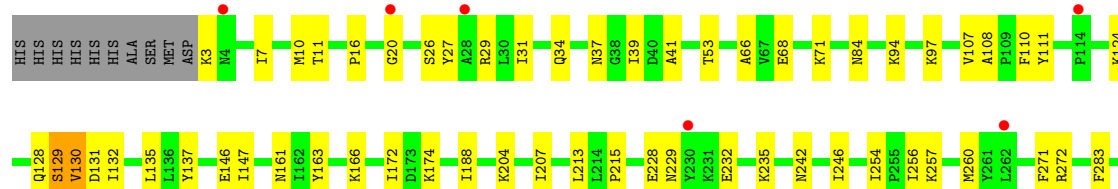
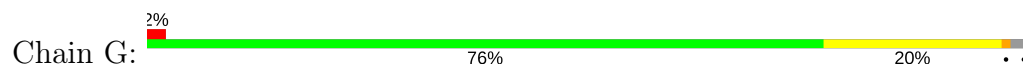
- Molecule 1: dihydrodipicolinate synthase



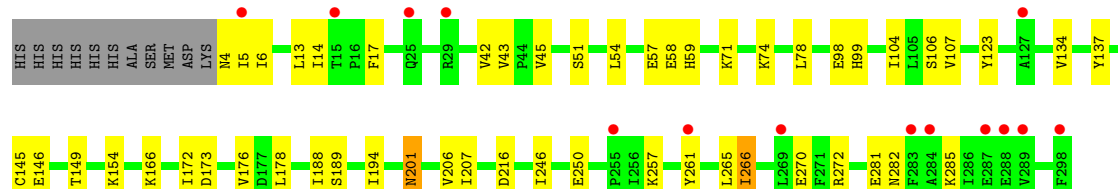
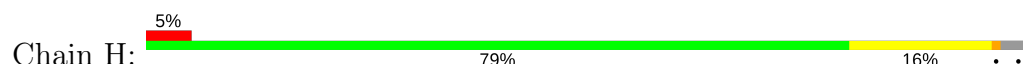
- Molecule 1: dihydrodipicolinate synthase



- Molecule 1: dihydrodipicolinate synthase



- Molecule 1: dihydrodipicolinate synthase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	90.98Å 97.61Å 131.40Å 90.00° 92.06° 90.00°	Depositor
Resolution (Å)	43.45 – 2.20 43.45 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.8 (43.45-2.20) 99.9 (43.45-2.20)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.39 (at 2.20Å)	Xtriage
Refinement program	PHENIX (phenix.refine: dev_1356)	Depositor
R, R_{free}	0.214 , 0.269 0.214 , 0.267	Depositor DCC
R_{free} test set	5833 reflections (5.01%)	DCC
Wilson B-factor (Å ²)	44.0	Xtriage
Anisotropy	0.070	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 48.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.000 for h,-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	18601	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.37% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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	0.38	0/2309	0.47	0/3121
1	B	0.41	0/2302	0.45	0/3111
1	C	0.38	0/2293	0.46	0/3100
1	D	0.51	0/2301	0.50	0/3109
1	E	0.38	0/2285	0.45	0/3089
1	F	0.37	0/2302	0.45	0/3111
1	G	0.42	0/2302	0.45	0/3111
1	H	0.36	0/2293	0.45	0/3100
All	All	0.40	0/18387	0.46	0/24852

There are no bond length outliers.

There are no bond angle outliers.

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	2282	0	2328	24	0
1	B	2278	0	2319	25	0
1	C	2269	0	2306	24	0
1	D	2278	0	2322	35	0
1	E	2261	0	2300	24	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	F	2278	0	2319	29	0
1	G	2278	0	2318	38	0
1	H	2269	0	2306	29	0
2	A	20	0	24	1	0
2	B	10	0	12	0	0
2	C	10	0	12	1	0
2	E	10	0	12	0	0
2	F	10	0	12	2	0
2	G	10	0	12	0	0
2	H	10	0	12	2	0
3	A	12	0	18	1	0
3	D	4	0	6	0	0
3	E	8	0	12	0	0
3	G	12	0	18	0	0
4	A	4	0	3	0	0
4	F	4	0	3	1	0
5	B	13	0	18	0	0
5	C	13	0	18	1	0
6	B	10	0	14	0	0
6	C	10	0	14	0	0
6	E	10	0	14	0	0
6	F	10	0	14	1	0
6	H	10	0	14	1	0
7	C	6	0	8	1	0
8	A	42	0	0	0	0
8	B	29	0	0	0	0
8	C	23	0	0	0	0
8	D	23	0	0	0	0
8	E	32	0	0	0	0
8	F	23	0	0	0	0
8	G	17	0	0	0	0
8	H	13	0	0	0	0
All	All	18601	0	18788	222	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 6.

All (222) 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:G:246:ILE:HG13	1:G:289:VAL:HG21	1.52	0.89

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:107:VAL:HA	1:B:137:TYR:HB3	1.60	0.83
1:A:107:VAL:HA	1:A:137:TYR:HB3	1.64	0.79
1:D:254:ILE:HB	1:D:255:PRO:HD3	1.65	0.78
1:G:107:VAL:HA	1:G:137:TYR:HB3	1.68	0.75
1:C:107:VAL:HA	1:C:137:TYR:HB3	1.69	0.74
1:A:113:LYS:HE3	1:D:142[A]:ARG:HD3	1.74	0.70
1:B:4:ASN:HB3	1:B:133:PRO:HG3	1.72	0.69
1:G:229:ASN:OD1	1:G:232:GLU:HG2	1.93	0.69
1:D:69:THR:O	1:D:69:THR:HG22	1.92	0.68
1:D:23:ASP:OD1	1:D:23:ASP:C	2.30	0.68
1:D:82:GLY:HA3	1:D:107:VAL:HG12	1.76	0.67
1:C:140:PRO:HB2	7:C:304:GOL:H2	1.76	0.67
1:F:56:HIS:H	4:F:303:ACT:H2	1.59	0.67
1:D:107:VAL:HA	1:D:137:TYR:HB3	1.79	0.65
1:G:166:KPI:HDA	1:G:207:ILE:HD12	1.78	0.64
1:A:10:MET:HG2	1:A:41:ALA:HB3	1.78	0.64
1:D:149:THR:HG23	1:D:178:LEU:HD23	1.80	0.64
1:E:107:VAL:HA	1:E:137:TYR:HB3	1.80	0.64
1:H:107:VAL:HA	1:H:137:TYR:HB3	1.80	0.64
1:F:40:ASP:OD1	1:F:74:LYS:NZ	2.31	0.63
1:G:10:MET:HG2	1:G:41:ALA:HB3	1.81	0.63
1:E:58:GLU:OE2	1:E:272:ARG:NH2	2.27	0.62
1:F:166:KPI:HDA	1:F:207:ILE:HD12	1.81	0.62
1:G:97:LYS:HE2	1:G:131:ASP:OD1	2.00	0.61
1:F:107:VAL:HA	1:F:137:TYR:HB3	1.82	0.61
1:D:166:KPI:HDA	1:D:207:ILE:HD12	1.82	0.60
1:A:235:LYS:O	1:A:239:GLU:HG2	2.02	0.60
1:B:246:ILE:HG21	1:B:289:VAL:HG21	1.83	0.60
1:H:149:THR:HG23	1:H:178:LEU:HD23	1.83	0.60
1:G:128:GLN:O	1:G:129:SER:HB3	2.02	0.60
1:G:94:LYS:HE3	1:G:129:SER:HB2	1.82	0.60
1:C:60:ARG:HB3	1:C:95:PHE:CZ	2.37	0.60
1:H:246:ILE:HD11	1:H:285:LYS:HB3	1.84	0.59
1:D:10:MET:HG2	1:D:41:ALA:HB3	1.84	0.59
1:B:113:LYS:HE2	1:B:145:CYS:HA	1.85	0.59
1:G:97:LYS:CE	1:G:131:ASP:OD1	2.50	0.59
1:G:147:ILE:O	1:G:174:LYS:NZ	2.35	0.58
1:G:110:PHE:O	1:G:111:TYR:HB3	2.03	0.58
1:G:16:PRO:HD2	1:G:27:TYR:HD1	1.69	0.58
1:D:160:GLU:CD	1:D:160:GLU:H	2.06	0.58
1:H:58:GLU:OE2	1:H:272:ARG:NH2	2.32	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:302:LYS:N	1:D:52:ALA:O	2.38	0.57
1:H:250:GLU:OE2	1:H:282:ASN:ND2	2.38	0.57
1:A:113:LYS:CE	1:D:142[A]:ARG:HD3	2.35	0.57
1:H:51:SER:O	2:H:301:LYS:NZ	2.38	0.56
1:D:69:THR:O	1:D:69:THR:CG2	2.54	0.56
1:E:216:ASP:OD1	1:E:216:ASP:N	2.37	0.56
1:F:10:MET:HG2	1:F:41:ALA:HB3	1.88	0.56
1:G:188:ILE:HG21	1:G:207:ILE:HG13	1.88	0.55
1:D:35:ILE:HG12	1:D:75:VAL:HG21	1.87	0.55
1:H:201:ASN:ND2	1:H:201:ASN:O	2.38	0.55
1:A:188:ILE:HG21	1:A:207:ILE:HG13	1.89	0.54
1:F:50:GLU:OE2	1:F:257:LYS:NZ	2.32	0.54
1:H:43:VAL:HG22	1:H:78:LEU:HB3	1.89	0.54
1:G:108:ALA:HB2	1:G:147:ILE:HD11	1.88	0.54
1:B:35:ILE:HG12	1:B:75:VAL:HG21	1.89	0.54
1:B:14:ILE:HD13	1:B:260:MET:HG3	1.90	0.53
1:G:26:SER:OG	1:G:29:ARG:NH2	2.41	0.53
1:E:188:ILE:HG21	1:E:207:ILE:HG13	1.90	0.53
1:C:87:HIS:CE1	1:C:88:GLU:HG3	2.44	0.53
1:D:188:ILE:HG21	1:D:207:ILE:HG13	1.90	0.53
1:C:37:ASN:OD1	1:C:296:LYS:NZ	2.37	0.53
1:C:189:SER:HB3	1:C:206:VAL:HG12	1.91	0.52
1:F:154:LYS:HD3	6:F:302:PGE:H62	1.91	0.52
1:A:166:KPI:HDA	1:A:207:ILE:HD12	1.90	0.52
1:C:20:GLY:N	1:C:270:GLU:OE2	2.43	0.52
1:H:166:KPI:HDA	1:H:207:ILE:HD12	1.90	0.52
1:H:45:VAL:HG13	1:H:54:LEU:HD12	1.92	0.52
1:B:246:ILE:HG12	1:B:289:VAL:HG11	1.92	0.52
1:C:149:THR:HG23	1:C:178:LEU:HD23	1.92	0.52
1:G:34:GLN:HE22	1:G:215:PRO:HG3	1.77	0.51
1:A:256:ILE:O	1:A:260:MET:HG2	2.11	0.50
1:E:149:THR:O	1:E:153:ILE:HG12	2.11	0.50
1:G:7:ILE:HB	1:G:204:LYS:O	2.11	0.50
1:B:67:VAL:HG11	1:B:100:GLY:HA3	1.92	0.50
1:H:189:SER:HB3	1:H:206:VAL:HG12	1.92	0.50
1:B:58:GLU:OE2	1:B:272:ARG:NH2	2.32	0.50
1:D:23:ASP:OD1	1:D:23:ASP:O	2.30	0.50
1:F:45:VAL:HG13	1:F:54:LEU:HD12	1.93	0.50
1:A:98:GLU:OE1	1:A:99:HIS:ND1	2.44	0.49
1:F:254:ILE:HA	1:F:271:PHE:CE1	2.46	0.49
1:E:76:LYS:HE2	1:E:102:ASP:HB3	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:35:ILE:HG12	1:F:75:VAL:HG21	1.93	0.49
1:G:68:GLU:O	1:G:71:LYS:HG2	2.13	0.49
1:E:47:THR:HG1	1:F:111:TYR:HH	1.60	0.49
1:C:52:ALA:O	2:C:301:LYS:N	2.45	0.49
1:G:20:GLY:O	1:G:272:ARG:NH2	2.44	0.49
1:A:106:SER:HB3	1:A:123:TYR:HE1	1.77	0.49
1:D:105:LEU:HD13	1:D:135:LEU:HD23	1.94	0.49
1:H:188:ILE:HG21	1:H:207:ILE:HG13	1.96	0.48
1:G:130:VAL:HG13	1:G:132:ILE:H	1.78	0.48
1:C:216:ASP:N	1:C:216:ASP:OD1	2.47	0.48
1:D:97:LYS:NZ	1:D:131:ASP:OD1	2.33	0.48
1:D:55:THR:HG22	1:D:57:GLU:H	1.79	0.48
1:G:84:ASN:OD1	2:H:301:LYS:N	2.47	0.48
1:F:192:ASP:OD2	1:F:192:ASP:N	2.47	0.48
1:G:124:LYS:O	1:G:128:GLN:HG3	2.14	0.48
1:C:213:LEU:HD11	1:C:295:ILE:HD13	1.96	0.47
1:G:242:ASN:OD1	1:G:292:LYS:NZ	2.43	0.47
1:F:45:VAL:HG12	1:F:51:SER:HB3	1.95	0.47
1:F:196:TYR:OH	1:F:230:TYR:HB3	2.13	0.47
1:C:97:LYS:NZ	1:C:131:ASP:OD1	2.35	0.47
1:G:37:ASN:OD1	1:G:296:LYS:NZ	2.45	0.47
1:B:97:LYS:HE3	1:B:131:ASP:H	1.79	0.47
1:E:256:ILE:O	1:E:260:MET:HG2	2.14	0.47
1:E:6:ILE:HG12	1:E:76:LYS:HD3	1.96	0.47
1:G:53:THR:HB	1:G:272:ARG:HD2	1.96	0.47
1:F:18:LYS:HB2	1:F:23:ASP:HB2	1.96	0.47
1:H:261:TYR:HA	1:H:266:ILE:HG22	1.98	0.46
1:B:166:KPI:HDA	1:B:207:ILE:HD12	1.98	0.46
1:F:232:GLU:O	1:F:235:LYS:HG2	2.15	0.46
1:D:45:VAL:HG21	1:D:59:HIS:CD2	2.50	0.46
1:E:200:SER:HA	1:E:230:TYR:HE1	1.80	0.46
1:B:138:ASN:ND2	1:B:167:GLU:OE2	2.37	0.46
1:C:260:MET:HA	1:C:260:MET:HE2	1.98	0.46
1:H:98:GLU:HG3	1:H:99:HIS:ND1	2.31	0.46
1:G:31:ILE:HD13	1:G:66:ALA:HA	1.98	0.46
1:C:45:VAL:HG13	1:C:54:LEU:HD12	1.98	0.46
1:C:97:LYS:HD3	1:C:132:ILE:HG23	1.97	0.46
1:D:37:ASN:HB3	1:D:215:PRO:HB2	1.98	0.46
1:D:216:ASP:OD1	1:D:216:ASP:N	2.46	0.46
1:F:59:HIS:NE2	1:F:80:GLY:O	2.46	0.46
1:H:216:ASP:N	1:H:216:ASP:OD1	2.49	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:107:VAL:HG23	1:E:137:TYR:CD2	2.51	0.45
1:E:45:VAL:HG21	1:E:59:HIS:CD2	2.51	0.45
1:B:10:MET:HG2	1:B:41:ALA:HB3	1.97	0.45
1:D:246:ILE:HD11	1:D:285:LYS:HB3	1.99	0.45
1:H:266:ILE:HD12	1:H:266:ILE:HA	1.87	0.45
1:B:166:KPI:HE	1:B:207:ILE:HB	1.99	0.45
1:C:154:LYS:HD2	5:C:302:PG4:H61	1.99	0.45
1:G:256:ILE:O	1:G:260:MET:HG2	2.17	0.45
1:B:216:ASP:OD1	1:B:216:ASP:N	2.50	0.45
1:C:254:ILE:HB	1:C:255:PRO:HD3	1.99	0.44
1:D:102:ASP:O	1:D:133:PRO:HD2	2.18	0.44
1:D:243:ILE:O	1:D:247:LEU:HG	2.17	0.44
1:B:6:ILE:HG23	1:B:76:LYS:HD2	1.99	0.44
1:C:35:ILE:HG12	1:C:75:VAL:HG21	2.00	0.44
1:D:21:LYS:HA	1:D:21:LYS:HD3	1.54	0.44
1:A:108:ALA:HB2	1:A:147:ILE:HD11	1.99	0.44
1:H:17:PHE:HB2	1:H:270:GLU:HB3	1.99	0.44
1:E:140:PRO:HG3	1:E:146:GLU:HG3	2.00	0.44
1:D:254:ILE:CB	1:D:255:PRO:HD3	2.41	0.44
1:G:97:LYS:HB2	1:G:97:LYS:HE2	1.88	0.44
1:F:51:SER:O	2:F:301:LYS:NZ	2.51	0.44
1:D:98:GLU:HG3	1:D:99:HIS:ND1	2.32	0.44
1:E:152:ILE:HG22	1:E:185:MET:HE1	1.99	0.44
1:F:52:ALA:O	2:F:301:LYS:N	2.51	0.44
1:G:97:LYS:HD2	1:G:130:VAL:HA	2.00	0.44
1:F:172:ILE:HG23	1:H:194:ILE:HG21	1.99	0.44
1:G:11:THR:HG22	1:G:39:ILE:HG12	2.00	0.43
1:A:243:ILE:O	1:A:247:LEU:HG	2.17	0.43
1:A:50:GLU:HG2	1:A:275:LEU:HD12	1.99	0.43
1:F:149:THR:HG23	1:F:178:LEU:HD13	2.00	0.43
1:A:149:THR:HG23	1:A:178:LEU:HD23	2.00	0.43
1:B:262:LEU:HD21	1:B:283:PHE:CZ	2.54	0.43
1:E:160:GLU:CD	1:E:160:GLU:H	2.21	0.43
1:B:149:THR:HG23	1:B:178:LEU:HG	1.99	0.43
1:C:34:GLN:OE1	1:C:39:ILE:HG13	2.18	0.43
1:F:110:PHE:O	1:F:111:TYR:HB3	2.18	0.43
1:G:3:LYS:HD3	1:G:163:TYR:HA	2.00	0.43
1:C:57:GLU:O	1:C:61:THR:OG1	2.23	0.43
1:B:45:VAL:HG21	1:B:59:HIS:CD2	2.53	0.43
1:F:178:LEU:HD13	1:F:178:LEU:HA	1.88	0.43
1:F:188:ILE:HG21	1:F:207:ILE:HG13	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:285:LYS:HB2	1:F:285:LYS:HE3	1.78	0.43
1:F:194:ILE:HG21	1:H:172:ILE:HG23	2.00	0.43
1:E:243:ILE:O	1:E:247:LEU:HG	2.19	0.43
1:H:13:LEU:HD11	1:H:42:VAL:HB	2.00	0.43
1:A:204:LYS:HD2	3:A:305:EDO:H22	2.00	0.43
1:D:56:HIS:NE2	1:D:88:GLU:OE1	2.47	0.43
1:B:135:LEU:HD13	1:B:163:TYR:CZ	2.54	0.42
1:D:71:LYS:HD3	1:D:71:LYS:HA	1.62	0.42
1:H:104:ILE:HG12	1:H:134:VAL:HG22	2.01	0.42
1:H:154:LYS:HD2	6:H:302:PGE:H6	2.00	0.42
1:B:102:ASP:OD2	1:B:102:ASP:N	2.50	0.42
1:H:166:KPI:HE	1:H:207:ILE:HB	2.02	0.42
1:A:3:LYS:HD3	1:A:163:TYR:HA	2.02	0.42
1:B:262:LEU:HD21	1:B:283:PHE:CE2	2.54	0.42
1:A:135:LEU:HD13	1:A:163:TYR:CZ	2.54	0.42
1:E:245:LYS:HD2	1:E:245:LYS:HA	1.79	0.42
1:G:34:GLN:NE2	1:G:215:PRO:HG3	2.35	0.42
1:H:45:VAL:HG11	1:H:59:HIS:CG	2.54	0.42
1:B:60:ARG:HG2	1:B:95:PHE:CD2	2.54	0.42
1:E:254:ILE:HB	1:E:255:PRO:HD3	2.01	0.42
1:H:45:VAL:HG12	1:H:51:SER:HB3	2.01	0.42
1:E:194:ILE:HG21	1:G:172:ILE:HG23	2.01	0.42
1:H:14:ILE:HB	1:H:257:LYS:HD3	2.02	0.42
1:G:135:LEU:HD13	1:G:163:TYR:CZ	2.55	0.42
1:D:43:VAL:HG22	1:D:78:LEU:HB3	2.02	0.42
1:F:138:ASN:HD21	1:F:146:GLU:HG3	1.84	0.42
1:A:189:SER:HB3	1:A:206:VAL:HG12	2.01	0.42
1:A:196:TYR:OH	1:A:230:TYR:HB3	2.20	0.41
1:B:256:ILE:O	1:B:260:MET:HG2	2.19	0.41
1:D:7:ILE:HB	1:D:204:LYS:O	2.21	0.41
1:C:14:ILE:HD13	1:C:260:MET:HG3	2.01	0.41
1:A:243:ILE:HA	1:A:246:ILE:HG22	2.02	0.41
1:E:82:GLY:HA3	1:E:107:VAL:HG12	2.02	0.41
1:H:173:ASP:HA	1:H:176:VAL:HG12	2.02	0.41
1:A:110:PHE:O	1:A:111:TYR:HB3	2.21	0.41
1:B:161:ASN:N	1:B:161:ASN:OD1	2.46	0.41
1:E:87:HIS:CE1	1:E:88:GLU:HG3	2.55	0.41
1:D:182:GLU:N	1:D:183:PRO:HD3	2.34	0.41
1:H:106:SER:HB3	1:H:123:TYR:HE1	1.85	0.41
1:G:213:LEU:HD11	1:G:295:ILE:HD13	2.02	0.41
1:F:12:ALA:HA	1:F:43:VAL:HB	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:254:ILE:HA	1:G:271:PHE:CE1	2.55	0.41
1:A:60:ARG:HG2	1:A:95:PHE:CD2	2.56	0.41
1:C:166:KPI:HDA	1:C:207:ILE:HD12	2.02	0.41
1:E:243:ILE:HB	1:E:293:TYR:CE2	2.56	0.41
1:H:281:GLU:N	1:H:281:GLU:OE1	2.36	0.41
1:D:261:TYR:HA	1:D:266:ILE:HG13	2.01	0.41
1:A:262:LEU:HD21	1:A:283:PHE:CE2	2.55	0.41
1:E:200:SER:HA	1:E:230:TYR:CE1	2.56	0.40
1:G:288:GLU:O	1:G:291:LYS:HG2	2.21	0.40
1:A:275:LEU:HD23	1:A:275:LEU:HA	1.92	0.40
1:C:74:LYS:H	1:C:74:LYS:HG3	1.66	0.40
1:D:257:LYS:HD2	1:D:271:PHE:CE1	2.56	0.40
1:G:257:LYS:HD2	1:G:271:PHE:CE1	2.56	0.40
1:C:120:TYR:OH	1:C:158:ASP:OD2	2.29	0.40
1:E:60:ARG:HB3	1:E:95:PHE:CZ	2.56	0.40
1:F:246:ILE:HG12	1:F:289:VAL:HG21	2.03	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	294/306 (96%)	287 (98%)	7 (2%)	0	100	100
1	B	293/306 (96%)	286 (98%)	7 (2%)	0	100	100
1	C	292/306 (95%)	284 (97%)	8 (3%)	0	100	100
1	D	293/306 (96%)	285 (97%)	8 (3%)	0	100	100
1	E	291/306 (95%)	284 (98%)	7 (2%)	0	100	100
1	F	293/306 (96%)	284 (97%)	9 (3%)	0	100	100
1	G	293/306 (96%)	285 (97%)	7 (2%)	1 (0%)	44	49

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	H	292/306 (95%)	280 (96%)	12 (4%)	0	100	100
All	All	2341/2448 (96%)	2275 (97%)	65 (3%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	G	129	SER

5.3.2 Protein sidechains [i](#)

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	249/257 (97%)	246 (99%)	3 (1%)	75	86
1	B	248/257 (96%)	239 (96%)	9 (4%)	40	50
1	C	247/257 (96%)	241 (98%)	6 (2%)	54	67
1	D	248/257 (96%)	240 (97%)	8 (3%)	44	56
1	E	246/257 (96%)	243 (99%)	3 (1%)	75	86
1	F	248/257 (96%)	241 (97%)	7 (3%)	49	61
1	G	248/257 (96%)	242 (98%)	6 (2%)	54	67
1	H	247/257 (96%)	236 (96%)	11 (4%)	32	39
All	All	1981/2056 (96%)	1928 (97%)	53 (3%)	50	62

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	106	SER
1	A	113	LYS
1	A	146	GLU
1	B	6	ILE
1	B	68	GLU
1	B	97	LYS
1	B	102	ASP

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Mol	Chain	Res	Type
1	B	106	SER
1	B	146	GLU
1	B	147	ILE
1	B	228	GLU
1	B	281	GLU
1	C	27	TYR
1	C	74	LYS
1	C	191	GLU
1	C	227	ASP
1	C	277	SER
1	C	288	GLU
1	D	21	LYS
1	D	71	LYS
1	D	73	THR
1	D	97	LYS
1	D	146	GLU
1	D	150	ASP
1	D	182	GLU
1	D	217	MET
1	E	36	GLU
1	E	227	ASP
1	E	235	LYS
1	F	18	LYS
1	F	173	ASP
1	F	178	LEU
1	F	192	ASP
1	F	231	LYS
1	F	281	GLU
1	F	294	LYS
1	G	130	VAL
1	G	146	GLU
1	G	161	ASN
1	G	228	GLU
1	G	235	LYS
1	G	283	PHE
1	H	4	ASN
1	H	5	ILE
1	H	6	ILE
1	H	57	GLU
1	H	71	LYS
1	H	74	LYS
1	H	145	CYS

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Mol	Chain	Res	Type
1	H	146	GLU
1	H	201	ASN
1	H	265	LEU
1	H	266	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	G	34	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

8 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	A	166	1	11,13,14	2.58	4 (36%)	7,15,17	1.52	1 (14%)
1	KPI	B	166	1	11,13,14	2.56	4 (36%)	7,15,17	1.33	1 (14%)
1	KPI	C	166	1	11,13,14	2.55	4 (36%)	7,15,17	1.20	0
1	KPI	D	166	1	11,13,14	2.54	4 (36%)	7,15,17	1.18	0
1	KPI	E	166	1	11,13,14	2.60	4 (36%)	7,15,17	1.43	1 (14%)
1	KPI	F	166	1	11,13,14	2.58	4 (36%)	7,15,17	1.12	0
1	KPI	G	166	1	11,13,14	2.55	4 (36%)	7,15,17	1.14	0
1	KPI	H	166	1	11,13,14	2.55	4 (36%)	7,15,17	1.34	1 (14%)

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	A	166	1	-	0/8/14/16	0/0/0/0
1	KPI	B	166	1	-	0/8/14/16	0/0/0/0
1	KPI	C	166	1	-	0/8/14/16	0/0/0/0
1	KPI	D	166	1	-	0/8/14/16	0/0/0/0
1	KPI	E	166	1	-	0/8/14/16	0/0/0/0
1	KPI	F	166	1	-	0/8/14/16	0/0/0/0
1	KPI	G	166	1	-	0/8/14/16	0/0/0/0
1	KPI	H	166	1	-	0/8/14/16	0/0/0/0

All (32) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	166	KPI	C1-CX1	-5.63	1.39	1.50
1	G	166	KPI	C1-CX1	-5.61	1.39	1.50
1	H	166	KPI	C1-CX1	-5.61	1.39	1.50
1	B	166	KPI	C1-CX1	-5.60	1.39	1.50
1	F	166	KPI	C1-CX1	-5.59	1.39	1.50
1	E	166	KPI	C1-CX1	-5.59	1.39	1.50
1	C	166	KPI	C1-CX1	-5.54	1.39	1.50
1	D	166	KPI	C1-CX1	-5.54	1.39	1.50
1	E	166	KPI	CB-CA	-3.54	1.48	1.53
1	H	166	KPI	CB-CA	-3.27	1.49	1.53
1	B	166	KPI	CB-CA	-3.24	1.49	1.53
1	C	166	KPI	CB-CA	-3.23	1.49	1.53
1	D	166	KPI	CB-CA	-3.20	1.49	1.53
1	A	166	KPI	CB-CA	-3.16	1.49	1.53
1	F	166	KPI	CB-CA	-3.14	1.49	1.53
1	G	166	KPI	CB-CA	-3.11	1.49	1.53
1	F	166	KPI	CA-C	-2.74	1.46	1.50
1	A	166	KPI	CA-C	-2.64	1.46	1.50
1	E	166	KPI	CA-C	-2.49	1.47	1.50
1	B	166	KPI	CA-C	-2.39	1.47	1.50
1	H	166	KPI	CA-C	-2.37	1.47	1.50
1	D	166	KPI	CA-C	-2.34	1.47	1.50
1	C	166	KPI	CA-C	-2.30	1.47	1.50
1	G	166	KPI	CA-C	-2.24	1.47	1.50
1	A	166	KPI	CX1-NZ	4.45	1.36	1.28
1	B	166	KPI	CX1-NZ	4.47	1.36	1.28
1	H	166	KPI	CX1-NZ	4.48	1.36	1.28
1	E	166	KPI	CX1-NZ	4.50	1.36	1.28
1	D	166	KPI	CX1-NZ	4.52	1.36	1.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	166	KPI	CX1-NZ	4.53	1.36	1.28
1	C	166	KPI	CX1-NZ	4.60	1.36	1.28
1	G	166	KPI	CX1-NZ	4.62	1.36	1.28

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	166	KPI	CD-CE-NZ	-2.85	106.24	110.71
1	E	166	KPI	CD-CE-NZ	-2.64	106.56	110.71
1	B	166	KPI	CD-CE-NZ	-2.26	107.17	110.71
1	H	166	KPI	CB-CA-C	-2.06	108.26	111.65

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

7 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	166	KPI	1	0
1	B	166	KPI	2	0
1	C	166	KPI	1	0
1	D	166	KPI	1	0
1	F	166	KPI	1	0
1	G	166	KPI	1	0
1	H	166	KPI	2	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

27 ligands 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	$\# Z > 2$
2	LYS	A	301	-	4,9,9	0.25	0	3,10,10	0.42	0
2	LYS	A	302	-	4,9,9	0.25	0	3,10,10	0.46	0
3	EDO	A	303	-	3,3,3	0.57	0	2,2,2	0.32	0
4	ACT	A	304	-	1,3,3	1.37	0	0,3,3	0.00	-
3	EDO	A	305	-	3,3,3	0.55	0	2,2,2	0.31	0
3	EDO	A	306	-	3,3,3	0.57	0	2,2,2	0.36	0
2	LYS	B	301	-	4,9,9	0.24	0	3,10,10	0.48	0
5	PG4	B	302	-	12,12,12	0.68	0	11,11,11	0.74	0
6	PGE	B	303	-	9,9,9	0.68	0	8,8,8	0.75	0
2	LYS	C	301	-	4,9,9	0.24	0	3,10,10	0.48	0
5	PG4	C	302	-	12,12,12	0.71	0	11,11,11	0.73	0
6	PGE	C	303	-	9,9,9	0.67	0	8,8,8	0.72	0
7	GOL	C	304	-	5,5,5	0.35	0	5,5,5	0.24	0
3	EDO	D	301	-	3,3,3	0.57	0	2,2,2	0.36	0
2	LYS	E	301	-	4,9,9	0.22	0	3,10,10	0.48	0
6	PGE	E	302	-	9,9,9	0.68	0	8,8,8	0.70	0
3	EDO	E	303	-	3,3,3	0.56	0	2,2,2	0.33	0
3	EDO	E	304	-	3,3,3	0.56	0	2,2,2	0.32	0
2	LYS	F	301	-	4,9,9	0.22	0	3,10,10	0.47	0
6	PGE	F	302	-	9,9,9	0.68	0	8,8,8	0.73	0
4	ACT	F	303	-	1,3,3	1.37	0	0,3,3	0.00	-
2	LYS	G	301	-	4,9,9	0.22	0	3,10,10	0.43	0
3	EDO	G	302	-	3,3,3	0.59	0	2,2,2	0.36	0
3	EDO	G	303	-	3,3,3	0.57	0	2,2,2	0.36	0
3	EDO	G	304	-	3,3,3	0.56	0	2,2,2	0.34	0
2	LYS	H	301	-	4,9,9	0.22	0	3,10,10	0.48	0
6	PGE	H	302	-	9,9,9	0.66	0	8,8,8	0.79	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	LYS	A	301	-	-	0/5/9/9	0/0/0/0
2	LYS	A	302	-	-	0/5/9/9	0/0/0/0
3	EDO	A	303	-	-	0/1/1/1	0/0/0/0
4	ACT	A	304	-	-	0/0/0/0	0/0/0/0
3	EDO	A	305	-	-	0/1/1/1	0/0/0/0
3	EDO	A	306	-	-	0/1/1/1	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	LYS	B	301	-	-	0/5/9/9	0/0/0/0
5	PG4	B	302	-	-	0/10/10/10	0/0/0/0
6	PGE	B	303	-	-	0/7/7/7	0/0/0/0
2	LYS	C	301	-	-	0/5/9/9	0/0/0/0
5	PG4	C	302	-	-	0/10/10/10	0/0/0/0
6	PGE	C	303	-	-	0/7/7/7	0/0/0/0
7	GOL	C	304	-	-	0/4/4/4	0/0/0/0
3	EDO	D	301	-	-	0/1/1/1	0/0/0/0
2	LYS	E	301	-	-	0/5/9/9	0/0/0/0
6	PGE	E	302	-	-	0/7/7/7	0/0/0/0
3	EDO	E	303	-	-	0/1/1/1	0/0/0/0
3	EDO	E	304	-	-	0/1/1/1	0/0/0/0
2	LYS	F	301	-	-	0/5/9/9	0/0/0/0
6	PGE	F	302	-	-	0/7/7/7	0/0/0/0
4	ACT	F	303	-	-	0/0/0/0	0/0/0/0
2	LYS	G	301	-	-	0/5/9/9	0/0/0/0
3	EDO	G	302	-	-	0/1/1/1	0/0/0/0
3	EDO	G	303	-	-	0/1/1/1	0/0/0/0
3	EDO	G	304	-	-	0/1/1/1	0/0/0/0
2	LYS	H	301	-	-	0/5/9/9	0/0/0/0
6	PGE	H	302	-	-	0/7/7/7	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

10 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	302	LYS	1	0
3	A	305	EDO	1	0
2	C	301	LYS	1	0
5	C	302	PG4	1	0
7	C	304	GOL	1	0
2	F	301	LYS	2	0
6	F	302	PGE	1	0
4	F	303	ACT	1	0
2	H	301	LYS	2	0
6	H	302	PGE	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	295/306 (96%)	0.05	0 100 100	29, 40, 57, 77	0
1	B	295/306 (96%)	0.22	3 (1%) 82 81	32, 45, 67, 88	1 (0%)
1	C	294/306 (96%)	0.13	1 (0%) 93 93	32, 45, 67, 85	0
1	D	294/306 (96%)	0.25	4 (1%) 75 73	32, 51, 73, 84	0
1	E	293/306 (95%)	0.13	2 (0%) 87 86	33, 46, 64, 77	1 (0%)
1	F	295/306 (96%)	0.12	3 (1%) 82 81	34, 47, 64, 72	1 (0%)
1	G	295/306 (96%)	0.30	6 (2%) 65 63	37, 52, 71, 85	0
1	H	294/306 (96%)	0.44	14 (4%) 31 30	40, 58, 80, 92	1 (0%)
All	All	2355/2448 (96%)	0.20	33 (1%) 75 73	29, 48, 70, 92	4 (0%)

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	298	PHE	4.2
1	H	127	ALA	4.1
1	F	287	GLU	3.6
1	H	287	GLU	3.5
1	D	292	LYS	3.4
1	H	289	VAL	3.3
1	E	298	PHE	3.2
1	H	261	TYR	3.2
1	F	234	LYS	3.1
1	H	288	GLU	3.1
1	G	20	GLY	3.0
1	H	15	THR	2.8
1	G	230	TYR	2.7
1	D	289	VAL	2.7
1	G	4	ASN	2.6
1	E	160	GLU	2.6

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Mol	Chain	Res	Type	RSRZ
1	H	29	ARG	2.6
1	H	284	ALA	2.6
1	B	95	PHE	2.5
1	H	5	ILE	2.3
1	B	129	SER	2.3
1	F	227	ASP	2.3
1	H	25	GLN	2.3
1	H	283	PHE	2.3
1	B	267	GLU	2.2
1	D	29	ARG	2.2
1	H	269	LEU	2.2
1	D	265	LEU	2.2
1	C	17	PHE	2.1
1	G	114	PRO	2.1
1	G	262	LEU	2.1
1	H	255	PRO	2.1
1	G	28	ALA	2.0

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

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 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	LLDF	B-factors(Å ²)	Q<0.9
1	KPI	G	166	14/15	0.90	0.13	-	45,55,70,72	0
1	KPI	D	166	14/15	0.89	0.18	-	32,46,63,67	0
1	KPI	A	166	14/15	0.90	0.15	-	26,37,54,55	0
1	KPI	F	166	14/15	0.85	0.16	-	29,42,55,55	0
1	KPI	C	166	14/15	0.95	0.13	-	26,44,61,63	0
1	KPI	B	166	14/15	0.93	0.16	-	30,38,57,57	0
1	KPI	H	166	14/15	0.89	0.20	-	35,45,78,89	0
1	KPI	E	166	14/15	0.94	0.20	-	29,44,74,82	0

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 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	LLDF	B-factors(Å ²)	Q<0.9
6	PGE	B	303	10/10	0.60	0.29	4.43	60,78,83,84	0
2	LYS	C	301	10/10	0.82	0.23	4.42	51,60,67,75	0
2	LYS	B	301	10/10	0.79	0.23	4.08	49,59,73,83	0
5	PG4	C	302	13/13	0.88	0.18	3.81	44,60,72,72	0
3	EDO	G	302	4/4	0.87	0.21	2.81	46,48,55,57	0
2	LYS	G	301	10/10	0.84	0.21	2.76	40,49,75,81	0
3	EDO	D	301	4/4	0.83	0.19	2.52	57,61,68,73	0
2	LYS	A	301	10/10	0.82	0.19	1.45	35,47,59,61	0
2	LYS	H	301	10/10	0.85	0.19	1.26	48,55,69,69	0
2	LYS	E	301	10/10	0.95	0.16	1.23	41,47,64,68	0
4	ACT	F	303	4/4	0.86	0.17	1.18	56,57,62,64	0
6	PGE	E	302	10/10	0.73	0.20	1.09	70,72,72,72	0
2	LYS	F	301	10/10	0.89	0.16	0.87	36,48,55,59	0
2	LYS	A	302	10/10	0.83	0.16	0.68	38,56,59,61	0
6	PGE	F	302	10/10	0.82	0.19	0.65	60,64,74,74	0
3	EDO	E	303	4/4	0.76	0.16	0.13	64,68,72,74	0
6	PGE	C	303	10/10	0.76	0.15	0.10	60,65,74,74	0
3	EDO	G	303	4/4	0.76	0.16	-0.29	58,59,60,62	0
5	PG4	B	302	13/13	0.87	0.14	-0.44	53,60,68,71	0
6	PGE	H	302	10/10	0.85	0.12	-1.26	53,72,80,80	0
3	EDO	A	305	4/4	0.60	0.21	-	57,64,66,67	0
3	EDO	G	304	4/4	0.70	0.14	-	60,61,62,63	0
3	EDO	E	304	4/4	0.77	0.17	-	54,57,59,59	0
4	ACT	A	304	4/4	0.67	0.26	-	71,72,73,75	0
3	EDO	A	303	4/4	0.72	0.22	-	47,49,53,56	0
3	EDO	A	306	4/4	0.79	0.16	-	57,60,61,62	0
7	GOL	C	304	6/6	0.66	0.30	-	58,66,66,70	0

6.5 Other polymers

There are no such residues in this entry.