



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2026 – 07:00 AM UTC

PDB ID : 7LOY / pdb_00007loy
Title : Dihydropicolinate synthase with pyruvate from Candidatus Liberibacter solanacearum
Authors : Gilkes, J.M.; Frampton, R.A.; Board, A.; Sheen, C.R.; Smith, G.R.; Dobson, R.C.J.
Deposited on : 2021-02-11
Resolution : 2.40 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

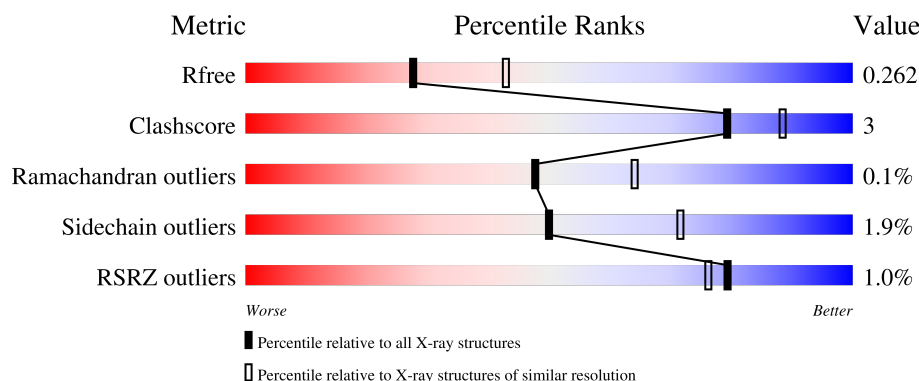
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

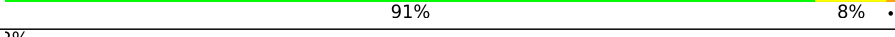
The reported resolution of this entry is 2.40 Å.

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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	295	 91% 9%
1	B	295	 90% 10%
1	C	295	 91% 9%
1	D	295	 91% 8%
1	E	295	 88% 12%

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Mol	Chain	Length	Quality of chain
1	F	295	 92%8%

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 13864 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	295	Total	C	N	O	S	0	0	0
			2236	1408	377	439	12			
1	B	295	Total	C	N	O	S	0	0	0
			2236	1408	377	439	12			
1	C	295	Total	C	N	O	S	0	0	0
			2236	1408	377	439	12			
1	D	295	Total	C	N	O	S	0	0	0
			2236	1408	377	439	12			
1	E	295	Total	C	N	O	S	0	0	0
			2236	1408	377	439	12			
1	F	295	Total	C	N	O	S	0	0	0
			2236	1408	377	439	12			

- Molecule 2 is water.

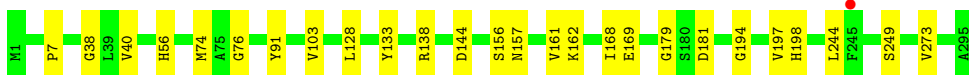
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	125	Total	O	0	0
			125	125		
2	B	108	Total	O	0	0
			108	108		
2	C	55	Total	O	0	0
			55	55		
2	D	64	Total	O	0	0
			64	64		
2	E	47	Total	O	0	0
			47	47		
2	F	49	Total	O	0	0
			49	49		

3 Residue-property plots

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

Chain A: 



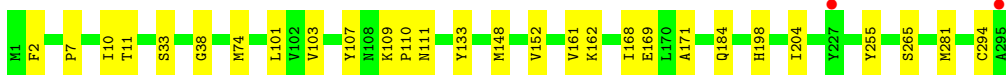
- Molecule 1: 4-hydroxy-tetrahydrodipicolinate synthase

Chain B: 

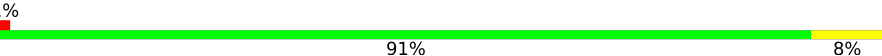


- Molecule 1: 4-hydroxy-tetrahydrodipicolinate synthase

Chain C: 



- Molecule 1: 4-hydroxy-tetrahydrodipicolinate synthase

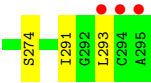
Chain D: 



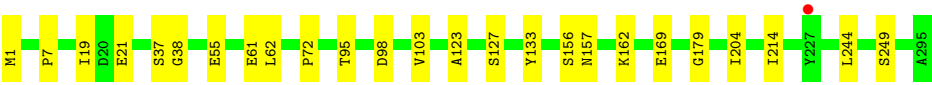
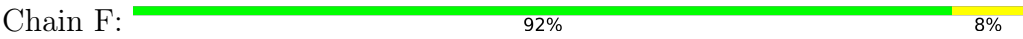
- Molecule 1: 4-hydroxy-tetrahydrodipicolinate synthase

Chain E: 





● Molecule 1: 4-hydroxy-tetrahydrodipicolinate synthase



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	101.78Å 133.58Å 155.99Å 90.00° 100.55° 90.00°	Depositor
Resolution (Å)	46.27 – 2.40 46.27 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.9 (46.27-2.40) 99.8 (46.27-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.30 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.201 , 0.252 (Not available) , 0.262	Depositor DCC
R_{free} test set	4050 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	33.8	Xtriage
Anisotropy	0.027	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 23.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	13864	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.28% 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: 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	1.03	0/2260	1.49	6/3066 (0.2%)
1	B	1.03	1/2260 (0.0%)	1.50	4/3066 (0.1%)
1	C	1.06	1/2260 (0.0%)	1.51	1/3066 (0.0%)
1	D	1.06	0/2260	1.51	0/3066
1	E	1.05	0/2260	1.55	1/3066 (0.0%)
1	F	1.07	0/2260	1.54	3/3066 (0.1%)
All	All	1.05	2/13560 (0.0%)	1.52	15/18396 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2
1	B	0	1
1	C	0	4
1	D	0	1
1	E	0	2
1	F	0	1
All	All	0	11

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	109	LYS	N-CA	5.70	1.50	1.46
1	C	109	LYS	N-CA	5.56	1.50	1.46

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	156	SER	CA-C-N	7.27	132.86	120.72
1	A	156	SER	C-N-CA	7.27	132.86	120.72
1	A	144	ASP	CA-CB-CG	6.75	119.35	112.60
1	F	156	SER	CA-C-N	6.62	130.04	120.38
1	F	156	SER	C-N-CA	6.62	130.04	120.38
1	A	273	VAL	CA-C-N	6.39	128.74	120.44
1	A	273	VAL	C-N-CA	6.39	128.74	120.44
1	C	11	THR	CB-CA-C	6.07	116.52	110.33
1	E	11	THR	CB-CA-C	5.98	116.43	110.33
1	B	179	GLY	CA-C-O	-5.85	117.55	122.29
1	F	179	GLY	CA-C-O	-5.28	118.59	122.23
1	B	22	ASP	CA-C-N	5.11	127.08	120.44
1	B	22	ASP	C-N-CA	5.11	127.08	120.44
1	B	144	ASP	CA-CB-CG	5.09	117.69	112.60
1	A	179	GLY	CA-C-O	-5.06	118.19	122.29

There are no chirality outliers.

All (11) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	161	VAL	Mainchain,Peptide
1	B	161	VAL	Mainchain
1	C	110	PRO	Peptide
1	C	161	VAL	Peptide
1	C	204	ILE	Peptide
1	C	294	CYS	Peptide
1	D	161	VAL	Peptide
1	E	162	KPI	Mainchain
1	E	204	ILE	Peptide
1	F	204	ILE	Peptide

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	2236	0	2220	13	0
1	B	2236	0	2221	13	0
1	C	2236	0	2220	10	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	2236	0	2221	11	0
1	E	2236	0	2220	15	0
1	F	2236	0	2221	13	0
2	A	125	0	0	2	0
2	B	108	0	0	0	0
2	C	55	0	0	0	0
2	D	64	0	0	0	0
2	E	47	0	0	0	0
2	F	49	0	0	1	0
All	All	13864	0	13323	73	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 (73) 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:1:MET:N	2:F:301:HOH:O	2.04	0.89
1:D:74:MET:CE	1:D:101:LEU:HD13	2.13	0.79
1:D:74:MET:HE2	1:D:101:LEU:HD13	1.71	0.73
1:C:255:TYR:CE1	1:C:281:MET:HE3	2.27	0.69
1:E:4:ARG:HA	1:E:219:GLN:HE21	1.57	0.69
1:B:168:ILE:HD12	1:B:198:HIS:CD2	2.29	0.68
1:B:74:MET:CE	1:B:101:LEU:HD13	2.23	0.68
1:F:123:ALA:O	1:F:157:ASN:ND2	2.31	0.62
1:E:103:VAL:HA	1:E:133:TYR:HB3	1.82	0.62
1:C:168:ILE:HD12	1:C:198:HIS:CD2	2.35	0.61
1:A:181:ASP:HB2	2:A:322:HOH:O	2.02	0.59
1:D:171:ALA:HA	1:D:184:GLN:HE22	1.68	0.58
1:A:56:HIS:HD2	2:A:366:HOH:O	1.86	0.57
1:B:197:VAL:HG12	1:B:227:TYR:CE1	2.39	0.57
1:C:103:VAL:HA	1:C:133:TYR:HB3	1.86	0.57
1:E:7:PRO:HD2	1:E:38:GLY:O	2.06	0.55
1:B:82:THR:OG1	1:B:118:HIS:HD2	1.89	0.55
1:B:7:PRO:HD2	1:B:38:GLY:O	2.07	0.55
1:E:291:ILE:HG22	1:E:293:LEU:HD23	1.89	0.54
1:C:171:ALA:HA	1:C:184:GLN:HE22	1.71	0.54
1:F:61:GLU:HA	1:F:95:THR:HG21	1.89	0.54
1:A:103:VAL:HA	1:A:133:TYR:HB3	1.91	0.53
1:F:103:VAL:HA	1:F:133:TYR:HB3	1.91	0.52
1:F:244:LEU:O	1:F:249:SER:HA	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:7:PRO:HD2	1:C:38:GLY:O	2.10	0.52
1:D:103:VAL:HA	1:D:133:TYR:HB3	1.91	0.52
1:E:57:CYS:HB3	1:E:91:TYR:CE2	2.45	0.51
1:B:103:VAL:HA	1:B:133:TYR:HB3	1.94	0.50
1:D:238:PHE:HB3	1:D:239:PRO:HD3	1.92	0.50
1:E:189:ASP:HB3	1:E:241:HIS:CE1	2.46	0.50
1:C:107:TYR:CD1	1:E:138:ARG:HB3	2.47	0.50
1:A:56:HIS:HE1	1:A:76:GLY:O	1.94	0.50
1:F:7:PRO:HD2	1:F:38:GLY:O	2.12	0.49
1:C:255:TYR:HE1	1:C:281:MET:HE3	1.77	0.49
1:A:244:LEU:O	1:A:249:SER:HA	2.13	0.49
1:C:2:PHE:CZ	1:C:74:MET:HB2	2.49	0.48
1:F:95:THR:HG22	1:F:95:THR:O	2.14	0.48
1:F:1:MET:CE	1:F:98:ASP:HB3	2.44	0.48
1:F:214:ILE:H	1:F:214:ILE:HD12	1.79	0.47
1:E:189:ASP:OD1	1:E:205:SER:OG	2.16	0.47
1:E:238:PHE:HB3	1:E:239:PRO:HD3	1.97	0.46
1:F:1:MET:HE1	1:F:98:ASP:HB3	1.98	0.45
1:D:197:VAL:HG12	1:D:227:TYR:CE1	2.51	0.45
1:D:7:PRO:HD2	1:D:38:GLY:O	2.17	0.45
1:B:189:ASP:HB3	1:B:241:HIS:CD2	2.52	0.45
1:E:28:ILE:HG23	1:E:39:LEU:HD11	1.99	0.45
1:B:54:GLU:CD	1:B:54:GLU:H	2.25	0.45
1:D:74:MET:HE2	1:D:101:LEU:HB2	1.98	0.44
1:E:291:ILE:CG2	1:E:293:LEU:HD23	2.46	0.44
1:C:148:MET:O	1:C:152:VAL:HG23	2.16	0.44
1:B:45:THR:HA	1:B:250:ILE:HD11	1.99	0.44
1:E:214:ILE:HG23	1:E:237:LEU:HD21	2.00	0.44
1:A:194:GLY:O	1:A:197:VAL:HG12	2.18	0.43
1:A:128:LEU:O	1:A:157:ASN:OD1	2.35	0.43
1:A:56:HIS:CE1	1:A:76:GLY:O	2.71	0.43
1:A:168:ILE:HD12	1:A:198:HIS:CD2	2.53	0.43
1:D:228:ARG:NE	1:D:228:ARG:HA	2.34	0.42
1:F:21:GLU:HB3	1:F:62:LEU:HD21	2.01	0.42
1:B:132:ILE:O	1:B:161:VAL:HA	2.19	0.42
1:C:101:LEU:HD11	1:C:133:TYR:HB2	2.02	0.42
1:F:19:ILE:HD11	1:F:55:GLU:HB3	2.01	0.42
1:A:91:TYR:C	1:A:91:TYR:CD1	2.97	0.42
1:F:37:SER:O	1:F:72:PRO:HD2	2.20	0.41
1:A:7:PRO:HD2	1:A:38:GLY:O	2.19	0.41
1:E:47:GLU:CD	1:E:250:ILE:HG21	2.45	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:260:LEU:HD23	1:B:293:LEU:HD13	2.02	0.41
1:E:2:PHE:CZ	1:E:74:MET:HB2	2.56	0.41
1:A:40:VAL:HG22	1:A:74:MET:HB3	2.02	0.41
1:A:138:ARG:HB3	1:B:107:TYR:CD1	2.56	0.41
1:D:247:GLU:HB3	1:D:248:PRO:HD2	2.02	0.41
1:D:238:PHE:CE2	1:D:242:GLN:HG3	2.56	0.40
1:E:8:ALA:HA	1:E:40:VAL:HB	2.03	0.40
1:B:74:MET:HE2	1:B:101:LEU:HD13	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	292/295 (99%)	282 (97%)	10 (3%)	0	100	100
1	B	292/295 (99%)	283 (97%)	9 (3%)	0	100	100
1	C	292/295 (99%)	284 (97%)	7 (2%)	1 (0%)	36	50
1	D	292/295 (99%)	278 (95%)	14 (5%)	0	100	100
1	E	292/295 (99%)	279 (96%)	13 (4%)	0	100	100
1	F	292/295 (99%)	280 (96%)	12 (4%)	0	100	100
All	All	1752/1770 (99%)	1686 (96%)	65 (4%)	1 (0%)	48	64

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	111	ASN

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	243/243 (100%)	242 (100%)	1 (0%)	84	92
1	B	243/243 (100%)	240 (99%)	3 (1%)	63	81
1	C	243/243 (100%)	239 (98%)	4 (2%)	55	76
1	D	243/243 (100%)	233 (96%)	10 (4%)	27	46
1	E	239/243 (98%)	232 (97%)	7 (3%)	37	60
1	F	188/243 (77%)	186 (99%)	2 (1%)	65	82
All	All	1399/1458 (96%)	1372 (98%)	27 (2%)	50	71

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	169	GLU
1	B	180	SER
1	B	251	SER
1	B	278	LYS
1	C	10	ILE
1	C	33	SER
1	C	169	GLU
1	C	265	SER
1	D	10	ILE
1	D	48	SER
1	D	62	LEU
1	D	65	LYS
1	D	168	ILE
1	D	169	GLU
1	D	197	VAL
1	D	214	ILE
1	D	248	PRO
1	D	265	SER
1	E	1	MET
1	E	16	ASP
1	E	95	THR
1	E	137	SER

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Mol	Chain	Res	Type
1	E	169	GLU
1	E	265	SER
1	E	274	SER
1	F	127	SER
1	F	169	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	56	HIS
1	A	174	GLN
1	A	196	ASN
1	A	229	GLN
1	A	242	GLN
1	B	90	GLN
1	B	118	HIS
1	B	196	ASN
1	B	241	HIS
1	B	290	HIS
1	C	174	GLN
1	C	184	GLN
1	C	196	ASN
1	D	27	HIS
1	D	135	ASN
1	D	184	GLN
1	D	196	ASN
1	D	242	GLN
1	E	27	HIS
1	E	108	ASN
1	E	219	GLN
1	E	229	GLN
1	F	27	HIS
1	F	108	ASN
1	F	174	GLN
1	F	229	GLN

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	D	162	1	11,13,14	1.14	1 (9%)	9,15,17	3.06	6 (66%)
1	KPI	A	162	1	11,13,14	1.40	1 (9%)	9,15,17	1.98	1 (11%)
1	KPI	B	162	1	11,13,14	1.15	1 (9%)	9,15,17	2.25	5 (55%)
1	KPI	C	162	1	11,13,14	1.27	2 (18%)	9,15,17	2.63	3 (33%)
1	KPI	F	162	1	11,13,14	1.38	2 (18%)	9,15,17	2.26	2 (22%)
1	KPI	E	162	1	11,13,14	1.47	1 (9%)	9,15,17	2.68	3 (33%)

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	D	162	1	-	8/13/14/16	-
1	KPI	A	162	1	-	1/13/14/16	-
1	KPI	B	162	1	-	0/13/14/16	-
1	KPI	C	162	1	-	2/13/14/16	-
1	KPI	F	162	1	-	0/13/14/16	-
1	KPI	E	162	1	-	0/13/14/16	-

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	162	KPI	CX2-CX1	-3.82	1.45	1.49
1	A	162	KPI	CX2-CX1	-3.37	1.45	1.49
1	F	162	KPI	CX2-CX1	-2.94	1.46	1.49
1	C	162	KPI	O-C	2.52	1.29	1.20
1	B	162	KPI	O-C	2.22	1.28	1.20
1	F	162	KPI	O-C	2.11	1.28	1.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	162	KPI	O1-CX2	-2.07	1.24	1.30
1	C	162	KPI	CX2-CX1	-2.01	1.47	1.49

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	162	KPI	C1-CX1-CX2	6.92	124.64	118.11
1	C	162	KPI	C1-CX1-CX2	6.29	124.05	118.11
1	F	162	KPI	C1-CX1-CX2	5.65	123.44	118.11
1	A	162	KPI	C1-CX1-CX2	5.19	123.00	118.11
1	D	162	KPI	CE-NZ-CX1	5.15	135.73	121.58
1	D	162	KPI	CX2-CX1-NZ	5.07	126.83	114.88
1	B	162	KPI	C1-CX1-CX2	4.06	121.94	118.11
1	C	162	KPI	CE-NZ-CX1	3.32	130.71	121.58
1	D	162	KPI	O2-CX2-CX1	-2.78	117.95	121.35
1	E	162	KPI	CD-CE-NZ	-2.76	105.79	110.67
1	D	162	KPI	O1-CX2-CX1	2.73	122.33	116.50
1	B	162	KPI	CE-NZ-CX1	2.59	128.68	121.58
1	D	162	KPI	CD-CE-NZ	-2.46	106.31	110.67
1	D	162	KPI	C1-CX1-CX2	2.41	120.39	118.11
1	B	162	KPI	CX2-CX1-NZ	2.31	120.31	114.88
1	C	162	KPI	O2-CX2-CX1	-2.29	118.55	121.35
1	B	162	KPI	O1-CX2-CX1	2.16	121.09	116.50
1	B	162	KPI	CD-CE-NZ	-2.13	106.90	110.67
1	F	162	KPI	CE-NZ-CX1	2.05	127.20	121.58
1	E	162	KPI	CE-NZ-CX1	2.04	127.17	121.58

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	D	162	KPI	N-CA-CB-CG
1	D	162	KPI	C-CA-CB-CG
1	D	162	KPI	O-C-CA-CB
1	D	162	KPI	CG-CD-CE-NZ
1	D	162	KPI	CA-CB-CG-CD
1	D	162	KPI	C1-CX1-NZ-CE
1	D	162	KPI	CE-CD-CG-CB
1	D	162	KPI	CX2-CX1-NZ-CE
1	A	162	KPI	CG-CD-CE-NZ
1	C	162	KPI	CG-CD-CE-NZ
1	C	162	KPI	NZ-CX1-CX2-O1

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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	294/295 (99%)	-0.47	1 (0%) 90 88	19, 27, 42, 64	0
1	B	294/295 (99%)	-0.41	3 (1%) 79 76	19, 29, 46, 65	0
1	C	294/295 (99%)	-0.16	2 (0%) 84 81	25, 37, 54, 74	0
1	D	294/295 (99%)	-0.11	4 (1%) 73 69	25, 39, 57, 77	0
1	E	294/295 (99%)	0.15	6 (2%) 65 60	32, 44, 60, 111	0
1	F	294/295 (99%)	0.16	1 (0%) 90 88	31, 46, 64, 81	0
All	All	1764/1770 (99%)	-0.14	17 (0%) 79 76	19, 38, 58, 111	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	295	ALA	4.1
1	E	295	ALA	3.8
1	D	245	PHE	3.7
1	E	293	LEU	3.2
1	B	227	TYR	3.1
1	A	245	PHE	2.9
1	B	295	ALA	2.7
1	F	227	TYR	2.6
1	D	77	ALA	2.6
1	E	228	ARG	2.4
1	E	1	MET	2.4
1	E	294	CYS	2.3
1	E	227	TYR	2.2
1	D	263	ASN	2.1
1	C	227	TYR	2.1
1	D	275	ILE	2.1
1	B	1	MET	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. 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	F	162	14/15	0.90	0.10	39,41,48,49	0
1	KPI	E	162	14/15	0.93	0.08	33,35,41,43	0
1	KPI	C	162	14/15	0.93	0.09	30,32,40,40	0
1	KPI	D	162	14/15	0.95	0.07	24,30,39,39	0
1	KPI	B	162	14/15	0.96	0.07	19,22,24,25	0
1	KPI	A	162	14/15	0.97	0.06	20,22,33,36	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.