



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 09:15 PM UTC

PDB ID : 4JX5 / pdb_00004jx5
Title : Structure of the carboxyl transferase domain from *Rhizobium etli* pyruvate carboxylase with pyruvate
Authors : Lietzan, A.D.; St Maurice, M.
Deposited on : 2013-03-27
Resolution : 2.55 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

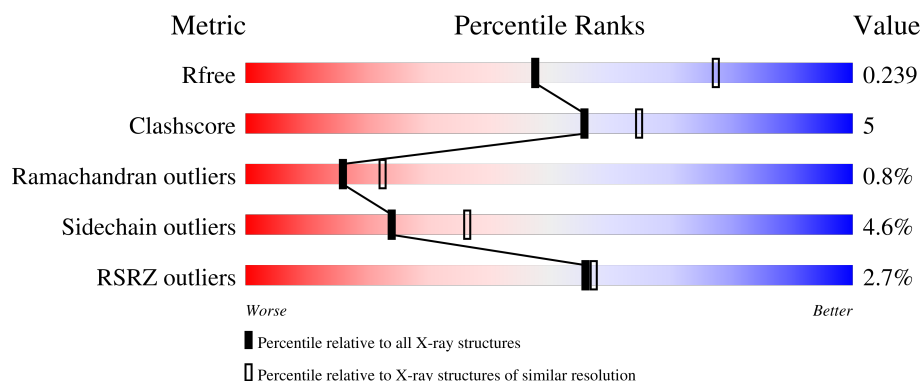
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.55 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	1091 (2.54-2.54)
Clashscore	190562	1120 (2.54-2.54)
Ramachandran outliers	187476	1106 (2.54-2.54)
Sidechain outliers	187428	1106 (2.54-2.54)
RSRZ outliers	180081	1091 (2.54-2.54)

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	632	82% 11% 6%
1	B	632	84% 10% 6%
1	C	632	5% 79% 13% 6%
1	D	632	3% 80% 12% 7%

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	CL	C	1102	-	-	X	-
6	GOL	B	1101	-	-	X	-

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 17941 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 Pyruvate carboxylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	597	Total	C	N	O	S	0	2	0
			4539	2889	758	869	23			
1	B	596	Total	C	N	O	S	0	1	0
			4468	2840	747	858	23			
1	C	591	Total	C	N	O	S	0	0	0
			4357	2758	739	838	22			
1	D	590	Total	C	N	O	S	0	0	0
			4322	2733	733	834	22			

There are 116 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	436	MET	-	expression tag	UNP Q2K340
A	437	GLY	-	expression tag	UNP Q2K340
A	438	SER	-	expression tag	UNP Q2K340
A	439	SER	-	expression tag	UNP Q2K340
A	440	HIS	-	expression tag	UNP Q2K340
A	441	HIS	-	expression tag	UNP Q2K340
A	442	HIS	-	expression tag	UNP Q2K340
A	443	HIS	-	expression tag	UNP Q2K340
A	444	HIS	-	expression tag	UNP Q2K340
A	445	HIS	-	expression tag	UNP Q2K340
A	446	HIS	-	expression tag	UNP Q2K340
A	447	HIS	-	expression tag	UNP Q2K340
A	448	ASP	-	expression tag	UNP Q2K340
A	449	TYR	-	expression tag	UNP Q2K340
A	450	ASP	-	expression tag	UNP Q2K340
A	451	ILE	-	expression tag	UNP Q2K340
A	452	PRO	-	expression tag	UNP Q2K340
A	453	THR	-	expression tag	UNP Q2K340
A	454	SER	-	expression tag	UNP Q2K340
A	455	GLU	-	expression tag	UNP Q2K340
A	456	ASN	-	expression tag	UNP Q2K340

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
A	457	LEU	-	expression tag	UNP Q2K340
A	458	TYR	-	expression tag	UNP Q2K340
A	459	PHE	-	expression tag	UNP Q2K340
A	460	GLN	-	expression tag	UNP Q2K340
A	461	GLY	-	expression tag	UNP Q2K340
A	462	LEU	-	expression tag	UNP Q2K340
A	463	LEU	-	expression tag	UNP Q2K340
A	464	HIS	-	expression tag	UNP Q2K340
B	436	MET	-	expression tag	UNP Q2K340
B	437	GLY	-	expression tag	UNP Q2K340
B	438	SER	-	expression tag	UNP Q2K340
B	439	SER	-	expression tag	UNP Q2K340
B	440	HIS	-	expression tag	UNP Q2K340
B	441	HIS	-	expression tag	UNP Q2K340
B	442	HIS	-	expression tag	UNP Q2K340
B	443	HIS	-	expression tag	UNP Q2K340
B	444	HIS	-	expression tag	UNP Q2K340
B	445	HIS	-	expression tag	UNP Q2K340
B	446	HIS	-	expression tag	UNP Q2K340
B	447	HIS	-	expression tag	UNP Q2K340
B	448	ASP	-	expression tag	UNP Q2K340
B	449	TYR	-	expression tag	UNP Q2K340
B	450	ASP	-	expression tag	UNP Q2K340
B	451	ILE	-	expression tag	UNP Q2K340
B	452	PRO	-	expression tag	UNP Q2K340
B	453	THR	-	expression tag	UNP Q2K340
B	454	SER	-	expression tag	UNP Q2K340
B	455	GLU	-	expression tag	UNP Q2K340
B	456	ASN	-	expression tag	UNP Q2K340
B	457	LEU	-	expression tag	UNP Q2K340
B	458	TYR	-	expression tag	UNP Q2K340
B	459	PHE	-	expression tag	UNP Q2K340
B	460	GLN	-	expression tag	UNP Q2K340
B	461	GLY	-	expression tag	UNP Q2K340
B	462	LEU	-	expression tag	UNP Q2K340
B	463	LEU	-	expression tag	UNP Q2K340
B	464	HIS	-	expression tag	UNP Q2K340
C	436	MET	-	expression tag	UNP Q2K340
C	437	GLY	-	expression tag	UNP Q2K340
C	438	SER	-	expression tag	UNP Q2K340
C	439	SER	-	expression tag	UNP Q2K340
C	440	HIS	-	expression tag	UNP Q2K340

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
C	441	HIS	-	expression tag	UNP Q2K340
C	442	HIS	-	expression tag	UNP Q2K340
C	443	HIS	-	expression tag	UNP Q2K340
C	444	HIS	-	expression tag	UNP Q2K340
C	445	HIS	-	expression tag	UNP Q2K340
C	446	HIS	-	expression tag	UNP Q2K340
C	447	HIS	-	expression tag	UNP Q2K340
C	448	ASP	-	expression tag	UNP Q2K340
C	449	TYR	-	expression tag	UNP Q2K340
C	450	ASP	-	expression tag	UNP Q2K340
C	451	ILE	-	expression tag	UNP Q2K340
C	452	PRO	-	expression tag	UNP Q2K340
C	453	THR	-	expression tag	UNP Q2K340
C	454	SER	-	expression tag	UNP Q2K340
C	455	GLU	-	expression tag	UNP Q2K340
C	456	ASN	-	expression tag	UNP Q2K340
C	457	LEU	-	expression tag	UNP Q2K340
C	458	TYR	-	expression tag	UNP Q2K340
C	459	PHE	-	expression tag	UNP Q2K340
C	460	GLN	-	expression tag	UNP Q2K340
C	461	GLY	-	expression tag	UNP Q2K340
C	462	LEU	-	expression tag	UNP Q2K340
C	463	LEU	-	expression tag	UNP Q2K340
C	464	HIS	-	expression tag	UNP Q2K340
D	436	MET	-	expression tag	UNP Q2K340
D	437	GLY	-	expression tag	UNP Q2K340
D	438	SER	-	expression tag	UNP Q2K340
D	439	SER	-	expression tag	UNP Q2K340
D	440	HIS	-	expression tag	UNP Q2K340
D	441	HIS	-	expression tag	UNP Q2K340
D	442	HIS	-	expression tag	UNP Q2K340
D	443	HIS	-	expression tag	UNP Q2K340
D	444	HIS	-	expression tag	UNP Q2K340
D	445	HIS	-	expression tag	UNP Q2K340
D	446	HIS	-	expression tag	UNP Q2K340
D	447	HIS	-	expression tag	UNP Q2K340
D	448	ASP	-	expression tag	UNP Q2K340
D	449	TYR	-	expression tag	UNP Q2K340
D	450	ASP	-	expression tag	UNP Q2K340
D	451	ILE	-	expression tag	UNP Q2K340
D	452	PRO	-	expression tag	UNP Q2K340
D	453	THR	-	expression tag	UNP Q2K340

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
D	454	SER	-	expression tag	UNP Q2K340
D	455	GLU	-	expression tag	UNP Q2K340
D	456	ASN	-	expression tag	UNP Q2K340
D	457	LEU	-	expression tag	UNP Q2K340
D	458	TYR	-	expression tag	UNP Q2K340
D	459	PHE	-	expression tag	UNP Q2K340
D	460	GLN	-	expression tag	UNP Q2K340
D	461	GLY	-	expression tag	UNP Q2K340
D	462	LEU	-	expression tag	UNP Q2K340
D	463	LEU	-	expression tag	UNP Q2K340
D	464	HIS	-	expression tag	UNP Q2K340

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Zn 1 1	0	0
2	B	1	Total Zn 1 1	0	0
2	C	1	Total Zn 1 1	0	0
2	D	1	Total Zn 1 1	0	0

- Molecule 3 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Cl 1 1	0	0
3	B	1	Total Cl 1 1	0	0
3	C	1	Total Cl 1 1	0	0
3	D	1	Total Cl 1 1	0	0

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

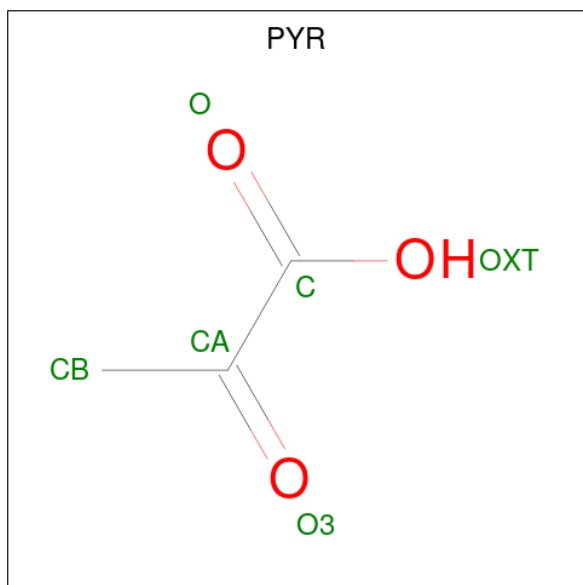
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total Mg 1 1	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Mg	0	0
			1	1		
4	C	1	Total	Mg	0	0
			1	1		
4	D	1	Total	Mg	0	0
			1	1		

- Molecule 5 is PYRUVIC ACID (CCD ID: PYR) (formula: $C_3H_4O_3$).



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

- Molecule 6 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

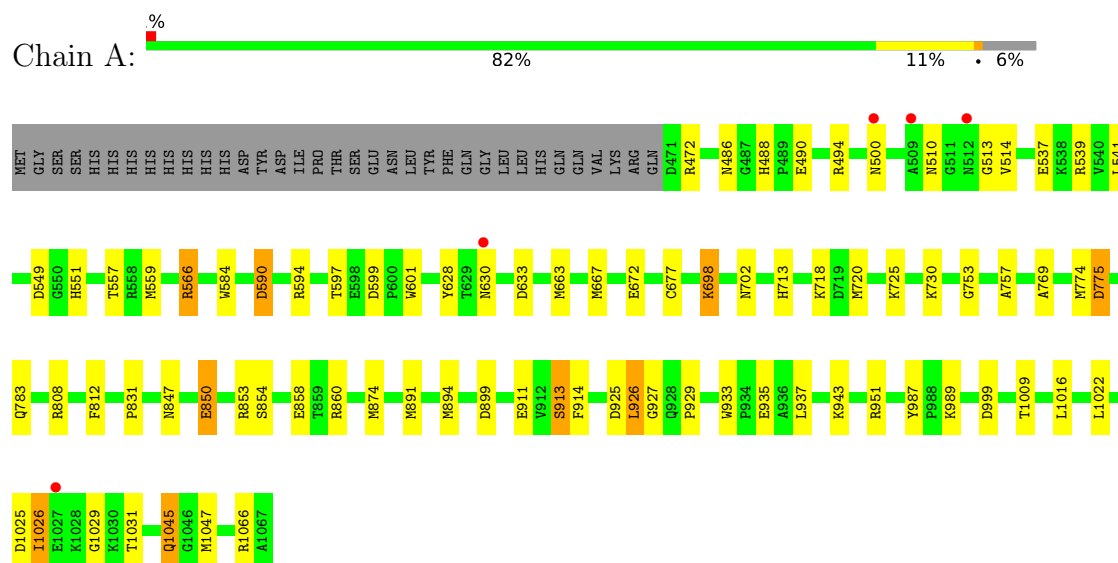
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	86	Total	O	0	0
			86	86		
7	B	46	Total	O	0	0
			46	46		
7	C	42	Total	O	0	0
			42	42		
7	D	33	Total	O	0	0
			33	33		

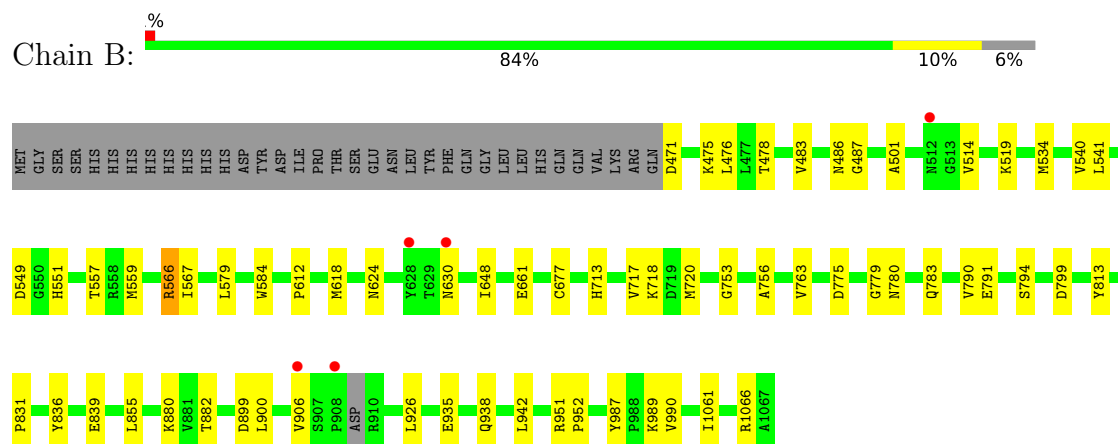
3 Residue-property plots [i](#)

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.

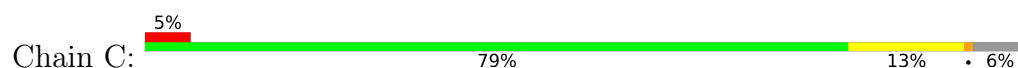
- Molecule 1: Pyruvate carboxylase

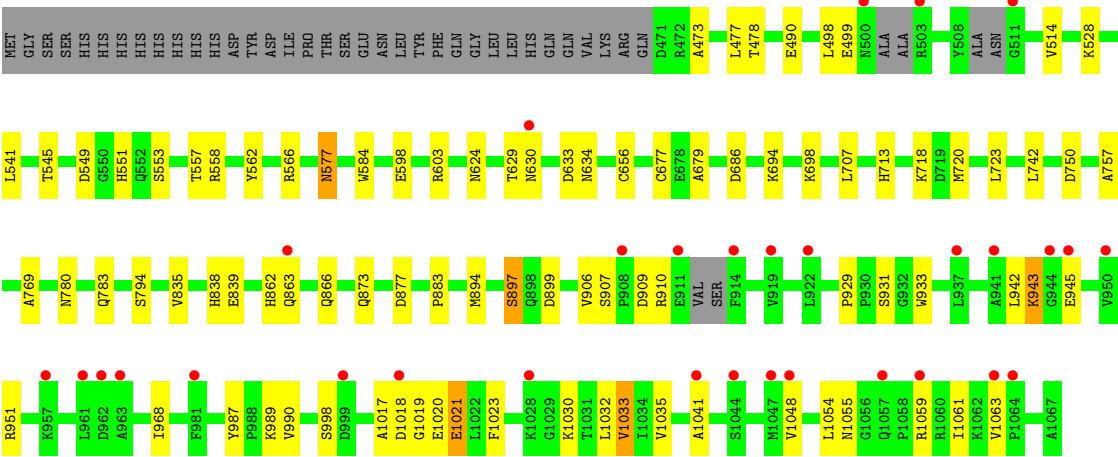


- Molecule 1: Pyruvate carboxylase

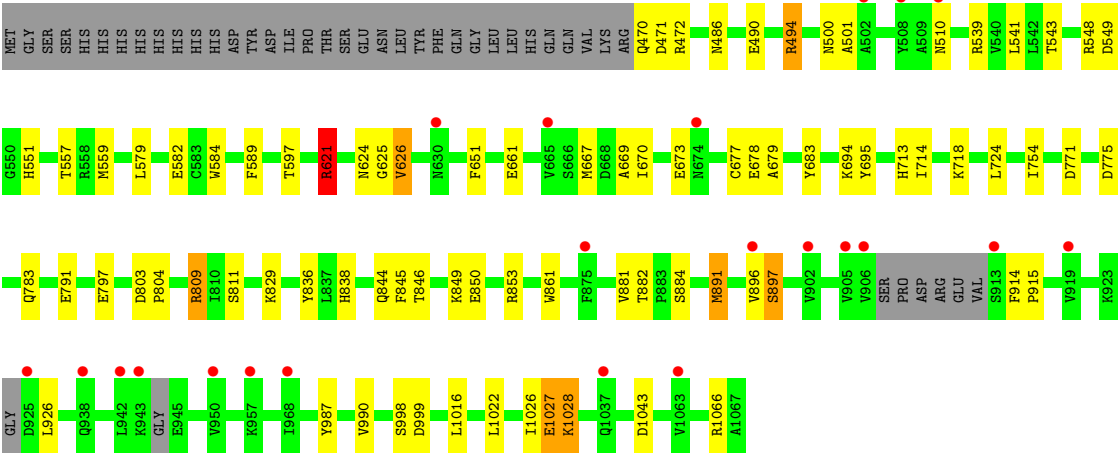
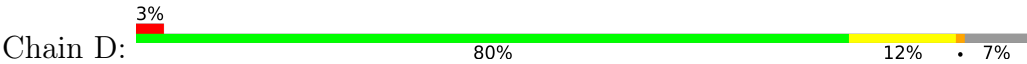


- Molecule 1: Pyruvate carboxylase





• Molecule 1: Pyruvate carboxylase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	85.13Å 157.25Å 245.32Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.88 – 2.55 46.88 – 2.55	Depositor EDS
% Data completeness (in resolution range)	99.5 (46.88-2.55) 99.7 (46.88-2.55)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.38 (at 2.54Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.187 , 0.232 (Not available) , 0.239	Depositor DCC
R_{free} test set	5406 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	55.9	Xtriage
Anisotropy	0.034	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 61.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	17941	wwPDB-VP
Average B, all atoms (Å ²)	71.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.22% 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, CL, MG, ZN, PYR, KCX

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	0.92	2/4630 (0.0%)	1.04	4/6301 (0.1%)
1	B	0.78	0/4554	0.99	0/6210
1	C	0.89	4/4438 (0.1%)	1.01	5/6056 (0.1%)
1	D	0.68	0/4402	0.95	1/6013 (0.0%)
All	All	0.82	6/18024 (0.0%)	1.00	10/24580 (0.0%)

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	1063	VAL	C-O	15.20	1.43	1.24
1	C	1048	VAL	C-O	11.63	1.35	1.24
1	C	910	ARG	CZ-NH1	6.90	1.42	1.32
1	C	1059	ARG	CZ-NH1	6.50	1.41	1.32
1	A	628	TYR	N-CA	5.43	1.53	1.45
1	A	858	GLU	C-O	-5.26	1.18	1.24

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	838	HIS	N-CA-C	5.88	117.77	111.36
1	A	1009	THR	CA-C-N	-5.85	112.55	119.05
1	A	1009	THR	C-N-CA	-5.85	112.55	119.05
1	C	1063	VAL	CA-C-O	5.57	127.42	119.95
1	A	935	GLU	N-CA-C	5.50	116.95	111.07
1	A	926	LEU	N-CA-C	-5.47	105.97	113.30
1	C	698	LYS	N-CA-C	-5.43	104.59	111.11
1	C	742	LEU	CA-C-N	-5.13	114.95	120.03
1	C	742	LEU	C-N-CA	-5.13	114.95	120.03
1	C	1030	LYS	N-CA-C	5.12	117.73	110.10

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4539	0	4406	36	0
1	B	4468	0	4278	34	0
1	C	4357	0	4067	46	0
1	D	4322	0	3994	51	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	2	0
3	D	1	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
5	A	6	0	0	0	0
5	B	6	0	0	0	0
5	C	6	0	0	0	0
5	D	6	0	0	1	0
6	B	6	0	8	5	0
6	C	6	0	8	0	0
7	A	86	0	0	4	0
7	B	46	0	0	0	0
7	C	42	0	0	1	0
7	D	33	0	0	1	0
All	All	17941	0	16761	162	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 5.

All (162) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:809:ARG:HG2	1:D:809:ARG:HH11	1.17	1.06
1:D:470:GLN:HG3	1:D:472:ARG:HB2	1.43	1.01
1:C:677:CYS:H	1:C:713:HIS:HD2	1.10	0.99
1:B:566:ARG:HH11	1:B:566:ARG:HG3	1.27	0.99
1:D:677:CYS:H	1:D:713:HIS:HD2	1.00	0.97
1:C:562:TYR:O	1:C:566:ARG:HG3	1.68	0.94
1:A:677:CYS:H	1:A:713:HIS:HD2	1.02	0.92
1:D:677:CYS:H	1:D:713:HIS:CD2	1.88	0.90
1:D:494:ARG:HH11	1:D:494:ARG:HG2	1.37	0.89
1:B:780:ASN:H	6:B:1101:GOL:H11	1.37	0.88
1:A:677:CYS:H	1:A:713:HIS:CD2	1.92	0.86
1:B:677:CYS:H	1:B:713:HIS:HD2	1.22	0.85
1:C:677:CYS:H	1:C:713:HIS:CD2	1.95	0.83
1:D:549:ASP:HB3	1:D:783:GLN:HE22	1.44	0.83
1:B:566:ARG:HH11	1:B:566:ARG:CG	1.97	0.76
1:C:490:GLU:OE1	1:C:558:ARG:NH2	2.21	0.73
1:D:677:CYS:N	1:D:713:HIS:HD2	1.82	0.72
1:D:470:GLN:CG	1:D:472:ARG:HB2	2.20	0.70
1:D:809:ARG:HG2	1:D:809:ARG:NH1	1.94	0.70
1:A:472:ARG:CB	1:A:1026:ILE:HD11	2.22	0.70
1:C:577:ASN:HB2	7:C:1208:HOH:O	1.92	0.70
1:C:990:VAL:N	3:C:1102:CL:CL	2.60	0.68
1:D:621:ARG:HB3	1:D:624:ASN:HB2	1.75	0.67
1:C:1021:GLU:HG2	1:C:1035:VAL:HG22	1.76	0.67
1:B:677:CYS:H	1:B:713:HIS:CD2	2.11	0.65
1:B:549:ASP:HB3	1:B:783:GLN:HE22	1.62	0.65
1:C:549:ASP:HB3	1:C:783:GLN:HE22	1.64	0.63
1:C:679:ALA:HB1	1:C:707:LEU:HD22	1.81	0.62
1:D:486:ASN:HD21	1:D:1066:ARG:H	1.47	0.61
1:A:730:LYS:HG2	7:A:1270:HOH:O	2.01	0.60
1:A:753:GLY:HA3	1:A:831:PRO:HB3	1.84	0.60
1:B:780:ASN:H	6:B:1101:GOL:C1	2.13	0.60
1:D:549:ASP:HB3	1:D:783:GLN:NE2	2.16	0.60
1:A:847:ASN:HB2	7:A:1242:HOH:O	2.03	0.59
1:A:677:CYS:N	1:A:713:HIS:HD2	1.87	0.58
1:A:850:GLU:OE2	1:A:853:ARG:NE	2.37	0.58
1:D:486:ASN:ND2	1:D:1066:ARG:H	2.02	0.58
1:B:566:ARG:CG	1:B:566:ARG:NH1	2.59	0.57
1:C:545:THR:OG1	1:C:783:GLN:NE2	2.38	0.57
1:D:589:PHE:CG	1:D:626:VAL:HG11	2.40	0.56
1:B:478:THR:HA	1:B:1061:ILE:HG21	1.87	0.56
1:C:677:CYS:N	1:C:713:HIS:HD2	1.93	0.56

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:850:GLU:O	1:D:853:ARG:HB3	2.06	0.56
1:D:539:ARG:NH1	7:D:1202:HOH:O	2.32	0.56
1:B:779:GLY:HA2	6:B:1101:GOL:H12	1.87	0.56
1:C:478:THR:HA	1:C:1061:ILE:HG21	1.88	0.55
1:D:667:MET:HE3	1:D:679:ALA:HB2	1.86	0.55
1:C:987:TYR:HB3	1:C:990:VAL:HB	1.89	0.55
1:A:549:ASP:HB3	1:A:783:GLN:HE22	1.70	0.55
1:D:494:ARG:HG2	1:D:494:ARG:NH1	2.10	0.55
1:B:519:LYS:HG2	1:B:612:PRO:O	2.07	0.55
1:B:486:ASN:ND2	1:B:1066:ARG:H	2.05	0.54
1:A:1045:GLN:HB3	1:A:1047:MET:HG2	1.88	0.54
1:A:1026:ILE:O	1:A:1026:ILE:HG12	2.08	0.54
1:C:624:ASN:HD21	1:C:630:ASN:ND2	2.06	0.54
1:C:1033:VAL:HG23	1:C:1055:ASN:OD1	2.07	0.54
1:A:488:HIS:CE1	1:A:490:GLU:HB2	2.42	0.53
1:A:894:MET:HE1	1:A:913:SER:O	2.08	0.53
1:B:791:GLU:HG2	1:D:836:TYR:CD2	2.44	0.53
1:D:683:TYR:CE1	1:D:724:LEU:HD13	2.43	0.53
1:D:694:LYS:HD3	1:D:695:TYR:CE2	2.44	0.53
1:C:694:LYS:NZ	1:C:931:SER:HB2	2.24	0.53
1:B:753:GLY:HA3	1:B:831:PRO:HB3	1.90	0.52
1:B:486:ASN:HD21	1:B:1066:ARG:H	1.55	0.52
1:D:621:ARG:HB2	1:D:625:GLY:O	2.09	0.52
1:D:651:PHE:HB2	1:D:670:ILE:HD13	1.91	0.52
1:B:618:MET:HB3	1:B:648:ILE:HD12	1.93	0.51
1:C:633:ASP:OD1	1:C:951:ARG:NH1	2.44	0.51
1:D:1026:ILE:HG23	1:D:1027:GLU:HG3	1.93	0.51
1:A:599:ASP:OD1	1:A:601:TRP:N	2.43	0.51
1:A:630:ASN:OD1	1:A:925:ASP:O	2.29	0.50
1:A:698:LYS:HG2	1:A:702:ASN:ND2	2.26	0.50
1:C:624:ASN:HD21	1:C:630:ASN:HD22	1.57	0.50
1:C:989:LYS:N	3:C:1102:CL:CL	2.81	0.50
1:A:590:ASP:HB2	1:A:987:TYR:CZ	2.47	0.49
1:B:753:GLY:HA2	6:B:1101:GOL:O1	2.11	0.49
1:A:590:ASP:OD1	1:A:594:ARG:NH2	2.45	0.49
1:A:1016:LEU:HD21	1:A:1022:LEU:HD22	1.94	0.49
1:A:537:GLU:HG3	1:A:539:ARG:HG2	1.94	0.49
1:A:730:LYS:CG	7:A:1270:HOH:O	2.60	0.49
1:D:548:ARG:HB3	1:D:582:GLU:OE1	2.13	0.49
1:D:809:ARG:HH11	1:D:809:ARG:CG	2.05	0.48
1:D:669:ALA:O	1:D:673:GLU:HG2	2.13	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:774:MET:O	1:A:775:ASP:C	2.57	0.48
1:C:598:GLU:OE2	1:C:603:ARG:NH2	2.43	0.48
1:C:1021:GLU:CG	1:C:1035:VAL:HG22	2.41	0.48
1:A:551:HIS:CE1	1:A:559:MET:HB3	2.49	0.48
1:C:473:ALA:HB1	1:C:1054:LEU:HD21	1.96	0.47
1:D:775:ASP:HB2	1:D:811:SER:OG	2.14	0.47
1:B:471:ASP:OD2	1:B:475:LYS:HE3	2.15	0.47
1:A:486:ASN:ND2	1:A:1066:ARG:H	2.12	0.47
1:B:541:LEU:HB3	1:B:579:LEU:HB2	1.95	0.47
1:C:897:SER:C	1:C:899:ASP:H	2.23	0.47
1:D:548:ARG:HE	1:D:582:GLU:CD	2.23	0.47
1:C:624:ASN:ND2	1:C:630:ASN:HD22	2.13	0.46
1:C:750:ASP:C	1:C:780:ASN:O	2.59	0.46
1:D:541:LEU:HB3	1:D:579:LEU:HB2	1.98	0.46
1:C:562:TYR:CE2	1:C:566:ARG:HD2	2.50	0.46
1:D:891:MET:HE3	1:D:891:MET:HA	1.98	0.46
1:C:942:LEU:O	1:C:943:LYS:C	2.59	0.45
1:D:543:THR:HB	1:D:771:ASP:OD1	2.17	0.45
1:A:927:GLY:HA3	7:A:1224:HOH:O	2.15	0.45
1:D:844:GLN:O	1:D:845:PHE:C	2.59	0.45
1:A:891[B]:MET:HE1	1:A:914:PHE:CE1	2.52	0.45
1:D:987:TYR:HB3	1:D:990:VAL:HB	1.98	0.45
1:B:836:TYR:CD2	1:D:791:GLU:HG2	2.51	0.45
1:C:473:ALA:O	1:C:477:LEU:HG	2.17	0.45
1:C:894:MET:HE3	1:C:894:MET:HB2	1.93	0.45
1:D:589:PHE:CD2	1:D:626:VAL:HG11	2.52	0.45
1:B:551:HIS:CE1	1:B:559:MET:HB3	2.51	0.45
1:B:624:ASN:ND2	1:B:630:ASN:OD1	2.50	0.45
1:D:551:HIS:CE1	1:D:559:MET:HB3	2.52	0.45
1:A:891[B]:MET:HE1	1:A:914:PHE:CD1	2.51	0.44
1:A:1025:ASP:OD2	1:A:1031:THR:OG1	2.25	0.44
1:C:624:ASN:ND2	1:C:630:ASN:ND2	2.65	0.44
1:B:677:CYS:N	1:B:713:HIS:HD2	2.02	0.44
1:D:803:ASP:HA	1:D:804:PRO:HD3	1.91	0.44
1:B:483:VAL:O	1:B:487:GLY:N	2.50	0.44
1:C:906:VAL:O	1:C:907:SER:C	2.60	0.44
1:D:849:LYS:HD2	1:D:861:TRP:CE2	2.53	0.44
1:D:1016:LEU:HD21	1:D:1022:LEU:HD22	1.98	0.44
1:C:723:LEU:HD11	1:C:839:GLU:HG2	1.99	0.44
1:C:551:HIS:ND1	1:C:557:THR:HA	2.32	0.43
1:C:541:LEU:O	1:C:769:ALA:HA	2.18	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1023:PHE:HA	1:C:1032:LEU:O	2.18	0.43
1:A:566:ARG:HH11	1:A:566:ARG:HG3	1.84	0.43
1:B:938:GLN:HG2	1:B:942:LEU:HD22	2.01	0.43
1:D:548:ARG:HD2	1:D:548:ARG:C	2.44	0.43
1:C:835:VAL:HA	1:C:838:HIS:CE1	2.54	0.43
1:D:621:ARG:HD2	5:D:1104:PYR:O	2.19	0.43
1:B:779:GLY:CA	6:B:1101:GOL:H12	2.48	0.43
1:D:914:PHE:HB3	1:D:915:PRO:HD2	2.01	0.43
1:C:1017:ALA:O	1:C:1020:GLU:HB3	2.19	0.43
1:C:838:HIS:O	1:C:839:GLU:HB2	2.20	0.42
1:C:968:ILE:HG13	1:C:968:ILE:O	2.19	0.42
1:C:929:PRO:HD3	1:C:933:TRP:CZ2	2.53	0.42
1:D:1027:GLU:HB2	1:D:1028:LYS:H	1.64	0.42
1:B:567:ILE:HD13	1:B:813:TYR:CG	2.54	0.42
1:C:634:ASN:OD1	1:C:634:ASN:N	2.51	0.42
1:A:633:ASP:OD1	1:A:951:ARG:NH1	2.49	0.42
1:A:663:MET:O	1:A:667:MET:HG3	2.20	0.42
1:A:929:PRO:HD3	1:A:933:TRP:CZ2	2.54	0.42
1:C:656:CYS:SG	1:C:883:PRO:HD2	2.59	0.42
1:B:799:ASP:OD1	1:B:799:ASP:C	2.63	0.41
1:A:874:MET:HG2	1:A:937:LEU:HD22	2.02	0.41
1:C:929:PRO:HD3	1:C:933:TRP:CE2	2.55	0.41
1:D:678:GLU:HA	1:D:714:ILE:O	2.20	0.41
1:D:1043:ASP:C	1:D:1043:ASP:OD1	2.63	0.41
1:B:567:ILE:HD13	1:B:813:TYR:CD2	2.55	0.41
1:A:541:LEU:O	1:A:769:ALA:HA	2.21	0.41
1:B:756:ALA:CB	1:D:754:ILE:HG22	2.51	0.41
1:A:812[B]:PHE:CE1	1:C:862:HIS:CD2	3.09	0.41
1:D:896:VAL:O	1:D:897:SER:C	2.64	0.41
1:A:757:ALA:HA	1:C:757:ALA:HB2	2.03	0.41
1:B:839:GLU:OE1	1:B:880:LYS:NZ	2.54	0.41
1:B:987:TYR:HB3	1:B:990:VAL:HB	2.02	0.41
1:D:891:MET:HA	1:D:891:MET:CE	2.50	0.41
1:D:494:ARG:NH1	1:D:494:ARG:CG	2.83	0.40
1:B:951:ARG:O	1:B:952:PRO:C	2.65	0.40
1:B:540:VAL:HG21	1:B:763:VAL:HG22	2.02	0.40
1:C:686:ASP:OD1	1:C:686:ASP:C	2.64	0.40
1:D:881:VAL:O	1:D:884:SER:N	2.55	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	596/632 (94%)	577 (97%)	16 (3%)	3 (0%)	24	34
1	B	592/632 (94%)	563 (95%)	26 (4%)	3 (0%)	24	34
1	C	582/632 (92%)	541 (93%)	33 (6%)	8 (1%)	9	11
1	D	581/632 (92%)	540 (93%)	36 (6%)	5 (1%)	14	20
All	All	2351/2528 (93%)	2221 (94%)	111 (5%)	19 (1%)	16	22

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	1041	ALA
1	D	500	ASN
1	D	510	ASN
1	D	1028	LYS
1	A	1029	GLY
1	B	501	ALA
1	B	935	GLU
1	C	943	LYS
1	C	1019	GLY
1	D	501	ALA
1	A	500	ASN
1	A	513	GLY
1	C	498	LEU
1	C	499	GLU
1	C	897	SER
1	D	621	ARG
1	C	877	ASP
1	C	945	GLU
1	B	906	VAL

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	462/519 (89%)	436 (94%)	26 (6%)	19	29
1	B	447/519 (86%)	429 (96%)	18 (4%)	28	42
1	C	422/519 (81%)	406 (96%)	16 (4%)	29	44
1	D	412/519 (79%)	392 (95%)	20 (5%)	22	34
All	All	1743/2076 (84%)	1663 (95%)	80 (5%)	24	36

All (80) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	494	ARG
1	A	510	ASN
1	A	514	VAL
1	A	557	THR
1	A	566	ARG
1	A	584	TRP
1	A	590	ASP
1	A	597	THR
1	A	672	GLU
1	A	698	LYS
1	A	720	MET
1	A	725	LYS
1	A	775	ASP
1	A	808	ARG
1	A	850	GLU
1	A	854	SER
1	A	860	ARG
1	A	899	ASP
1	A	911	GLU
1	A	913	SER
1	A	926	LEU
1	A	943	LYS
1	A	989	LYS
1	A	999	ASP

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	1026	ILE
1	A	1045	GLN
1	B	476	LEU
1	B	514	VAL
1	B	534	MET
1	B	557	THR
1	B	566	ARG
1	B	584	TRP
1	B	661	GLU
1	B	717	VAL
1	B	720	MET
1	B	775	ASP
1	B	790	VAL
1	B	794	SER
1	B	855	LEU
1	B	882	THR
1	B	899	ASP
1	B	900	LEU
1	B	926	LEU
1	B	989	LYS
1	C	514	VAL
1	C	528	LYS
1	C	553	SER
1	C	577	ASN
1	C	584	TRP
1	C	629	THR
1	C	720	MET
1	C	794	SER
1	C	863	GLN
1	C	866	GLN
1	C	873	GLN
1	C	909	ASP
1	C	998	SER
1	C	1018	ASP
1	C	1021	GLU
1	C	1033	VAL
1	D	471	ASP
1	D	490	GLU
1	D	494	ARG
1	D	557	THR
1	D	584	TRP
1	D	597	THR

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	D	621	ARG
1	D	626	VAL
1	D	661	GLU
1	D	797	GLU
1	D	809	ARG
1	D	829	LYS
1	D	846	THR
1	D	882	THR
1	D	891	MET
1	D	897	SER
1	D	926	LEU
1	D	998	SER
1	D	999	ASP
1	D	1027	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (33) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	486	ASN
1	A	577	ASN
1	A	630	ASN
1	A	642	GLN
1	A	713	HIS
1	A	783	GLN
1	A	820	GLN
1	A	863	GLN
1	A	866	GLN
1	B	486	ASN
1	B	577	ASN
1	B	624	ASN
1	B	713	HIS
1	B	783	GLN
1	B	820	GLN
1	B	1045	GLN
1	C	486	ASN
1	C	577	ASN
1	C	617	GLN
1	C	624	ASN
1	C	630	ASN
1	C	642	GLN
1	C	674	ASN
1	C	713	HIS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	C	783	GLN
1	C	1037	GLN
1	C	1057	GLN
1	D	486	ASN
1	D	577	ASN
1	D	624	ASN
1	D	713	HIS
1	D	783	GLN
1	D	873	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

4 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	KCX	C	718	2,1	10,11,12	0.75	1 (10%)	6,12,14	1.37	1 (16%)
1	KCX	D	718	2,1	10,11,12	0.52	0	6,12,14	2.87	2 (33%)
1	KCX	B	718	2,1	10,11,12	0.38	0	6,12,14	3.32	2 (33%)
1	KCX	A	718	2,1	10,11,12	1.01	0	6,12,14	2.32	2 (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	KCX	C	718	2,1	-	0/9/10/12	-
1	KCX	D	718	2,1	-	4/9/10/12	-

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	KCX	B	718	2,1	-	6/9/10/12	-
1	KCX	A	718	2,1	-	1/9/10/12	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	718	KCX	O-C	2.07	1.27	1.20

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	718	KCX	CE-NZ-CX	7.51	134.72	121.98
1	D	718	KCX	CE-NZ-CX	5.68	131.61	121.98
1	A	718	KCX	CE-NZ-CX	3.98	128.73	121.98
1	D	718	KCX	OQ1-CX-NZ	-3.60	119.45	124.92
1	A	718	KCX	OQ1-CX-NZ	3.41	130.10	124.92
1	C	718	KCX	CE-NZ-CX	2.24	125.79	121.98
1	B	718	KCX	OQ1-CX-NZ	2.17	128.22	124.92

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	718	KCX	O-C-CA-CB
1	B	718	KCX	N-CA-CB-CG
1	B	718	KCX	C-CA-CB-CG
1	B	718	KCX	O-C-CA-CB
1	B	718	KCX	CD-CE-NZ-CX
1	D	718	KCX	N-CA-CB-CG
1	D	718	KCX	C-CA-CB-CG
1	D	718	KCX	O-C-CA-CB
1	B	718	KCX	CA-CB-CG-CD
1	D	718	KCX	CA-CB-CG-CD
1	B	718	KCX	CG-CD-CE-NZ

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](#)

Of 18 ligands modelled in this entry, 12 are monoatomic - leaving 6 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	PYR	D	1104	-	5,5,5	1.48	2 (40%)	3,6,6	1.37	1 (33%)
6	GOL	C	1101	-	5,5,5	0.44	0	5,5,5	1.66	2 (40%)
5	PYR	A	1104	-	5,5,5	1.28	0	3,6,6	2.01	2 (66%)
6	GOL	B	1101	-	5,5,5	0.27	0	5,5,5	1.32	1 (20%)
5	PYR	B	1105	-	5,5,5	1.48	1 (20%)	3,6,6	1.94	1 (33%)
5	PYR	C	1105	-	5,5,5	1.55	1 (20%)	3,6,6	1.35	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
5	PYR	D	1104	-	-	1/4/4/4	-
6	GOL	C	1101	-	-	2/4/4/4	-
5	PYR	A	1104	-	-	3/4/4/4	-
6	GOL	B	1101	-	-	0/4/4/4	-
5	PYR	B	1105	-	-	2/4/4/4	-
5	PYR	C	1105	-	-	0/4/4/4	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	1105	PYR	CA-C	-2.77	1.45	1.54

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	C	1105	PYR	CA-C	-2.50	1.46	1.54
5	D	1104	PYR	CA-C	-2.13	1.47	1.54
5	D	1104	PYR	OXT-C	-2.12	1.24	1.30

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	1105	PYR	OXT-C-CA	3.27	122.64	113.59
5	A	1104	PYR	OXT-C-CA	2.55	120.65	113.59
5	A	1104	PYR	OXT-C-O	-2.26	118.52	123.90
6	C	1101	GOL	O1-C1-C2	-2.18	100.54	110.38
6	C	1101	GOL	C3-C2-C1	-2.13	104.00	111.80
6	B	1101	GOL	C3-C2-C1	-2.03	104.34	111.80
5	D	1104	PYR	OXT-C-CA	2.03	119.21	113.59

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	B	1105	PYR	OXT-C-CA-CB
5	D	1104	PYR	OXT-C-CA-CB
6	C	1101	GOL	O1-C1-C2-C3
5	A	1104	PYR	OXT-C-CA-O3
5	A	1104	PYR	OXT-C-CA-CB
6	C	1101	GOL	O1-C1-C2-O2
5	A	1104	PYR	O-C-CA-CB
5	B	1105	PYR	O-C-CA-CB

There are no ring outliers.

2 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	D	1104	PYR	1	0
6	B	1101	GOL	5	0

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	596/632 (94%)	-0.50	5 (0%)	82 84	22, 48, 72, 130	21 (3%)
1	B	595/632 (94%)	-0.16	5 (0%)	82 84	42, 62, 106, 149	18 (3%)
1	C	590/632 (93%)	0.20	31 (5%)	32 31	30, 81, 162, 211	14 (2%)
1	D	589/632 (93%)	0.20	22 (3%)	45 46	47, 81, 122, 140	14 (2%)
All	All	2370/2528 (93%)	-0.07	63 (2%)	56 57	22, 65, 130, 211	67 (2%)

All (63) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	630	ASN	5.0
1	D	913	SER	4.3
1	B	630	ASN	4.1
1	D	925	ASP	4.1
1	D	950	VAL	3.3
1	C	944	GLY	3.3
1	C	500	ASN	3.2
1	C	1063	VAL	3.2
1	D	875	PHE	3.1
1	C	962	ASP	3.1
1	C	1018	ASP	3.1
1	C	1048	VAL	3.0
1	C	999	ASP	3.0
1	B	908	PRO	3.0
1	D	938	GLN	2.8
1	D	510	ASN	2.8
1	A	500	ASN	2.8
1	C	914	PHE	2.7
1	B	906	VAL	2.7
1	D	968	ILE	2.6
1	D	902	VAL	2.6

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	D	905	VAL	2.6
1	C	630	ASN	2.5
1	D	508	TYR	2.5
1	A	630	ASN	2.5
1	C	919	VAL	2.5
1	A	509	ALA	2.5
1	A	1027	GLU	2.5
1	C	1059	ARG	2.4
1	B	512	ASN	2.4
1	C	950	VAL	2.4
1	D	906	VAL	2.4
1	C	1044	SER	2.4
1	B	628	TYR	2.4
1	C	511	GLY	2.4
1	C	941	ALA	2.4
1	C	945	GLU	2.3
1	C	963	ALA	2.3
1	C	911	GLU	2.3
1	D	502	ALA	2.3
1	C	1057	GLN	2.3
1	A	512	ASN	2.3
1	C	503	ARG	2.3
1	C	961	LEU	2.3
1	D	665	VAL	2.3
1	D	943	LYS	2.2
1	D	919	VAL	2.2
1	C	1041	ALA	2.2
1	C	937	LEU	2.2
1	C	908	PRO	2.2
1	D	957	LYS	2.2
1	C	1028	LYS	2.2
1	C	1064	PRO	2.2
1	C	981	PHE	2.2
1	D	1063	VAL	2.1
1	C	863	GLN	2.1
1	D	942	LEU	2.1
1	C	1047	MET	2.1
1	D	1037	GLN	2.1
1	C	922	LEU	2.1
1	D	674	ASN	2.0
1	D	896	VAL	2.0
1	C	957	LYS	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(Å ²)	Q<0.9
1	KCX	D	718	12/13	0.94	0.11	62,66,73,74	0
1	KCX	B	718	12/13	0.97	0.08	55,57,60,61	0
1	KCX	C	718	12/13	0.97	0.07	54,59,66,70	0
1	KCX	A	718	12/13	0.97	0.07	37,40,46,47	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(Å ²)	Q<0.9
5	PYR	D	1104	6/6	0.84	0.21	46,51,60,60	3
5	PYR	C	1105	6/6	0.91	0.24	30,32,35,38	6
4	MG	B	1103	1/1	0.93	0.08	52,52,52,52	0
3	CL	B	1102	1/1	0.94	0.07	66,66,66,66	0
3	CL	D	1101	1/1	0.95	0.08	79,79,79,79	0
4	MG	C	1103	1/1	0.95	0.10	41,41,41,41	0
5	PYR	B	1105	6/6	0.96	0.11	34,41,44,48	3
3	CL	C	1102	1/1	0.96	0.07	76,76,76,76	0
4	MG	D	1102	1/1	0.96	0.17	57,57,57,57	0
6	GOL	B	1101	6/6	0.96	0.10	49,56,59,64	0
6	GOL	C	1101	6/6	0.96	0.08	45,47,51,51	0
5	PYR	A	1104	6/6	0.97	0.07	44,51,52,54	0
3	CL	A	1102	1/1	0.98	0.05	53,53,53,53	0
2	ZN	D	1103	1/1	0.98	0.03	60,60,60,60	0
4	MG	A	1103	1/1	0.98	0.15	45,45,45,45	0
2	ZN	C	1104	1/1	0.99	0.04	54,54,54,54	0
2	ZN	A	1101	1/1	1.00	0.04	43,43,43,43	0
2	ZN	B	1104	1/1	1.00	0.04	51,51,51,51	0

6.5 Other polymers [i](#)

There are no such residues in this entry.