



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 10:07 AM GMT

PDB ID : 3KX6
Title : Crystal structure of fructose-1,6-bisphosphate aldolase from Babesia bovis at 2.1Å resolution
Authors : Seattle Structural Genomics Center for Infectious Disease (SSGCID)
Deposited on : 2009-12-02
Resolution : 2.10 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 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
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) : trunk26865

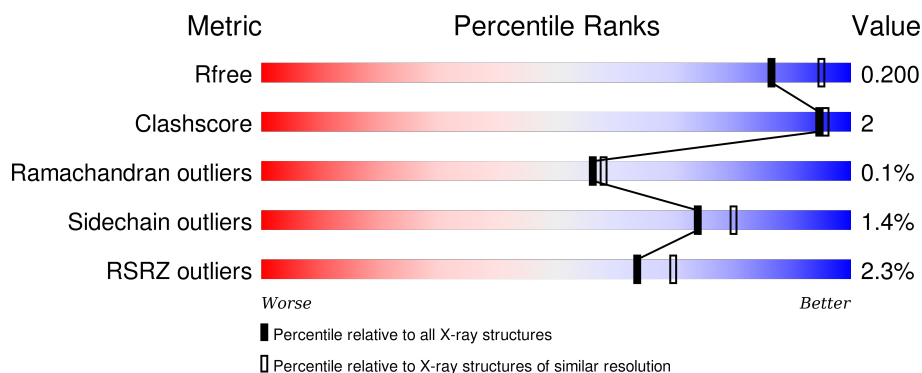
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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}	91344	3939 (2.10-2.10)
Clashscore	102246	4460 (2.10-2.10)
Ramachandran outliers	100387	4413 (2.10-2.10)
Sidechain outliers	100360	4414 (2.10-2.10)
RSRZ outliers	91569	3948 (2.10-2.10)

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	379	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 0%, red 1%, orange 1%, orange 88%, yellow 88%, yellow 9%, grey 9%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 88% 9% </div> </div>
1	B	379	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 0%, red 1%, orange 1%, orange 86%, yellow 86%, yellow 9%, grey 9%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 86% 9% </div> </div>
1	C	379	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 0%, red 3%, orange 3%, orange 87%, yellow 87%, yellow 9%, grey 9%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 3% 87% 9% </div> </div>
1	D	379	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 0%, red 3%, orange 3%, orange 87%, yellow 87%, yellow 9%, grey 9%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 3% 87% 9% </div> </div>

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	CIT	A	400	-	-	-	X
2	CIT	B	400	-	-	-	X
3	PEG	A	401	-	-	-	X
3	PEG	B	401	-	-	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 11265 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 Fructose-bisphosphate aldolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	344	Total	C	N	O	S	0	5	0
			2603	1634	450	507	12			
1	B	343	Total	C	N	O	S	0	4	0
			2607	1634	454	507	12			
1	C	343	Total	C	N	O	S	0	3	0
			2574	1614	445	503	12			
1	D	343	Total	C	N	O	S	0	4	0
			2584	1622	448	502	12			

There are 84 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-20	MET	-	EXPRESSION TAG	UNP A7AV51
A	-19	ALA	-	EXPRESSION TAG	UNP A7AV51
A	-18	HIS	-	EXPRESSION TAG	UNP A7AV51
A	-17	HIS	-	EXPRESSION TAG	UNP A7AV51
A	-16	HIS	-	EXPRESSION TAG	UNP A7AV51
A	-15	HIS	-	EXPRESSION TAG	UNP A7AV51
A	-14	HIS	-	EXPRESSION TAG	UNP A7AV51
A	-13	HIS	-	EXPRESSION TAG	UNP A7AV51
A	-12	MET	-	EXPRESSION TAG	UNP A7AV51
A	-11	GLY	-	EXPRESSION TAG	UNP A7AV51
A	-10	THR	-	EXPRESSION TAG	UNP A7AV51
A	-9	LEU	-	EXPRESSION TAG	UNP A7AV51
A	-8	GLU	-	EXPRESSION TAG	UNP A7AV51
A	-7	ALA	-	EXPRESSION TAG	UNP A7AV51
A	-6	GLN	-	EXPRESSION TAG	UNP A7AV51
A	-5	THR	-	EXPRESSION TAG	UNP A7AV51
A	-4	GLN	-	EXPRESSION TAG	UNP A7AV51
A	-3	GLY	-	EXPRESSION TAG	UNP A7AV51
A	-2	PRO	-	EXPRESSION TAG	UNP A7AV51
A	-1	GLY	-	EXPRESSION TAG	UNP A7AV51
A	0	SER	-	EXPRESSION TAG	UNP A7AV51

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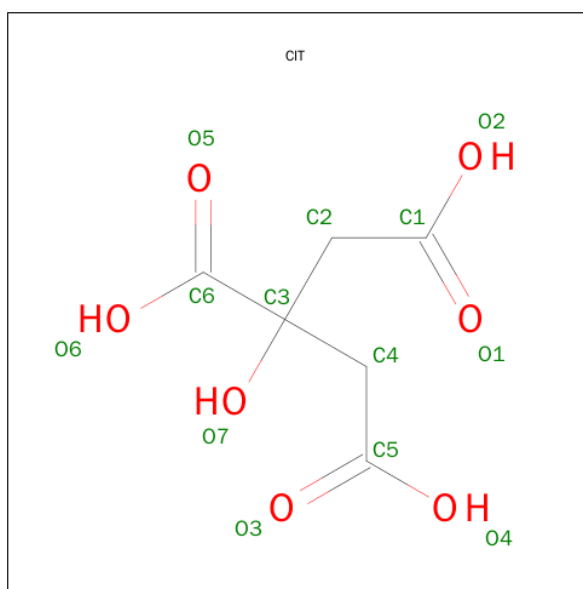
Chain	Residue	Modelled	Actual	Comment	Reference
B	-20	MET	-	EXPRESSION TAG	UNP A7AV51
B	-19	ALA	-	EXPRESSION TAG	UNP A7AV51
B	-18	HIS	-	EXPRESSION TAG	UNP A7AV51
B	-17	HIS	-	EXPRESSION TAG	UNP A7AV51
B	-16	HIS	-	EXPRESSION TAG	UNP A7AV51
B	-15	HIS	-	EXPRESSION TAG	UNP A7AV51
B	-14	HIS	-	EXPRESSION TAG	UNP A7AV51
B	-13	HIS	-	EXPRESSION TAG	UNP A7AV51
B	-12	MET	-	EXPRESSION TAG	UNP A7AV51
B	-11	GLY	-	EXPRESSION TAG	UNP A7AV51
B	-10	THR	-	EXPRESSION TAG	UNP A7AV51
B	-9	LEU	-	EXPRESSION TAG	UNP A7AV51
B	-8	GLU	-	EXPRESSION TAG	UNP A7AV51
B	-7	ALA	-	EXPRESSION TAG	UNP A7AV51
B	-6	GLN	-	EXPRESSION TAG	UNP A7AV51
B	-5	THR	-	EXPRESSION TAG	UNP A7AV51
B	-4	GLN	-	EXPRESSION TAG	UNP A7AV51
B	-3	GLY	-	EXPRESSION TAG	UNP A7AV51
B	-2	PRO	-	EXPRESSION TAG	UNP A7AV51
B	-1	GLY	-	EXPRESSION TAG	UNP A7AV51
B	0	SER	-	EXPRESSION TAG	UNP A7AV51
C	-20	MET	-	EXPRESSION TAG	UNP A7AV51
C	-19	ALA	-	EXPRESSION TAG	UNP A7AV51
C	-18	HIS	-	EXPRESSION TAG	UNP A7AV51
C	-17	HIS	-	EXPRESSION TAG	UNP A7AV51
C	-16	HIS	-	EXPRESSION TAG	UNP A7AV51
C	-15	HIS	-	EXPRESSION TAG	UNP A7AV51
C	-14	HIS	-	EXPRESSION TAG	UNP A7AV51
C	-13	HIS	-	EXPRESSION TAG	UNP A7AV51
C	-12	MET	-	EXPRESSION TAG	UNP A7AV51
C	-11	GLY	-	EXPRESSION TAG	UNP A7AV51
C	-10	THR	-	EXPRESSION TAG	UNP A7AV51
C	-9	LEU	-	EXPRESSION TAG	UNP A7AV51
C	-8	GLU	-	EXPRESSION TAG	UNP A7AV51
C	-7	ALA	-	EXPRESSION TAG	UNP A7AV51
C	-6	GLN	-	EXPRESSION TAG	UNP A7AV51
C	-5	THR	-	EXPRESSION TAG	UNP A7AV51
C	-4	GLN	-	EXPRESSION TAG	UNP A7AV51
C	-3	GLY	-	EXPRESSION TAG	UNP A7AV51
C	-2	PRO	-	EXPRESSION TAG	UNP A7AV51
C	-1	GLY	-	EXPRESSION TAG	UNP A7AV51
C	0	SER	-	EXPRESSION TAG	UNP A7AV51

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-20	MET	-	EXPRESSION TAG	UNP A7AV51
D	-19	ALA	-	EXPRESSION TAG	UNP A7AV51
D	-18	HIS	-	EXPRESSION TAG	UNP A7AV51
D	-17	HIS	-	EXPRESSION TAG	UNP A7AV51
D	-16	HIS	-	EXPRESSION TAG	UNP A7AV51
D	-15	HIS	-	EXPRESSION TAG	UNP A7AV51
D	-14	HIS	-	EXPRESSION TAG	UNP A7AV51
D	-13	HIS	-	EXPRESSION TAG	UNP A7AV51
D	-12	MET	-	EXPRESSION TAG	UNP A7AV51
D	-11	GLY	-	EXPRESSION TAG	UNP A7AV51
D	-10	THR	-	EXPRESSION TAG	UNP A7AV51
D	-9	LEU	-	EXPRESSION TAG	UNP A7AV51
D	-8	GLU	-	EXPRESSION TAG	UNP A7AV51
D	-7	ALA	-	EXPRESSION TAG	UNP A7AV51
D	-6	GLN	-	EXPRESSION TAG	UNP A7AV51
D	-5	THR	-	EXPRESSION TAG	UNP A7AV51
D	-4	GLN	-	EXPRESSION TAG	UNP A7AV51
D	-3	GLY	-	EXPRESSION TAG	UNP A7AV51
D	-2	PRO	-	EXPRESSION TAG	UNP A7AV51
D	-1	GLY	-	EXPRESSION TAG	UNP A7AV51
D	0	SER	-	EXPRESSION TAG	UNP A7AV51

- Molecule 2 is CITRIC ACID (three-letter code: CIT) (formula: C₆H₈O₇).



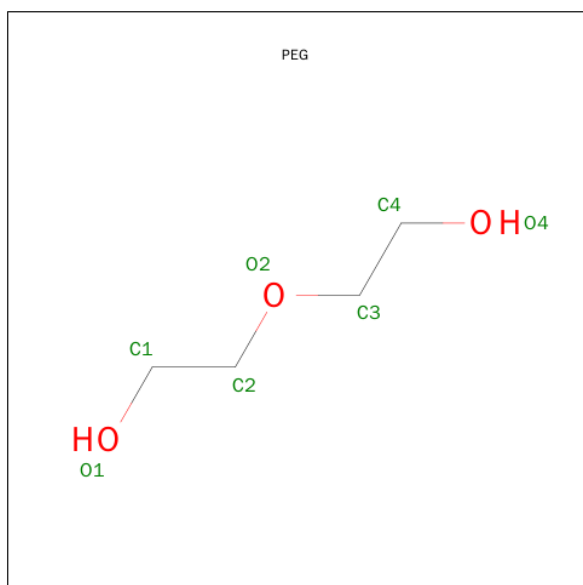
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			13	6	7		

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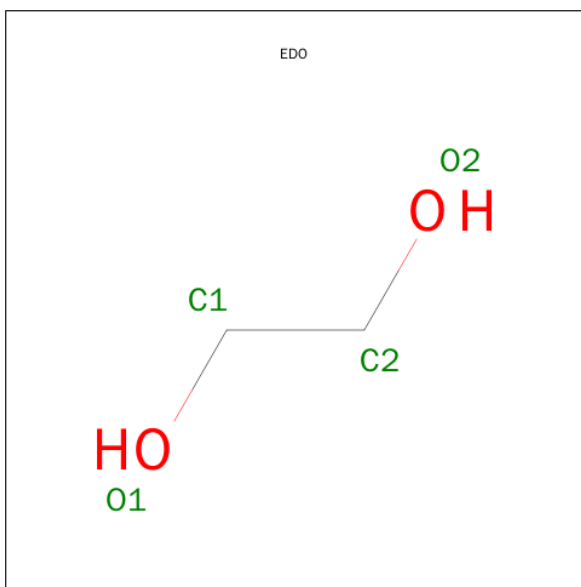
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	C	O	0	0
			13	6	7		
2	C	1	Total	C	O	0	0
			13	6	7		

- Molecule 3 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: C₄H₁₀O₃).



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

- Molecule 4 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



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

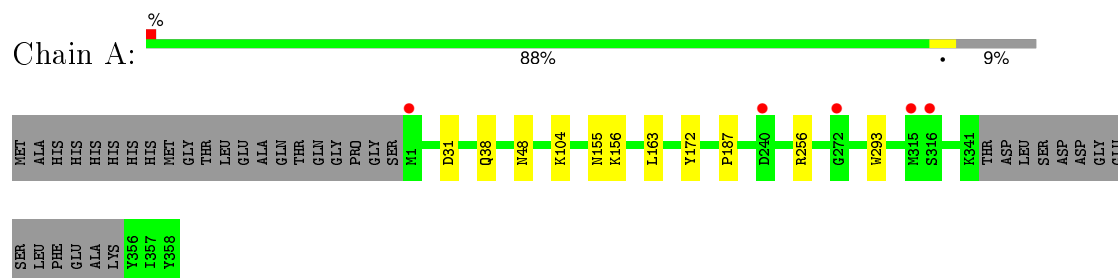
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	205	Total	O	0	0
			205	205		
5	B	226	Total	O	0	0
			226	226		
5	C	182	Total	O	0	0
			182	182		
5	D	192	Total	O	0	0
			192	192		

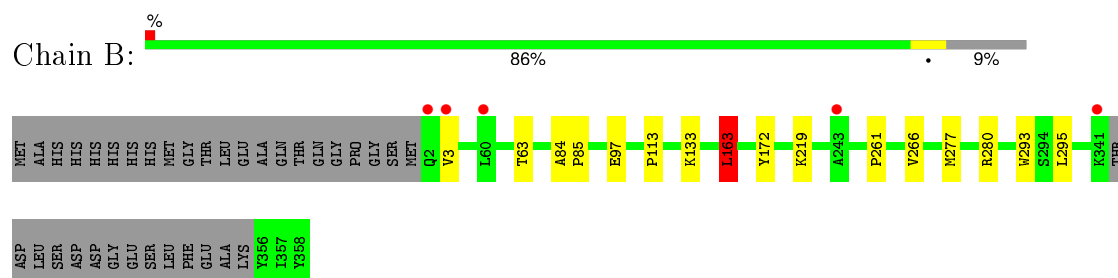
3 Residue-property plots [i](#)

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

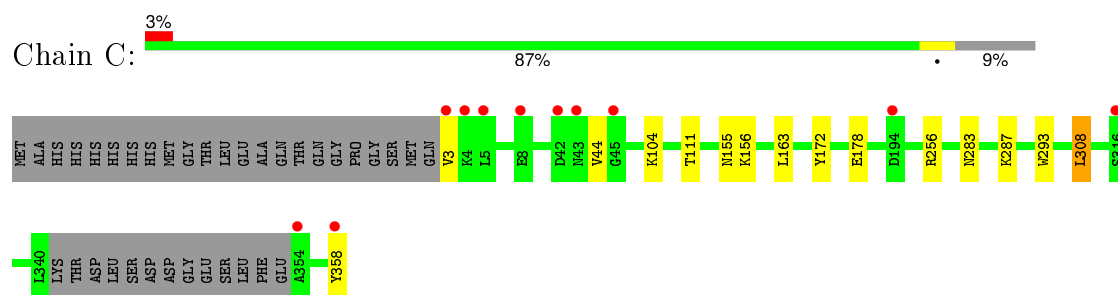
- Molecule 1: Fructose-bisphosphate aldolase



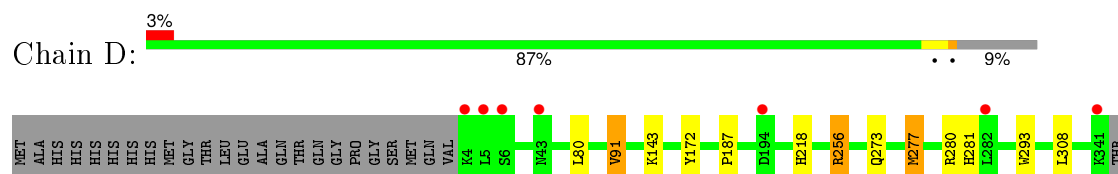
- Molecule 1: Fructose-bisphosphate aldolase



- Molecule 1: Fructose-bisphosphate aldolase



- Molecule 1: Fructose-bisphosphate aldolase



ASP		
LEU		
SER		
ASP		
ASP		
GLY		
GLU		
SER		
LEU		
PHE		
GLU		
A354		
Y357		
Y358		

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	85.88Å 114.16Å 174.54Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.10 19.96 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.4 (20.00-2.10) 99.4 (19.96-2.10)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.07 (at 2.09Å)	Xtriage
Refinement program	REFMAC 5.5.0104	Depositor
R, R_{free}	0.160 , 0.193 0.166 , 0.200	Depositor DCC
R_{free} test set	4992 reflections (5.26%)	DCC
Wilson B-factor (Å ²)	26.3	Xtriage
Anisotropy	0.644	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 59.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 99911 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	11265	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.42% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: PEG, EDO, CIT

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	0.73	0/2659	0.70	1/3608 (0.0%)
1	B	0.76	0/2660	0.69	1/3606 (0.0%)
1	C	0.74	1/2624 (0.0%)	0.74	1/3562 (0.0%)
1	D	0.72	0/2637	0.70	2/3577 (0.1%)
All	All	0.74	1/10580 (0.0%)	0.71	5/14353 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	178	GLU	CG-CD	5.73	1.60	1.51

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	163	LEU	CA-CB-CG	9.87	138.00	115.30
1	C	163	LEU	CA-CB-CG	9.65	137.50	115.30
1	B	163	LEU	CA-CB-CG	5.87	128.80	115.30
1	D	256[A]	ARG	NE-CZ-NH2	-5.17	117.72	120.30
1	D	256[B]	ARG	NE-CZ-NH2	-5.17	117.72	120.30

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	2603	0	2579	10	0
1	B	2607	0	2590	14	0
1	C	2574	0	2530	8	0
1	D	2584	0	2560	6	0
2	A	13	0	5	1	0
2	B	13	0	5	0	0
2	C	13	0	5	1	0
3	A	7	0	10	0	0
3	B	14	0	20	0	0
3	C	14	0	20	0	0
3	D	14	0	20	0	0
4	D	4	0	6	0	0
5	A	205	0	0	5	0
5	B	226	0	0	3	0
5	C	182	0	0	4	0
5	D	192	0	0	2	0
All	All	11265	0	10350	35	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 2.

All (35) 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:277:MET:HE3	1:D:280:ARG:HE	1.47	0.79
1:B:277:MET:CE	1:B:280:ARG:HE	2.10	0.65
1:A:256[A]:ARG:HD3	5:A:639:HOH:O	1.98	0.64
1:B:133:LYS:NZ	5:B:647:HOH:O	2.31	0.63
1:C:44:VAL:HG21	1:C:308:LEU:HD13	1.80	0.63
1:D:277:MET:CE	1:D:280:ARG:HE	2.17	0.58
1:C:283:ASN:O	1:C:287:LYS:HG3	2.06	0.56
1:B:277:MET:HE2	1:B:280:ARG:HE	1.73	0.52
1:D:218:HIS:HD2	5:D:373:HOH:O	1.95	0.49
1:C:358:TYR:OH	5:C:899:HOH:O	2.06	0.48
1:A:104:LYS:NZ	5:A:415:HOH:O	2.45	0.48
1:D:80:LEU:HD12	1:D:91[B]:VAL:CG1	2.44	0.47
1:B:277:MET:HE1	1:B:280:ARG:HE	1.78	0.47
1:C:256[A]:ARG:HD3	5:C:612:HOH:O	2.15	0.47
1:A:156:LYS:NZ	5:A:500:HOH:O	2.36	0.46
1:A:104:LYS:HE2	5:A:395:HOH:O	2.13	0.46
1:D:273:GLN:NE2	1:D:281:HIS:ND1	2.64	0.46
1:B:277:MET:CE	1:B:280:ARG:HH11	2.28	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:84:ALA:HB1	1:B:85:PRO:HD2	1.98	0.45
1:C:156:LYS:NZ	5:C:386:HOH:O	2.45	0.45
2:C:400:CIT:H41	5:C:836:HOH:O	2.15	0.45
1:A:31:ASP:OD2	1:A:104:LYS:HE3	2.17	0.44
1:B:63:THR:O	1:B:97:GLU:OE2	2.35	0.44
1:A:155[B]:ASN:ND2	1:C:155:ASN:OD1	2.51	0.44
1:D:256[A]:ARG:HD3	5:D:658:HOH:O	2.18	0.44
1:B:113:PRO:HA	1:C:3:VAL:HG13	2.00	0.43
1:A:104:LYS:HD3	5:A:792:HOH:O	2.18	0.43
1:B:219:LYS:NZ	5:B:814:HOH:O	2.52	0.43
1:A:38:GLN:NE2	1:A:48:ASN:HD22	2.17	0.43
1:B:3:VAL:HG21	1:C:111[A]:THR:HG21	2.01	0.42
1:B:163:LEU:HD13	5:B:535:HOH:O	2.18	0.42
1:B:266:VAL:HB	1:B:295:LEU:HD23	2.00	0.42
1:A:256[A]:ARG:HA	1:B:261:PRO:HG2	2.03	0.41
1:A:256[B]:ARG:HA	1:B:261:PRO:HG2	2.03	0.40
2:A:400:CIT:C6	2:A:400:CIT:O2	2.69	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	345/379 (91%)	340 (99%)	4 (1%)	1 (0%)	46	45
1	B	343/379 (90%)	337 (98%)	6 (2%)	0	100	100
1	C	342/379 (90%)	336 (98%)	6 (2%)	0	100	100
1	D	343/379 (90%)	338 (98%)	4 (1%)	1 (0%)	46	45
All	All	1373/1516 (91%)	1351 (98%)	20 (2%)	2 (0%)	56	58

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	187	PRO
1	D	187	PRO

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	276/312 (88%)	274 (99%)	2 (1%)	88	92
1	B	278/312 (89%)	275 (99%)	3 (1%)	80	85
1	C	270/312 (86%)	266 (98%)	4 (2%)	72	78
1	D	273/312 (88%)	266 (97%)	7 (3%)	54	58
All	All	1097/1248 (88%)	1081 (98%)	16 (2%)	74	78

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	172	TYR
1	A	293	TRP
1	B	163	LEU
1	B	172	TYR
1	B	293	TRP
1	C	104	LYS
1	C	172	TYR
1	C	293	TRP
1	C	308	LEU
1	D	91[A]	VAL
1	D	91[B]	VAL
1	D	143	LYS
1	D	172	TYR
1	D	277	MET
1	D	293	TRP
1	D	308	LEU

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

Mol	Chain	Res	Type
1	A	16	ASN
1	A	38	GLN
1	A	82	GLN
1	A	230	ASN
1	A	312	ASN
1	A	317	ASN
1	B	230	ASN
1	C	38	GLN
1	C	155	ASN
1	C	331	GLN
1	D	218	HIS
1	D	230	ASN
1	D	273	GLN
1	D	330	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

11 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	CIT	A	400	-	3,12,12	1.26	0	3,17,17	0.34	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	PEG	A	401	-	6,6,6	0.50	0	5,5,5	0.60	0
2	CIT	B	400	-	3,12,12	1.12	0	3,17,17	2.03	1 (33%)
3	PEG	B	401	-	6,6,6	0.49	0	5,5,5	0.70	0
3	PEG	B	402	-	6,6,6	0.58	0	5,5,5	0.40	0
2	CIT	C	400	-	3,12,12	1.31	1 (33%)	3,17,17	0.39	0
3	PEG	C	401	-	6,6,6	0.37	0	5,5,5	0.33	0
3	PEG	C	402	-	6,6,6	0.53	0	5,5,5	0.36	0
4	EDO	D	400	-	3,3,3	0.69	0	2,2,2	0.12	0
3	PEG	D	401	-	6,6,6	0.35	0	5,5,5	0.51	0
3	PEG	D	402	-	6,6,6	0.51	0	5,5,5	0.31	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	CIT	A	400	-	-	0/6/16/16	0/0/0/0
3	PEG	A	401	-	-	0/4/4/4	0/0/0/0
2	CIT	B	400	-	-	0/6/16/16	0/0/0/0
3	PEG	B	401	-	-	0/4/4/4	0/0/0/0
3	PEG	B	402	-	-	0/4/4/4	0/0/0/0
2	CIT	C	400	-	-	0/6/16/16	0/0/0/0
3	PEG	C	401	-	-	0/4/4/4	0/0/0/0
3	PEG	C	402	-	-	0/4/4/4	0/0/0/0
4	EDO	D	400	-	-	0/1/1/1	0/0/0/0
3	PEG	D	401	-	-	0/4/4/4	0/0/0/0
3	PEG	D	402	-	-	0/4/4/4	0/0/0/0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	400	CIT	O7-C3	2.02	1.46	1.43

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	400	CIT	C3-C4-C5	-3.25	109.77	114.96

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	400	CIT	1	0
2	C	400	CIT	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	344/379 (90%)	-0.29	5 (1%) 76 81	11, 19, 33, 46	0
1	B	343/379 (90%)	-0.35	5 (1%) 76 81	11, 18, 31, 38	0
1	C	343/379 (90%)	-0.28	11 (3%) 51 60	12, 19, 33, 42	0
1	D	343/379 (90%)	-0.22	10 (2%) 55 63	12, 19, 32, 47	0
All	All	1373/1516 (90%)	-0.28	31 (2%) 64 70	11, 19, 33, 47	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1	MET	6.3
1	C	3	VAL	6.1
1	D	358	TYR	5.8
1	D	5	LEU	5.1
1	D	354	ALA	4.0
1	C	45	GLY	3.9
1	C	354	ALA	3.4
1	B	341	LYS	3.3
1	C	5	LEU	3.3
1	C	316	SER	3.3
1	D	4	LYS	3.2
1	C	358	TYR	3.2
1	C	43	ASN	3.0
1	B	2	GLN	2.8
1	A	240	ASP	2.7
1	D	43	ASN	2.6
1	D	357	ILE	2.6
1	A	316	SER	2.6
1	A	272	GLY	2.6
1	C	8	GLU	2.5
1	D	194	ASP	2.5

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Mol	Chain	Res	Type	RSRZ
1	D	6	SER	2.5
1	D	341	LYS	2.5
1	A	315	MET	2.5
1	B	3	VAL	2.3
1	C	42	ASP	2.3
1	C	4	LYS	2.2
1	D	282	LEU	2.2
1	C	194	ASP	2.1
1	B	60	LEU	2.0
1	B	243	ALA	2.0

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates 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. 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
2	CIT	A	400	13/13	0.77	0.26	3.61	38,44,55,55	0
2	CIT	B	400	13/13	0.90	0.17	3.23	30,35,37,40	0
3	PEG	A	401	7/7	0.82	0.20	2.87	43,44,47,48	0
3	PEG	B	401	7/7	0.89	0.16	2.61	38,39,41,42	0
2	CIT	C	400	13/13	0.76	0.20	1.78	35,44,46,48	0
4	EDO	D	400	4/4	0.73	0.13	0.74	43,46,48,48	0
3	PEG	C	401	7/7	0.94	0.12	0.10	43,44,46,47	0
3	PEG	D	401	7/7	0.91	0.13	0.04	42,43,45,47	0
3	PEG	D	402	7/7	0.90	0.19	-	41,42,43,43	0
3	PEG	C	402	7/7	0.74	0.29	-	43,44,46,46	0
3	PEG	B	402	7/7	0.73	0.17	-	37,38,38,39	0

6.5 Other polymers

There are no such residues in this entry.