



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 9, 2026 – 10:51 AM UTC

PDB ID : 1JCH / pdb_00001jch
Title : Crystal Structure of Colicin E3 in Complex with its Immunity Protein
Authors : Soelaiman, S.; Jakes, K.; Wu, N.; Li, C.; Shoham, M.
Deposited on : 2001-06-09
Resolution : 3.02 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

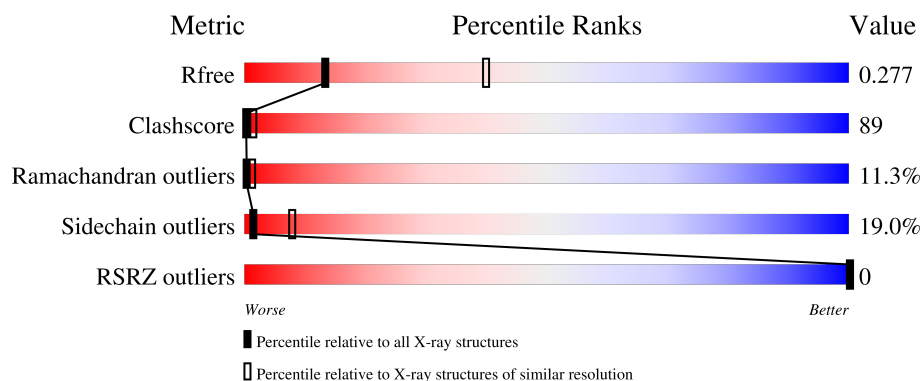
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.02 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	3131 (3.04-3.00)
Clashscore	190562	3444 (3.04-3.00)
Ramachandran outliers	187476	3319 (3.04-3.00)
Sidechain outliers	187428	3322 (3.04-3.00)
RSRZ outliers	180081	3130 (3.04-3.00)

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	551	17% 47% 17% • 15%
1	C	551	17% 46% 18% • 15%
2	B	84	32% 51% 14% •
2	D	84	31% 54% 13% •

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8998 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 COLICIN E3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	468	Total	C	N	O	S	0	0	0
			3570	2206	659	697	8			
1	C	468	Total	C	N	O	S	0	0	0
			3570	2206	659	697	8			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	260	PHE	GLY	conflict	UNP P00646
A	261	GLY	PHE	conflict	UNP P00646
C	260	PHE	GLY	conflict	UNP P00646
C	261	GLY	PHE	conflict	UNP P00646

- Molecule 2 is a protein called COLICIN E3 IMMUNITY PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	84	Total	C	N	O	S	0	0	0
			693	445	104	142	2			
2	D	84	Total	C	N	O	S	0	0	0
			693	445	104	142	2			

- Molecule 3 is CITRIC ACID (CCD ID: CIT) (formula: C₆H₈O₇).



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

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

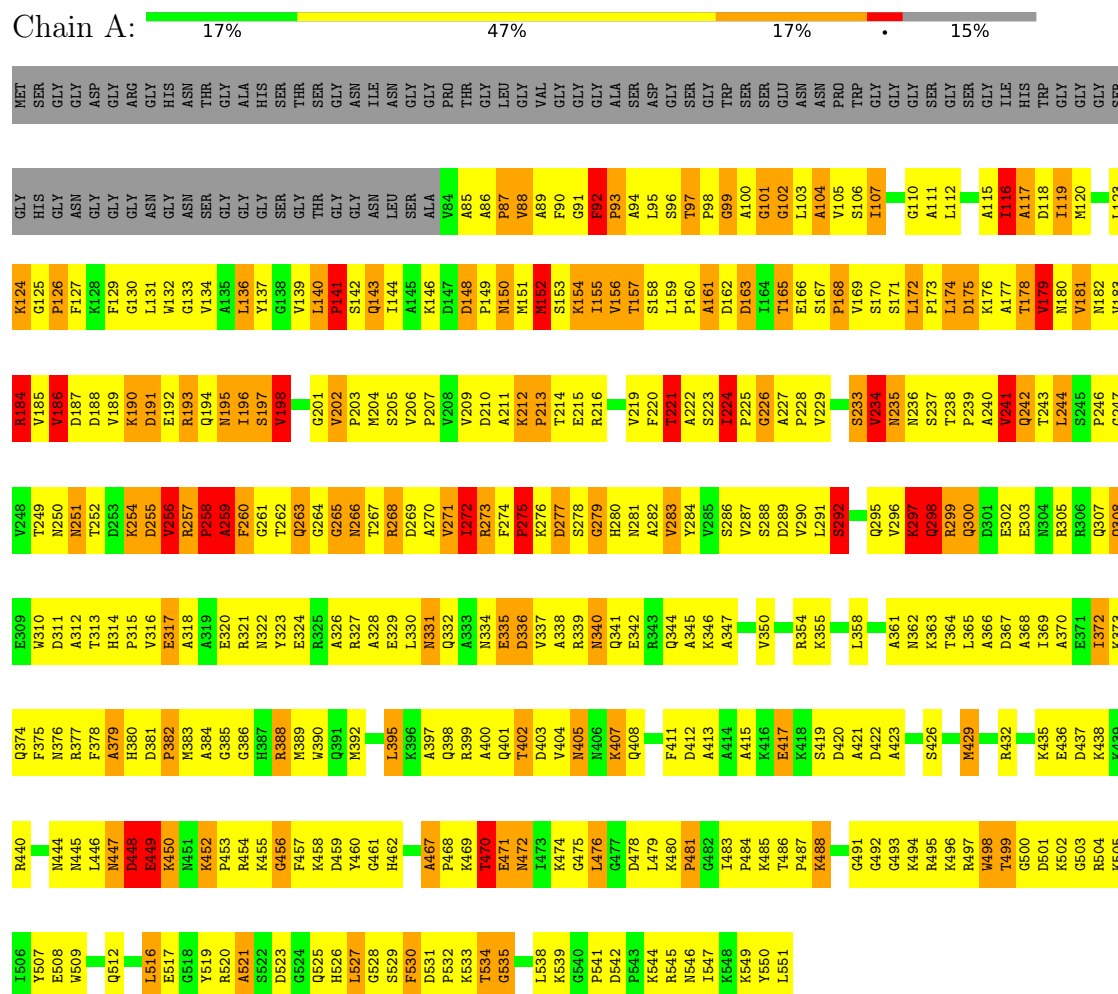
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	169	Total	O	0	0
			169	169		
5	B	28	Total	O	0	0
			28	28		
5	C	170	Total	O	0	0
			170	170		
5	D	29	Total	O	0	0
			29	29		

3 Residue-property plots

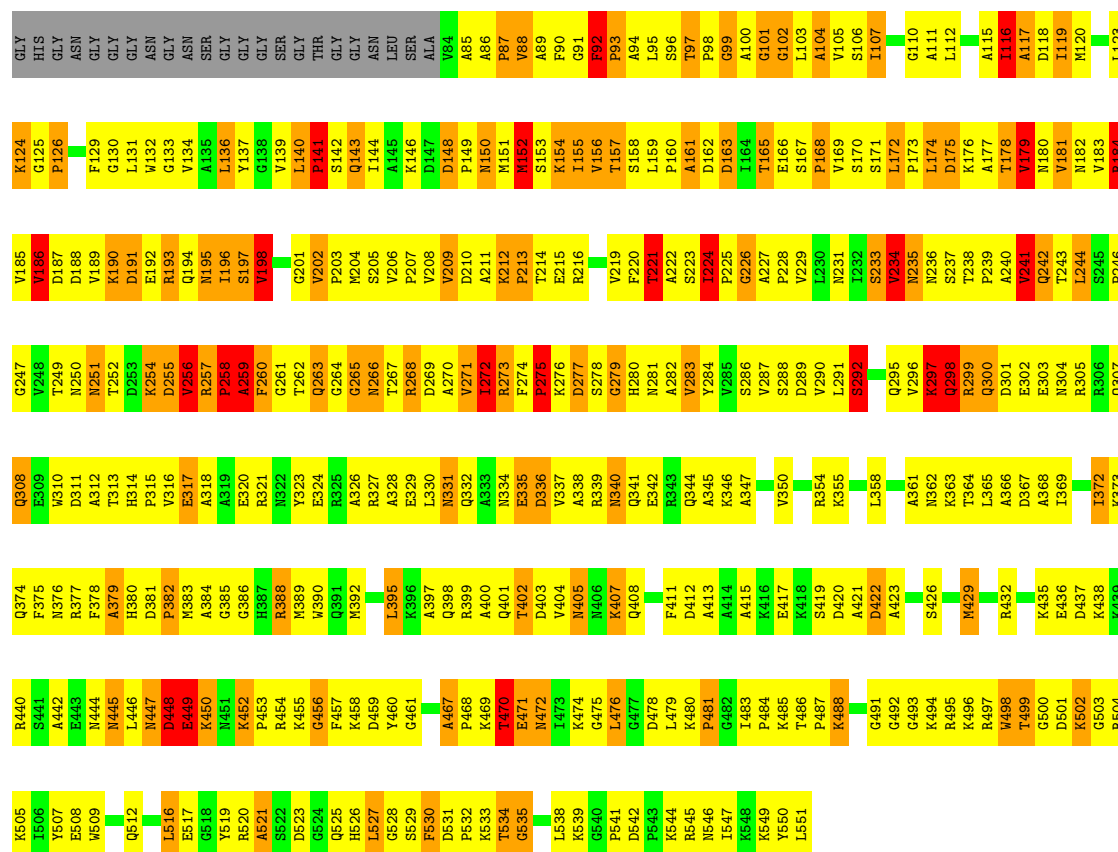
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: COLICIN E3



• Molecule 1: COLICIN E3





• Molecule 2: COLICIN E3 IMMUNITY PROTEIN

Chain B: 32% 51% 14%



• Molecule 2: COLICIN E3 IMMUNITY PROTEIN

Chain D: 31% 54% 13%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	66.99Å 195.73Å 85.12Å 90.00° 113.22° 90.00°	Depositor
Resolution (Å)	20.17 – 3.02 20.17 – 3.02	Depositor EDS
% Data completeness (in resolution range)	94.1 (20.17-3.02) 94.2 (20.17-3.02)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.26 (at 3.04Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.237 , 0.283 0.236 , 0.277	Depositor DCC
R_{free} test set	1127 reflections (3.04%)	wwPDB-VP
Wilson B-factor (Å ²)	72.3	Xtriage
Anisotropy	0.423	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 87.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	0.450 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8998	wwPDB-VP
Average B, all atoms (Å ²)	74.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.91% 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

5.1 Standard geometry

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

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	0.88	10/3639 (0.3%)	1.27	30/4922 (0.6%)
1	C	0.88	10/3639 (0.3%)	1.27	32/4922 (0.7%)
2	B	0.83	1/714 (0.1%)	1.23	5/967 (0.5%)
2	D	0.83	1/714 (0.1%)	1.23	5/967 (0.5%)
All	All	0.88	22/8706 (0.3%)	1.26	72/11778 (0.6%)

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
2	B	0	1
2	D	0	1
All	All	0	2

The worst 5 of 22 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	256	VAL	N-CA	7.07	1.55	1.46
1	C	256	VAL	N-CA	7.07	1.55	1.46
1	A	260	PHE	N-CA	-6.74	1.37	1.46
1	C	260	PHE	N-CA	-6.70	1.37	1.46
1	C	392	MET	SD-CE	6.54	1.95	1.79

The worst 5 of 72 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	92	PHE	CA-C-N	11.18	133.82	119.84
1	C	92	PHE	C-N-CA	11.18	133.82	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	92	PHE	CA-C-N	11.15	133.77	119.84
1	A	92	PHE	C-N-CA	11.15	133.77	119.84
1	A	222	ALA	N-CA-C	10.01	124.88	112.87

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	61	TYR	Sidechain
2	D	61	TYR	Sidechain

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	3570	0	3532	688	8
1	C	3570	0	3532	695	9
2	B	693	0	617	78	0
2	D	693	0	617	78	0
3	A	26	0	10	1	0
3	C	26	0	10	1	0
4	A	6	0	8	0	0
4	B	6	0	8	1	0
4	C	6	0	8	0	0
4	D	6	0	8	1	0
5	A	169	0	0	42	1
5	B	28	0	0	2	0
5	C	170	0	0	44	0
5	D	29	0	0	2	0
All	All	8998	0	8350	1502	9

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 89.

The worst 5 of 1502 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:C:257:ARG:HB3	1:C:258:PRO:CD	1.41	1.39
1:A:257:ARG:HB3	1:A:258:PRO:CD	1.41	1.35
1:C:179:VAL:HB	1:C:286:SER:HB2	1.22	1.17
1:C:272:ILE:HB	1:C:282:ALA:HA	1.22	1.16
1:A:257:ARG:HB2	1:A:303:GLU:OE2	1.45	1.16

The worst 5 of 9 symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:234:VAL:CG2	1:C:324:GLU:OE2[2_655]	1.62	0.58
1:A:324:GLU:OE2	1:C:234:VAL:CG2[2_555]	1.86	0.34
1:A:126:PRO:O	1:C:239:PRO:CD[2_555]	1.88	0.32
1:A:239:PRO:CD	1:C:126:PRO:O[2_655]	1.94	0.26
1:A:127:PHE:CD1	1:C:236:ASN:O[2_555]	2.00	0.20

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	466/551 (85%)	294 (63%)	115 (25%)	57 (12%)	0	1
1	C	466/551 (85%)	294 (63%)	115 (25%)	57 (12%)	0	1
2	B	82/84 (98%)	63 (77%)	14 (17%)	5 (6%)	1	6
2	D	82/84 (98%)	63 (77%)	14 (17%)	5 (6%)	1	6
All	All	1096/1270 (86%)	714 (65%)	258 (24%)	124 (11%)	0	1

5 of 124 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	101	GLY
1	A	102	GLY

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Mol	Chain	Res	Type
1	A	142	SER
1	A	152	MET
1	A	155	ILE

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	377/422 (89%)	302 (80%)	75 (20%)	1	7
1	C	377/422 (89%)	302 (80%)	75 (20%)	1	7
2	B	76/76 (100%)	65 (86%)	11 (14%)	3	14
2	D	76/76 (100%)	65 (86%)	11 (14%)	3	14
All	All	906/996 (91%)	734 (81%)	172 (19%)	1	8

5 of 172 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	234	VAL
1	C	402	THR
1	C	252	THR
1	C	298	GLN
1	C	470	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 45 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	266	ASN
1	C	374	GLN
1	C	298	GLN
1	C	308	GLN
1	C	401	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

8 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$
3	CIT	A	601	-	12,12,12	3.21	4 (33%)	17,17,17	1.67	4 (23%)
4	GOL	C	701	-	5,5,5	1.21	1 (20%)	5,5,5	0.59	0
4	GOL	A	5701	-	5,5,5	1.21	1 (20%)	5,5,5	0.59	0
4	GOL	B	702	-	5,5,5	1.07	0	5,5,5	0.24	0
3	CIT	C	5601	-	12,12,12	3.20	4 (33%)	17,17,17	1.66	4 (23%)
3	CIT	C	5602	-	12,12,12	3.06	3 (25%)	17,17,17	1.79	5 (29%)
3	CIT	A	602	-	12,12,12	3.06	3 (25%)	17,17,17	1.79	5 (29%)
4	GOL	D	5702	-	5,5,5	1.07	0	5,5,5	0.24	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
3	CIT	A	601	-	-	2/16/16/16	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	GOL	C	701	-	-	2/4/4/4	-
4	GOL	A	5701	-	-	2/4/4/4	-
4	GOL	B	702	-	-	4/4/4/4	-
3	CIT	C	5601	-	-	2/16/16/16	-
3	CIT	C	5602	-	-	2/16/16/16	-
3	CIT	A	602	-	-	2/16/16/16	-
4	GOL	D	5702	-	-	4/4/4/4	-

The worst 5 of 16 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	601	CIT	O1-C1	7.06	1.45	1.22
3	C	5601	CIT	O1-C1	7.05	1.45	1.22
3	C	5602	CIT	O1-C1	6.58	1.43	1.22
3	A	602	CIT	O1-C1	6.57	1.43	1.22
3	A	602	CIT	C2-C3	5.55	1.60	1.54

The worst 5 of 18 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	601	CIT	O6-C6-O5	-3.93	111.28	123.86
3	C	5602	CIT	O6-C6-O5	-3.93	111.28	123.86
3	A	601	CIT	O6-C6-O5	-3.68	112.07	123.86
3	C	5601	CIT	O6-C6-O5	-3.68	112.08	123.86
3	A	601	CIT	O7-C3-C2	2.61	115.32	109.38

There are no chirality outliers.

5 of 20 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	601	CIT	C2-C3-C6-O5
3	A	601	CIT	O7-C3-C6-O5
3	A	602	CIT	C2-C3-C6-O5
3	A	602	CIT	O7-C3-C6-O5
3	C	5601	CIT	C2-C3-C6-O5

There are no ring outliers.

4 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	601	CIT	1	0
4	B	702	GOL	1	0
3	C	5601	CIT	1	0
4	D	5702	GOL	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	468/551 (84%)	-1.56	0 100 100	25, 73, 122, 147	0
1	C	468/551 (84%)	-1.57	0 100 100	25, 73, 122, 147	0
2	B	84/84 (100%)	-1.60	0 100 100	26, 58, 87, 95	0
2	D	84/84 (100%)	-1.59	0 100 100	26, 58, 87, 95	0
All	All	1104/1270 (86%)	-1.57	0 100 100	25, 70, 120, 147	0

There are no RSRZ outliers to report.

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 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
3	CIT	C	5602	13/13	0.97	0.03	103,143,168,170	0
3	CIT	A	602	13/13	0.98	0.03	103,143,168,170	0
3	CIT	A	601	13/13	0.98	0.04	67,122,144,151	0
4	GOL	A	5701	6/6	0.98	0.06	63,76,79,81	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	GOL	C	701	6/6	0.98	0.07	63,76,79,81	0
4	GOL	B	702	6/6	0.99	0.03	76,101,102,106	0
3	CIT	C	5601	13/13	0.99	0.04	67,122,144,151	0
4	GOL	D	5702	6/6	0.99	0.03	76,101,102,106	0

6.5 Other polymers [i](#)

There are no such residues in this entry.