



wwPDB X-ray Structure Validation Summary Report

Feb 28, 2014 – 12:19 PM GMT

PDB ID : 1T5E
Title : The structure of MexA
Authors : Higgins, M.K.; Bokma, E.; Koronakis, E.; Hughes, C.; Koronakis, V.
Deposited on : 2004-05-04
Resolution : 3.00 Å(reported)

This is a wwPDB validation summary 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/ValidationPDFNotes.html>

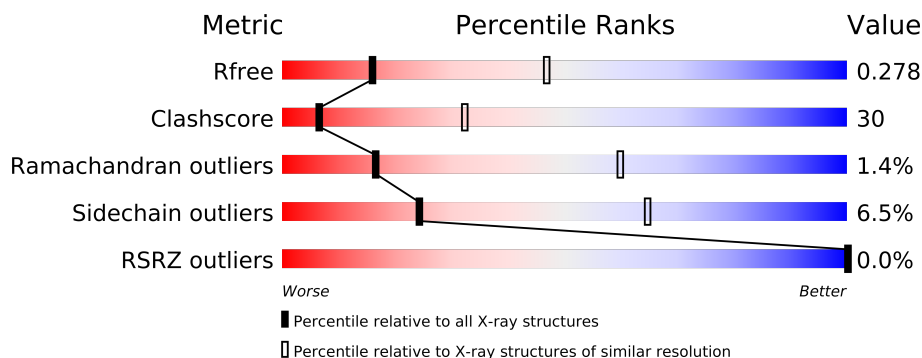
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.15 2013
Xtriage (Phenix)	:	dev-1323
EDS	:	stable22639
Percentile statistics	:	21963
Refmac	:	5.8.0049
CCP4	:	6.3.0 (Settle)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP)	:	stable22683

1 Overall quality at a glance

The reported resolution of this entry is 3.00 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}	66092	1216 (3.00-3.00)
Clashscore	79885	1594 (3.00-3.00)
Ramachandran outliers	78287	1537 (3.00-3.00)
Sidechain outliers	78261	1540 (3.00-3.00)
RSRZ outliers	66119	1217 (3.00-3.00)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	360	
1	B	360	
1	C	360	
1	D	360	
1	E	360	
1	F	360	
1	G	360	
1	H	360	
1	I	360	
1	J	360	
1	K	360	
1	L	360	
1	M	360	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
2	3GR	E	361	-	X
2	3GR	H	361	-	X
2	3GR	K	361	-	X
2	3GR	L	361	-	X
3	GOL	J	361	-	X

2 Entry composition

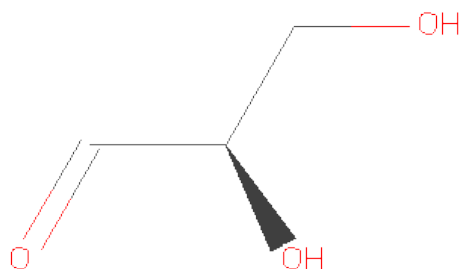
There are 3 unique types of molecules in this entry. The entry contains 23101 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Multidrug resistance protein mexA.

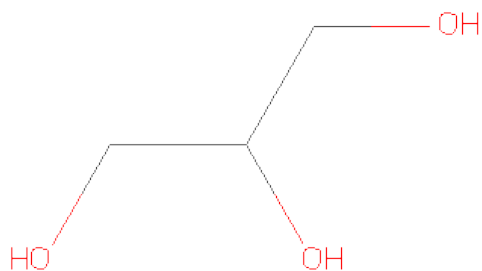
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	231	Total	C	N	O	S	0	0	0
			1771	1101	319	349	2			
1	B	231	Total	C	N	O	S	0	0	0
			1771	1101	319	349	2			
1	C	231	Total	C	N	O	S	0	0	0
			1771	1101	319	349	2			
1	D	231	Total	C	N	O	S	0	0	0
			1771	1101	319	349	2			
1	E	231	Total	C	N	O	S	0	0	0
			1771	1101	319	349	2			
1	F	231	Total	C	N	O	S	0	0	0
			1771	1101	319	349	2			
1	G	231	Total	C	N	O	S	0	0	0
			1771	1101	319	349	2			
1	H	231	Total	C	N	O	S	0	0	0
			1771	1101	319	349	2			
1	I	231	Total	C	N	O	S	0	0	0
			1771	1101	319	349	2			
1	J	231	Total	C	N	O	S	0	0	0
			1771	1101	319	349	2			
1	K	231	Total	C	N	O	S	0	0	0
			1771	1101	319	349	2			
1	L	231	Total	C	N	O	S	0	0	0
			1771	1101	319	349	2			
1	M	231	Total	C	N	O	S	0	0	0
			1771	1101	319	349	2			

- Molecule 2 is GLYCERALDEHYDE (three-letter code: 3GR) (formula: C₃H₆O₃).

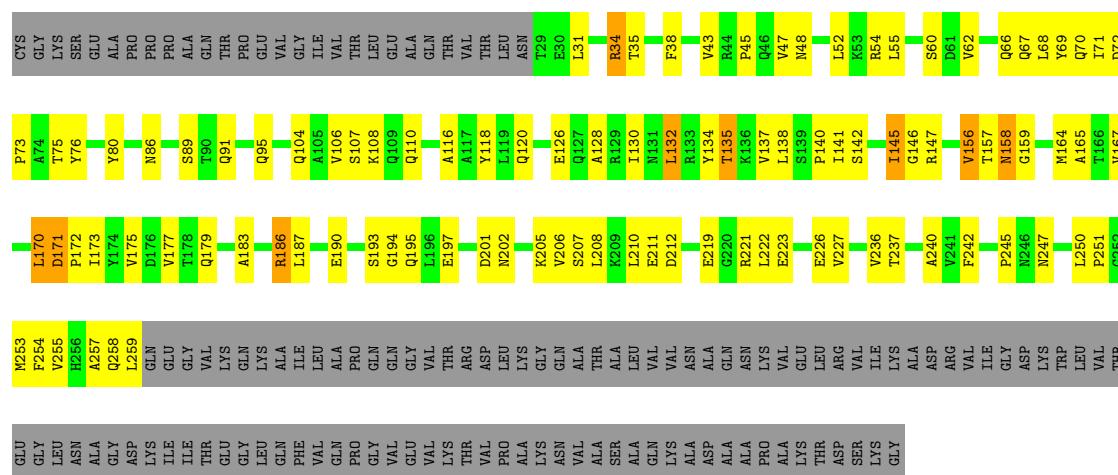


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	C	1	Total	C	O	0	0
			6	3	3		
2	D	1	Total	C	O	0	0
			6	3	3		
2	E	1	Total	C	O	0	0
			6	3	3		
2	F	1	Total	C	O	0	0
			6	3	3		
2	G	1	Total	C	O	0	0
			6	3	3		
2	H	1	Total	C	O	0	0
			6	3	3		
2	I	1	Total	C	O	0	0
			6	3	3		
2	K	1	Total	C	O	0	0
			6	3	3		
2	L	1	Total	C	O	0	0
			6	3	3		

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).

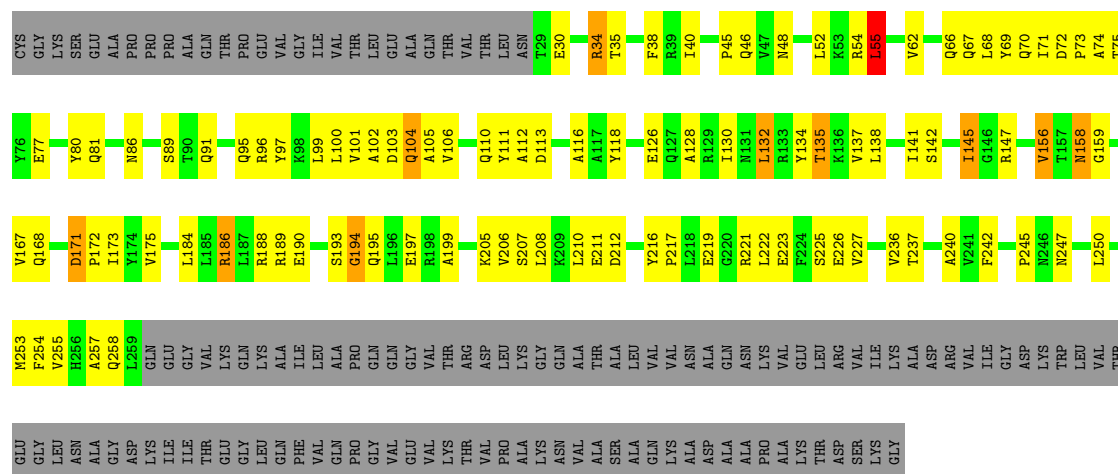


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	J	1	Total	C	O	0	0
			6	3	3		
3	M	1	Total	C	O	0	0
			6	3	3		



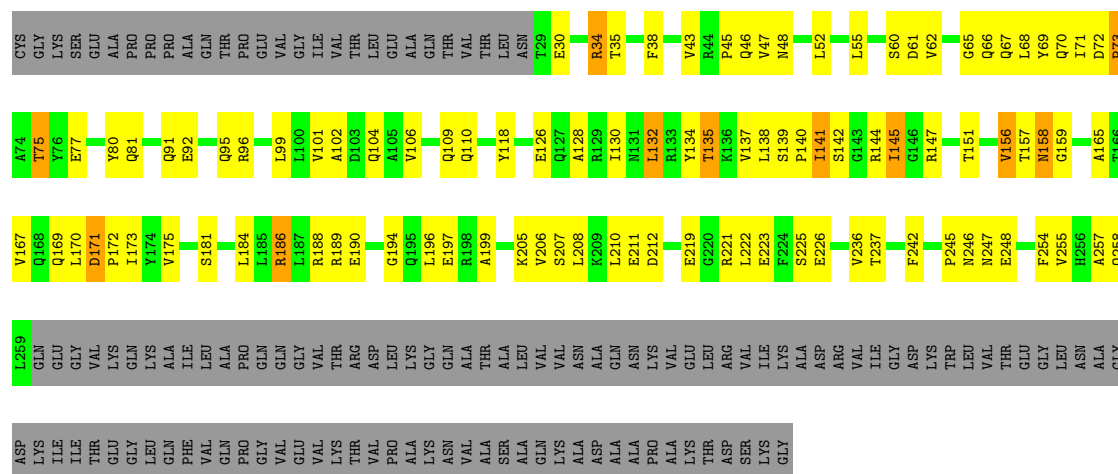
• Molecule 1: Multidrug resistance protein mexA

Chain D:

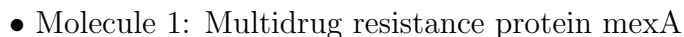


• Molecule 1: Multidrug resistance protein mexA

Chain E:

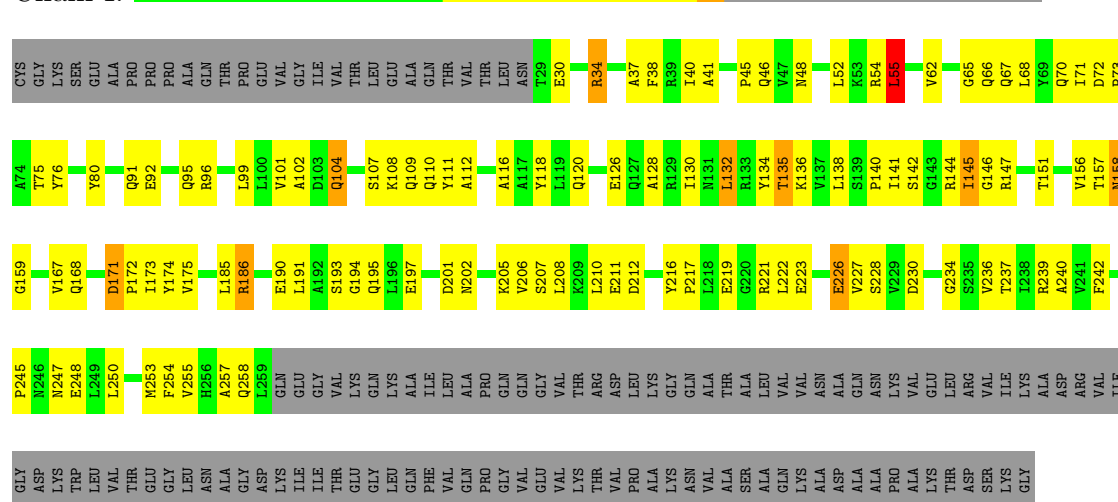


• Molecule 1: Multidrug resistance protein mexA



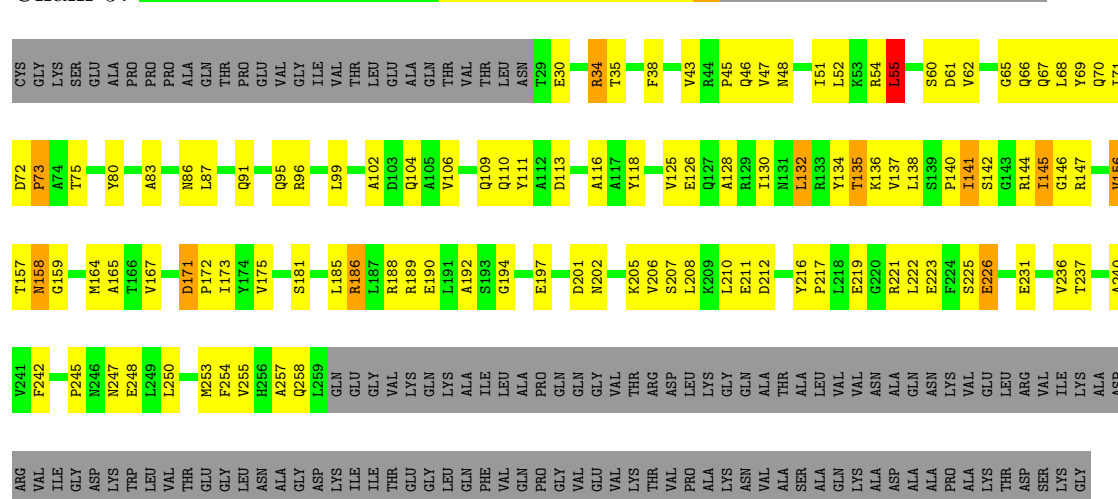
- Molecule 1: Multidrug resistance protein mexA

Chain I:



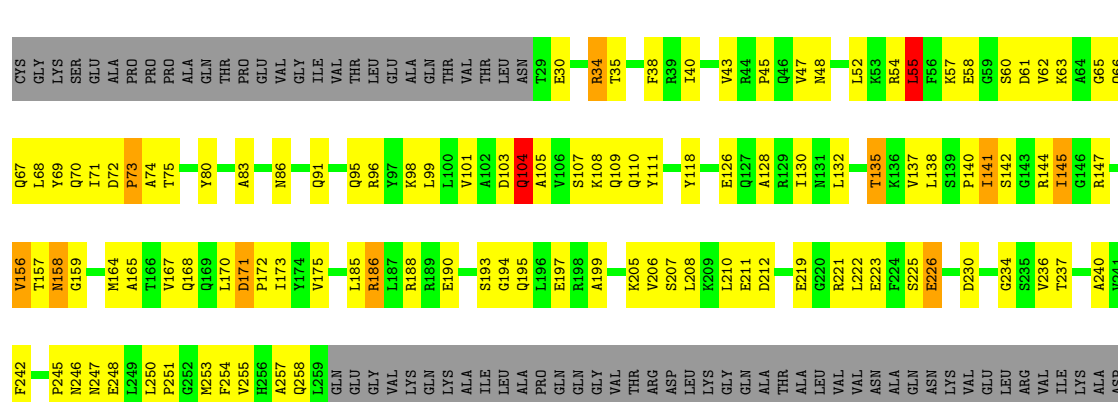
- Molecule 1: Multidrug resistance protein mexA

Chain J:



- Molecule 1: Multidrug resistance protein mexA

Chain K:



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	130.55Å 183.59Å 213.31Å 90.00° 107.38° 90.00°	Depositor
Resolution (Å)	95.00 – 3.00 83.48 – 2.99	Depositor EDS
% Data completeness (in resolution range)	97.8 (95.00-3.00) 98.3 (83.48-2.99)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.64 (at 3.01Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.273 , 0.285 0.268 , 0.278	Depositor DCC
R_{free} test set	9310 reflections (4.91%)	DCC
Wilson B-factor (Å ²)	75.0	Xtriage
Anisotropy	0.023	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 33.9	EDS
Estimated twinning fraction	0.005 for h,-k,-h-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 190354 reflections	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	23101	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 6.83% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, 3GR

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	0.48	1/1795 (0.1%)	0.62	0/2434
1	B	0.39	0/1795	0.65	0/2434
1	C	0.40	0/1795	0.66	0/2434
1	D	0.42	0/1795	0.69	0/2434
1	E	0.47	0/1795	0.71	0/2434
1	F	0.38	0/1795	0.67	0/2434
1	G	0.37	0/1795	0.63	0/2434
1	H	0.45	0/1795	0.70	0/2434
1	I	0.49	0/1795	0.71	0/2434
1	J	0.56	0/1795	0.75	0/2434
1	K	0.52	0/1795	0.73	0/2434
1	L	0.48	0/1795	0.71	0/2434
1	M	0.39	0/1795	0.66	0/2434
All	All	0.45	1/23335 (0.0%)	0.68	0/31642

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	1
1	I	0	1
All	All	0	2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	29	THR	C-N	-14.50	1.00	1.34

There are no bond angle outliers.

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	29	THR	Mainchain
1	I	111	TYR	Sidechain

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1771	0	1774	109	1
1	B	1771	0	1775	105	0
1	C	1771	0	1775	98	0
1	D	1771	0	1775	110	1
1	E	1771	0	1775	102	0
1	F	1771	0	1775	109	0
1	G	1771	0	1775	120	0
1	H	1771	0	1775	109	0
1	I	1771	0	1775	115	0
1	J	1771	0	1775	110	0
1	K	1771	0	1775	120	0
1	L	1771	0	1775	127	0
1	M	1771	0	1775	110	0
2	C	6	0	5	2	0
2	D	6	0	5	0	0
2	E	6	0	5	2	0
2	F	6	0	5	4	0
2	G	6	0	5	1	0
2	H	6	0	5	2	0
2	I	6	0	5	4	0
2	K	6	0	5	3	0
2	L	6	0	5	1	0
3	A	6	0	8	5	0
3	B	6	0	5	3	0
3	J	6	0	5	5	0
3	M	6	0	5	2	0
All	All	23101	0	23142	1365	1

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 30.

The worst 5 of 1365 close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:F:146:GLY:HA3	2:F:361:3GR:O2	1.37	1.24
1:M:70:GLN:HE22	1:M:135:THR:HG23	1.08	1.15
1:A:70:GLN:HE22	1:A:135:THR:HG23	1.18	1.07
1:F:91:GLN:HG2	1:F:95:GLN:HE21	1.21	1.05
1:J:91:GLN:HG2	1:J:95:GLN:HE21	1.16	1.04

All (1) 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	Distance(Å)	Clash(Å)
1:A:189:ARG:NH2	1:D:194:GLY:O[2_656]	2.19	0.01

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	229/360 (64%)	207 (90%)	21 (9%)	1 (0%)	43	87
1	B	229/360 (64%)	203 (89%)	24 (10%)	2 (1%)	25	73
1	C	229/360 (64%)	209 (91%)	18 (8%)	2 (1%)	25	73
1	D	229/360 (64%)	200 (87%)	24 (10%)	5 (2%)	10	45
1	E	229/360 (64%)	211 (92%)	15 (7%)	3 (1%)	18	62
1	F	229/360 (64%)	209 (91%)	17 (7%)	3 (1%)	18	62
1	G	229/360 (64%)	205 (90%)	20 (9%)	4 (2%)	14	54
1	H	229/360 (64%)	205 (90%)	19 (8%)	5 (2%)	10	45
1	I	229/360 (64%)	210 (92%)	16 (7%)	3 (1%)	18	62
1	J	229/360 (64%)	208 (91%)	17 (7%)	4 (2%)	14	54
1	K	229/360 (64%)	205 (90%)	20 (9%)	4 (2%)	14	54
1	L	229/360 (64%)	205 (90%)	20 (9%)	4 (2%)	14	54

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	M	229/360 (64%)	204 (89%)	22 (10%)	3 (1%)	18 62
All	All	2977/4680 (64%)	2681 (90%)	253 (8%)	43 (1%)	16 60

5 of 43 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	171	ASP
1	B	171	ASP
1	C	171	ASP
1	D	171	ASP
1	E	171	ASP

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	186/287 (65%)	176 (95%)	10 (5%)	31 74
1	B	186/287 (65%)	175 (94%)	11 (6%)	28 70
1	C	186/287 (65%)	176 (95%)	10 (5%)	31 74
1	D	186/287 (65%)	173 (93%)	13 (7%)	21 62
1	E	186/287 (65%)	170 (91%)	16 (9%)	15 50
1	F	186/287 (65%)	173 (93%)	13 (7%)	21 62
1	G	186/287 (65%)	175 (94%)	11 (6%)	28 70
1	H	186/287 (65%)	174 (94%)	12 (6%)	24 65
1	I	186/287 (65%)	176 (95%)	10 (5%)	31 74
1	J	186/287 (65%)	171 (92%)	15 (8%)	17 53
1	K	186/287 (65%)	171 (92%)	15 (8%)	17 53
1	L	186/287 (65%)	174 (94%)	12 (6%)	24 65
1	M	186/287 (65%)	177 (95%)	9 (5%)	35 79
All	All	2418/3731 (65%)	2261 (94%)	157 (6%)	24 65

5 of 157 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	F	189	ARG
1	H	103	ASP
1	L	186	ARG
1	G	34	ARG
1	G	145	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 97 such sidechains are listed below:

Mol	Chain	Res	Type
1	G	84	GLN
1	H	95	GLN
1	M	70	GLN
1	G	110	GLN
1	G	169	GLN

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

13 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	GOL	A	361	-	5,5,5	0.19	0	5,5,5	0.49	0
3	GOL	B	361	-	5,5,5	4.61	1 (20%)	5,5,5	4.19	1 (20%)
2	3GR	C	361	-	5,5,5	2.08	1 (20%)	5,5,5	5.35	1 (20%)
2	3GR	D	361	-	5,5,5	2.08	1 (20%)	5,5,5	5.35	1 (20%)
2	3GR	E	361	-	5,5,5	2.09	1 (20%)	5,5,5	5.34	1 (20%)
2	3GR	F	361	-	5,5,5	2.08	1 (20%)	5,5,5	5.34	1 (20%)
2	3GR	G	361	-	5,5,5	2.09	1 (20%)	5,5,5	5.34	1 (20%)
2	3GR	H	361	-	5,5,5	2.08	1 (20%)	5,5,5	5.35	1 (20%)
2	3GR	I	361	-	5,5,5	2.09	1 (20%)	5,5,5	5.34	1 (20%)
3	GOL	J	361	1	5,5,5	4.61	1 (20%)	5,5,5	4.19	1 (20%)
2	3GR	K	361	-	5,5,5	2.08	1 (20%)	5,5,5	5.35	1 (20%)
2	3GR	L	361	-	5,5,5	2.08	1 (20%)	5,5,5	5.35	1 (20%)
3	GOL	M	361	1	5,5,5	4.61	1 (20%)	5,5,5	4.20	1 (20%)

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	GOL	A	361	-	-	0/4/4/4	0/0/0/0
3	GOL	B	361	-	-	0/4/4/4	0/0/0/0
2	3GR	C	361	-	1/1/1/2	0/3/4/4	0/0/0/0
2	3GR	D	361	-	1/1/1/2	0/3/4/4	0/0/0/0
2	3GR	E	361	-	1/1/1/2	0/3/4/4	0/0/0/0
2	3GR	F	361	-	1/1/1/2	0/3/4/4	0/0/0/0
2	3GR	G	361	-	1/1/1/2	0/3/4/4	0/0/0/0
2	3GR	H	361	-	1/1/1/2	0/3/4/4	0/0/0/0
2	3GR	I	361	-	1/1/1/2	0/3/4/4	0/0/0/0
3	GOL	J	361	1	-	0/4/4/4	0/0/0/0
2	3GR	K	361	-	1/1/1/2	0/3/4/4	0/0/0/0
2	3GR	L	361	-	1/1/1/2	0/3/4/4	0/0/0/0
3	GOL	M	361	1	-	0/4/4/4	0/0/0/0

The worst 5 of 12 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	M	361	GOL	O3-C3	-10.26	0.97	1.42
3	B	361	GOL	O3-C3	-10.25	0.97	1.42
3	J	361	GOL	O3-C3	-10.25	0.97	1.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	G	361	3GR	O3-C3	-4.44	0.97	1.19
2	I	361	3GR	O3-C3	-4.44	0.97	1.19

The worst 5 of 12 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	361	3GR	O3-C3-C2	11.90	155.34	125.34
2	D	361	3GR	O3-C3-C2	11.89	155.32	125.34
2	K	361	3GR	O3-C3-C2	11.90	155.32	125.34
2	H	361	3GR	O3-C3-C2	11.89	155.32	125.34
2	L	361	3GR	O3-C3-C2	11.89	155.31	125.34

5 of 9 chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	E	361	3GR	C2
2	I	361	3GR	C2
2	C	361	3GR	C2
2	G	361	3GR	C2
2	H	361	3GR	C2

There are no torsion outliers.

There are no ring outliers.

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	231/360 (64%)	0.15	1 (0%)	90	41	56, 93, 132, 143	2 (0%)
1	B	231/360 (64%)	-0.04	0	100	100	37, 80, 110, 123	2 (0%)
1	C	231/360 (64%)	-0.06	0	100	100	39, 77, 111, 116	2 (0%)
1	D	231/360 (64%)	-0.03	0	100	100	30, 82, 109, 123	2 (0%)
1	E	231/360 (64%)	-0.07	0	100	100	22, 61, 95, 111	2 (0%)
1	F	231/360 (64%)	-0.05	0	100	100	53, 80, 106, 111	2 (0%)
1	G	231/360 (64%)	0.01	0	100	100	51, 86, 118, 126	2 (0%)
1	H	231/360 (64%)	-0.02	0	100	100	28, 64, 98, 109	2 (0%)
1	I	231/360 (64%)	0.01	0	100	100	24, 62, 96, 109	2 (0%)
1	J	231/360 (64%)	-0.05	0	100	100	15, 48, 84, 100	2 (0%)
1	K	231/360 (64%)	-0.07	0	100	100	20, 59, 92, 107	2 (0%)
1	L	231/360 (64%)	-0.07	0	100	100	25, 63, 102, 115	2 (0%)
1	M	231/360 (64%)	0.02	0	100	100	40, 93, 117, 130	2 (0%)
All	All	3003/4680 (64%)	-0.02	1 (0%)	100	100	15, 75, 114, 143	26 (0%)

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	232	GLY	2.2

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	3GR	E	361	6/6	0.36	8.17	93,95,98,102	0
3	GOL	J	361	6/6	0.40	6.63	92,94,94,96	0
2	3GR	K	361	6/6	0.35	4.35	79,80,83,84	0
2	3GR	L	361	6/6	0.33	2.69	81,82,83,86	0
2	3GR	H	361	6/6	0.32	2.28	87,88,89,90	0
2	3GR	D	361	6/6	0.32	1.85	93,93,94,95	0
2	3GR	C	361	6/6	0.28	1.85	100,100,101,102	0
3	GOL	A	361	6/6	0.30	1.53	97,98,99,102	0
3	GOL	M	361	6/6	0.26	1.43	99,100,101,101	0
2	3GR	I	361	6/6	0.28	0.98	87,87,88,88	0
3	GOL	B	361	6/6	0.29	0.82	95,96,96,96	0
2	3GR	F	361	6/6	0.25	0.31	99,100,101,103	0
2	3GR	G	361	6/6	0.20	-0.54	103,104,105,106	0

6.5 Other polymers ⓘ

There are no such residues in this entry.