



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 27, 2016 – 11:58 PM EDT

PDB ID : 5LGD
Title : The CIDRa domain from MCvar1 PfEMP1 bound to CD36
Authors : Hsieh, F.L.; Higgins, M.K.
Deposited on : 2016-07-06
Resolution : 2.07 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 : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20027939
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) : rb-20027939

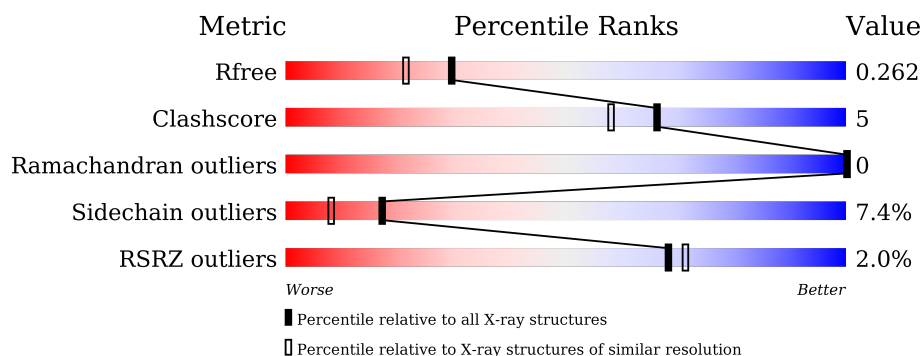
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.07 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	1799 (2.08-2.04)
Clashscore	102246	1910 (2.08-2.04)
Ramachandran outliers	100387	1893 (2.08-2.04)
Sidechain outliers	100360	1893 (2.08-2.04)
RSRZ outliers	91569	1802 (2.08-2.04)

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	472	<div> <div>%</div> <div> <div></div> <div>70%</div> <div>14%</div> <div>•</div> <div>15%</div> </div> </div>
2	B	179	<div> <div>3%</div> <div> <div></div> <div>72%</div> <div>10%</div> <div>•</div> <div>16%</div> </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
5	PLM	A	512	-	-	X	X
5	PLM	A	513	-	-	-	X
6	GOL	A	515	-	-	-	X
6	GOL	A	517	-	-	-	X
6	GOL	B	801	-	-	-	X
7	TRS	A	518	-	-	-	X
7	TRS	B	802	-	-	-	X

2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 4875 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 Platelet glycoprotein 4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	400	Total	C	N	O	S	0	0	0
			3218	2071	533	602	12			

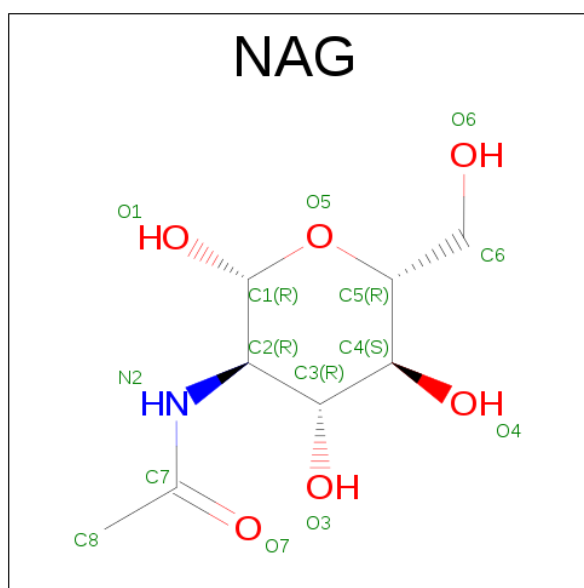
- Molecule 2 is a protein called PfEMP1 variant 1 of strain MC.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	151	Total	C	N	O	S	0	2	0
			1272	805	220	237	10			

There is a discrepancy between the modelled and reference sequences:

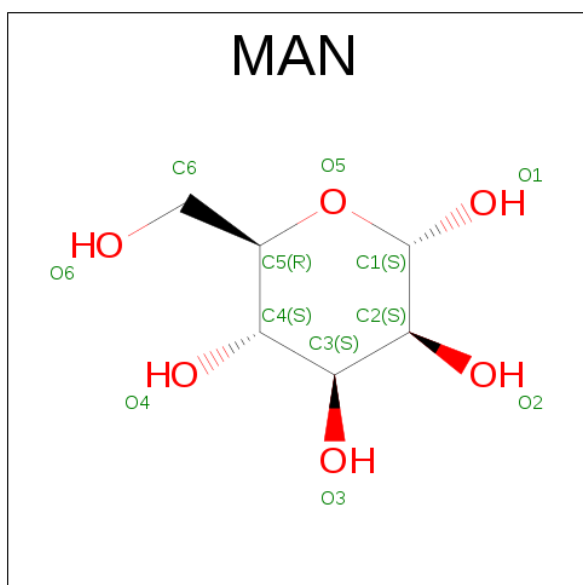
Chain	Residue	Modelled	Actual	Comment	Reference
B	626	SER	CYS	conflict	UNP Q25733

- Molecule 3 is N-ACETYL-D-GLUCOSAMINE (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



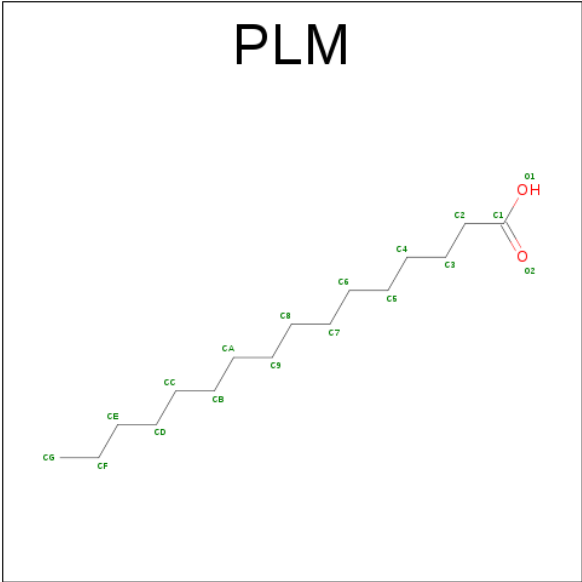
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	A	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 4 is ALPHA-D-MANNOSE (three-letter code: MAN) (formula: $C_6H_{12}O_6$).



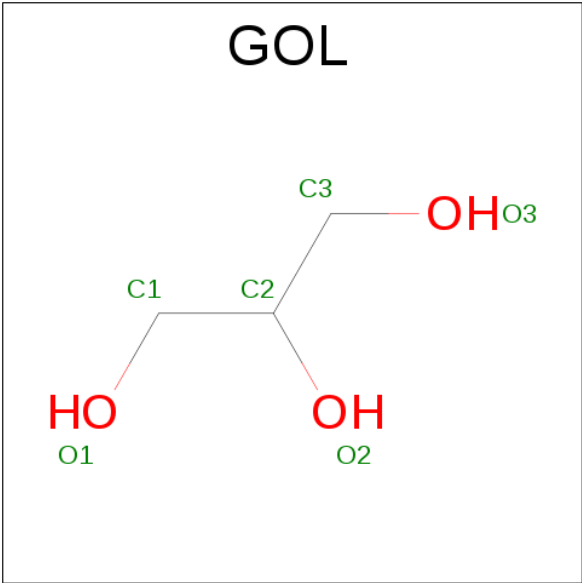
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			11	6	5		

- Molecule 5 is PALMITIC ACID (three-letter code: PLM) (formula: $C_{16}H_{32}O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			18	16	2		
5	A	1	Total	C	O	0	0
			18	16	2		

- Molecule 6 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



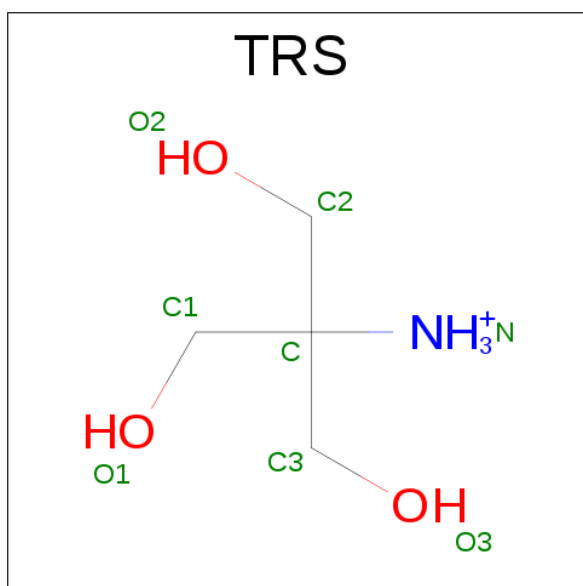
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			6	3	3		
6	A	1	Total	C	O	0	0
			6	3	3		

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

- Molecule 7 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (three-letter code: TRS) (formula: C₄H₁₂NO₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	A	1	Total	C	N	O	0	0
			8	4	1	3		
7	B	1	Total	C	N	O	0	0
			8	4	1	3		

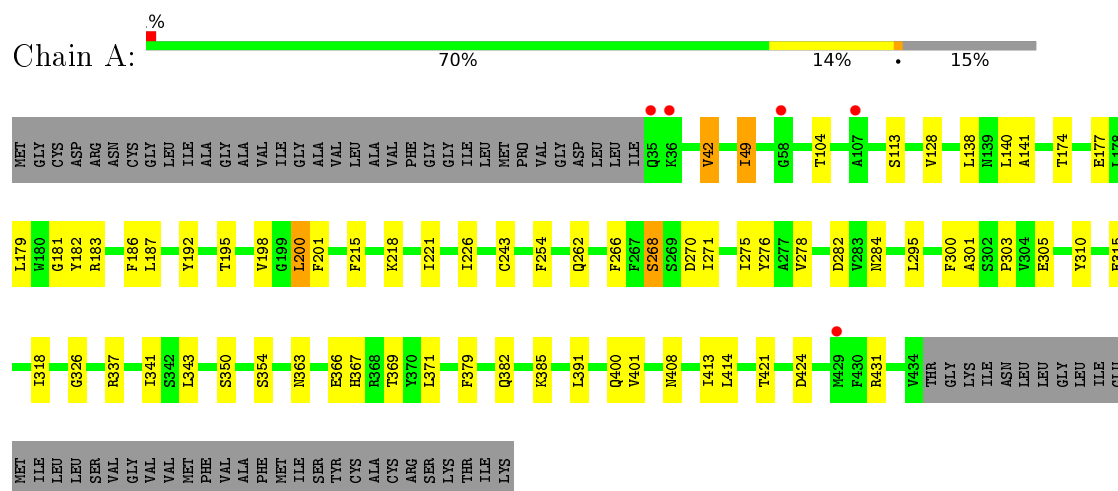
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	119	Total	O	0	0
			119	119		
8	B	33	Total	O	0	0
			33	33		

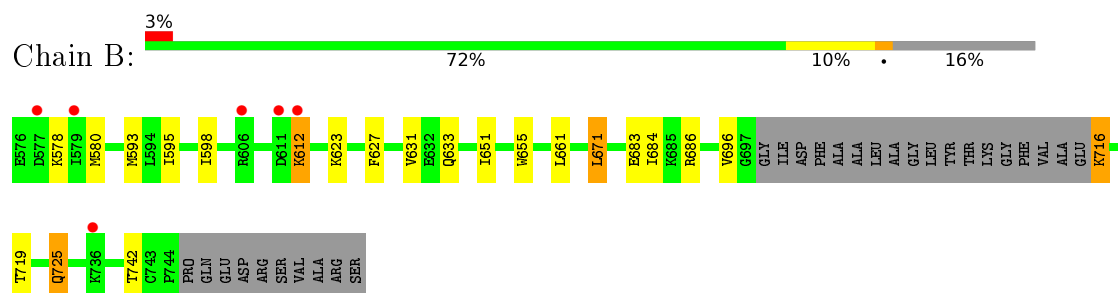
3 Residue-property plots [i](#)

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: Platelet glycoprotein 4



• Molecule 2: PfEMP1 variant 1 of strain MC



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	129.83 Å 40.73 Å 138.84 Å 90.00° 114.85° 90.00°	Depositor
Resolution (Å)	62.99 – 2.07 58.90 – 2.07	Depositor EDS
% Data completeness (in resolution range)	99.2 (62.99-2.07) 99.3 (58.90-2.07)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.66 (at 2.07 Å)	Xtriage
Refinement program	BUSTER 2.10.2	Depositor
R, R_{free}	0.212 , 0.253 0.219 , 0.262	Depositor DCC
R_{free} test set	2057 reflections (5.35%)	DCC
Wilson B-factor (Å ²)	33.0	Xtriage
Anisotropy	0.351	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 48.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.013 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	4875	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.50% 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: GOL, TRS, MAN, NAG, PLM

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	0.53	0/3296	0.73	0/4478
2	B	0.52	0/1300	0.68	0/1741
All	All	0.53	0/4596	0.72	0/6219

There are no bond length outliers.

There are no bond angle outliers.

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	3218	0	3180	38	0
2	B	1272	0	1257	9	0
3	A	140	0	128	1	0
4	A	11	0	10	0	0
5	A	36	0	62	10	0
6	A	24	0	32	1	0
6	B	6	0	8	0	0
7	A	8	0	12	2	0
7	B	8	0	12	0	0
8	A	119	0	0	1	0
8	B	33	0	0	1	0
All	All	4875	0	4701	46	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 (46) 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:A:337:ARG:HD3	5:A:512:PLM:H31	1.70	0.74
1:A:141:ALA:HA	5:A:512:PLM:H42	1.70	0.73
2:B:716:LYS:HA	2:B:725:GLN:HE22	1.61	0.66
2:B:595:ILE:HD11	2:B:683:GLU:HG3	1.79	0.65
1:A:42:VAL:HG12	1:A:221:ILE:HD11	1.81	0.63
2:B:598:ILE:HG21	2:B:686[A]:ARG:HE	1.64	0.62
1:A:198:VAL:HG11	5:A:512:PLM:H41	1.85	0.58
1:A:266:PHE:HZ	1:A:295:LEU:HD11	1.67	0.57
1:A:301:ALA:HB1	1:A:305:GLU:HB2	1.87	0.57
2:B:612:LYS:HE3	2:B:612:LYS:H	1.70	0.57
1:A:363:ASN:H	1:A:367:HIS:HD2	1.52	0.56
1:A:187:LEU:HD11	5:A:512:PLM:H21	1.86	0.56
1:A:408:ASN:HB2	7:A:518:TRS:H21	1.87	0.55
1:A:303:PRO:HG3	1:A:310:TYR:HA	1.89	0.54
1:A:201:PHE:HE1	5:A:512:PLM:HG3	1.71	0.54
1:A:382:GLN:HE21	1:A:421:THR:HB	1.75	0.52
1:A:104:THR:HG23	1:A:113:SER:HB2	1.92	0.51
1:A:181:GLY:HA2	1:A:198:VAL:O	2.11	0.50
1:A:243:CYS:HA	3:A:504:NAG:H83	1.93	0.50
1:A:49:ILE:HD11	1:A:379:PHE:HE2	1.76	0.50
1:A:200:LEU:HD23	1:A:414:LEU:HD21	1.95	0.49
1:A:182:TYR:CZ	1:A:198:VAL:HG23	2.49	0.48
1:A:200:LEU:HD23	1:A:414:LEU:CD2	2.45	0.46
1:A:192:TYR:HB3	2:B:655:TRP:CD2	2.51	0.46
2:B:627:PHE:O	2:B:631:VAL:HG23	2.16	0.46
1:A:363:ASN:HB3	1:A:366:GLU:HB3	1.99	0.45
1:A:198:VAL:CG1	5:A:512:PLM:H41	2.46	0.44
1:A:337:ARG:CD	5:A:512:PLM:H31	2.43	0.44
1:A:215:PHE:HB3	8:A:632:HOH:O	2.17	0.44
1:A:270:ASP:O	5:A:512:PLM:HF2	2.18	0.44
1:A:363:ASN:H	1:A:367:HIS:CD2	2.35	0.43
1:A:271:ILE:HG22	1:A:341:ILE:HG21	2.00	0.43
1:A:337:ARG:HD2	5:A:512:PLM:H52	2.01	0.43
1:A:200:LEU:HD21	1:A:413:ILE:HG21	2.00	0.43
1:A:300:PHE:HB3	1:A:326:GLY:HA2	2.01	0.43
2:B:696:VAL:HG11	2:B:719:THR:HG23	2.01	0.42
1:A:254:PHE:HD1	6:A:517:GOL:H32	1.84	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:671:LEU:HG	2:B:684:ILE:HG23	2.02	0.42
2:B:580:MET:HB2	8:B:916:HOH:O	2.19	0.42
1:A:318:ILE:HG23	7:A:518:TRS:H22	2.01	0.41
1:A:140:LEU:HB3	5:A:512:PLM:H51	2.03	0.41
1:A:268:SER:HB3	1:A:271:ILE:HG12	2.03	0.41
1:A:369:THR:HG23	1:A:385:LYS:HA	2.02	0.41
1:A:275:ILE:HG12	1:A:276:TYR:H	1.86	0.41
1:A:174:THR:OG1	1:A:177:GLU:HG3	2.21	0.40
1:A:200:LEU:HD21	1:A:413:ILE:CG2	2.51	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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	398/472 (84%)	385 (97%)	13 (3%)	0	100	100
2	B	149/179 (83%)	146 (98%)	3 (2%)	0	100	100
All	All	547/651 (84%)	531 (97%)	16 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	363/419 (87%)	337 (93%)	26 (7%)	18	9
2	B	141/159 (89%)	130 (92%)	11 (8%)	16	7
All	All	504/578 (87%)	467 (93%)	37 (7%)	17	8

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	42	VAL
1	A	49	ILE
1	A	128	VAL
1	A	138	LEU
1	A	179	LEU
1	A	183	ARG
1	A	186	PHE
1	A	195	THR
1	A	200	LEU
1	A	218	LYS
1	A	226	ILE
1	A	262	GLN
1	A	268	SER
1	A	278	VAL
1	A	282	ASP
1	A	284	ASN
1	A	315	GLU
1	A	343	LEU
1	A	350	SER
1	A	354	SER
1	A	371	LEU
1	A	391	LEU
1	A	400	GLN
1	A	401	VAL
1	A	424	ASP
1	A	431	ARG
2	B	578	LYS
2	B	593	MET
2	B	612	LYS
2	B	623	LYS
2	B	633	GLN
2	B	651	ILE
2	B	661	LEU
2	B	671	LEU
2	B	716	LYS

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Mol	Chain	Res	Type
2	B	725	GLN
2	B	742	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	367	HIS
1	A	382	GLN
2	B	693	GLN
2	B	725	GLN
2	B	726	HIS

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](#)

20 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	NAG	A	501	1,3	14,14,15	0.33	0	15,19,21	0.78	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	NAG	A	502	3,4	14,14,15	0.32	0	15,19,21	0.99	1 (6%)
4	MAN	A	503	3	11,11,12	0.46	0	15,15,17	2.02	2 (13%)
3	NAG	A	504	1	14,14,15	0.28	0	15,19,21	0.66	0
3	NAG	A	505	1	14,14,15	0.30	0	15,19,21	1.25	1 (6%)
3	NAG	A	506	1	14,14,15	0.31	0	15,19,21	0.54	0
3	NAG	A	507	1	14,14,15	0.26	0	15,19,21	0.44	0
3	NAG	A	508	1	14,14,15	0.25	0	15,19,21	0.85	1 (6%)
3	NAG	A	509	1	14,14,15	0.30	0	15,19,21	0.63	0
3	NAG	A	510	1	14,14,15	0.30	0	15,19,21	0.73	1 (6%)
3	NAG	A	511	1	14,14,15	0.32	0	15,19,21	0.49	0
5	PLM	A	512	-	14,17,17	0.58	0	14,17,17	0.59	0
5	PLM	A	513	-	14,17,17	0.17	0	14,17,17	0.48	0
6	GOL	A	514	-	5,5,5	0.14	0	5,5,5	0.24	0
6	GOL	A	515	-	5,5,5	0.19	0	5,5,5	0.27	0
6	GOL	A	516	-	5,5,5	0.12	0	5,5,5	0.28	0
6	GOL	A	517	-	5,5,5	0.13	0	5,5,5	0.22	0
7	TRS	A	518	-	7,7,7	0.33	0	9,9,9	0.38	0
6	GOL	B	801	-	5,5,5	0.15	0	5,5,5	0.45	0
7	TRS	B	802	-	7,7,7	0.31	0	9,9,9	0.26	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	NAG	A	501	1,3	-	0/6/23/26	0/1/1/1
3	NAG	A	502	3,4	-	0/6/23/26	0/1/1/1
4	MAN	A	503	3	-	0/2/19/22	0/1/1/1
3	NAG	A	504	1	-	0/6/23/26	0/1/1/1
3	NAG	A	505	1	-	0/6/23/26	0/1/1/1
3	NAG	A	506	1	-	0/6/23/26	0/1/1/1
3	NAG	A	507	1	-	0/6/23/26	0/1/1/1
3	NAG	A	508	1	-	0/6/23/26	0/1/1/1
3	NAG	A	509	1	-	0/6/23/26	0/1/1/1
3	NAG	A	510	1	-	0/6/23/26	0/1/1/1
3	NAG	A	511	1	-	0/6/23/26	0/1/1/1
5	PLM	A	512	-	-	0/13/15/15	0/0/0/0
5	PLM	A	513	-	-	0/13/15/15	0/0/0/0
6	GOL	A	514	-	-	0/4/4/4	0/0/0/0
6	GOL	A	515	-	-	0/4/4/4	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	GOL	A	516	-	-	0/4/4/4	0/0/0/0
6	GOL	A	517	-	-	0/4/4/4	0/0/0/0
7	TRS	A	518	-	-	0/9/9/9	0/0/0/0
6	GOL	B	801	-	-	0/4/4/4	0/0/0/0
7	TRS	B	802	-	-	0/9/9/9	0/0/0/0

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
4	A	503	MAN	C1-C2-C3	2.05	112.03	109.55
3	A	510	NAG	C1-O5-C5	2.58	115.94	112.14
3	A	508	NAG	C1-O5-C5	2.99	116.53	112.14
3	A	502	NAG	C1-O5-C5	3.48	117.25	112.14
3	A	505	NAG	C1-O5-C5	4.39	118.60	112.14
4	A	503	MAN	C1-O5-C5	7.04	122.50	112.14

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 14 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	504	NAG	1	0
5	A	512	PLM	10	0
6	A	517	GOL	1	0
7	A	518	TRS	2	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	400/472 (84%)	-0.12	5 (1%) 79 81	25, 42, 72, 105	0
2	B	151/179 (84%)	0.17	6 (3%) 42 46	29, 51, 94, 115	0
All	All	551/651 (84%)	-0.04	11 (1%) 68 71	25, 44, 81, 115	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	579	ILE	4.8
2	B	612	LYS	4.5
1	A	35	GLN	3.7
1	A	36	LYS	3.1
1	A	58	GLY	2.5
2	B	736	LYS	2.5
2	B	606	ARG	2.5
2	B	611	ASP	2.3
1	A	429	MET	2.2
1	A	107	ALA	2.2
2	B	577	ASP	2.1

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
7	TRS	A	518	8/8	0.86	0.30	13.56	84,84,86,87	0
7	TRS	B	802	8/8	0.91	0.23	10.42	67,71,75,76	0
6	GOL	A	515	6/6	0.76	0.23	7.02	52,68,70,72	0
5	PLM	A	512	18/18	0.81	0.22	5.45	33,39,59,61	0
6	GOL	A	517	6/6	0.92	0.23	5.08	60,65,69,71	0
5	PLM	A	513	18/18	0.71	0.24	3.98	53,61,72,75	0
6	GOL	B	801	6/6	0.95	0.19	2.46	39,45,53,59	0
3	NAG	A	507	14/15	0.96	0.14	0.26	33,39,43,46	0
3	NAG	A	506	14/15	0.92	0.12	0.06	49,61,72,77	0
3	NAG	A	505	14/15	0.88	0.15	0.02	50,55,64,68	0
3	NAG	A	504	14/15	0.94	0.10	-0.29	30,42,46,48	0
3	NAG	A	501	14/15	0.95	0.09	-0.81	39,42,46,50	0
6	GOL	A	514	6/6	0.94	0.11	-0.97	34,51,57,58	0
3	NAG	A	508	14/15	0.88	0.17	-	54,66,75,76	0
6	GOL	A	516	6/6	0.84	0.18	-	60,64,66,66	0
3	NAG	A	510	14/15	0.78	0.35	-	79,86,91,92	0
3	NAG	A	511	14/15	0.79	0.21	-	91,95,99,100	0
3	NAG	A	509	14/15	0.67	0.36	-	97,101,106,108	0
3	NAG	A	502	14/15	0.92	0.10	-	47,55,62,70	0
4	MAN	A	503	11/12	0.79	0.14	-	78,87,90,95	0

6.5 Other polymers [i](#)

There are no such residues in this entry.