



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 10:27 PM GMT

PDB ID : 4XUP
Title : Structure of the N-terminal CBM22-1-CBM22-2 tandem domain from *Paenibacillus barcinonensis* Xyn10C
Authors : Sainz-Polo, M.A.; Sanz-Aparicio, J.
Deposited on : 2015-01-26
Resolution : 2.43 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

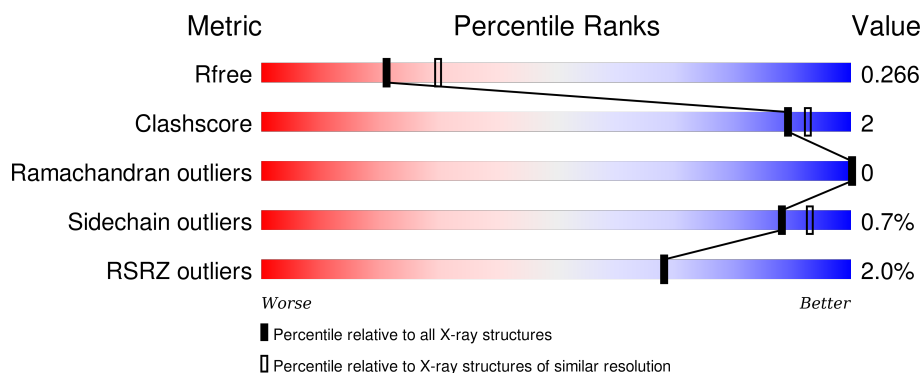
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.43 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	1003 (2.46-2.42)
Clashscore	102246	1071 (2.46-2.42)
Ramachandran outliers	100387	1065 (2.46-2.42)
Sidechain outliers	100360	1065 (2.46-2.42)
RSRZ outliers	91569	1005 (2.46-2.42)

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	334	<div> <div>2%</div> <div>92%</div> <div>7% ..</div> </div>
1	B	334	<div> <div>2%</div> <div>90%</div> <div>6% .</div> </div>
1	C	334	<div> <div>%</div> <div>44%</div> <div>52%</div> </div>
1	D	334	<div> <div>2%</div> <div>92%</div> <div>. .</div> </div>
1	E	334	<div> <div>46%</div> <div>52%</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	334	<div><div></div><div>2%</div><div>93%</div><div>5% •</div></div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 12540 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 Endo-1,4-beta-xylanase C.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	330	Total	C	N	O	S	0	0	0
			2549	1595	431	520	3			
1	B	321	Total	C	N	O	S	0	0	0
			2489	1561	420	505	3			
1	C	159	Total	C	N	O		0	0	0
			1221	762	208	251				
1	D	321	Total	C	N	O	S	0	0	0
			2489	1561	421	504	3			
1	E	159	Total	C	N	O		0	0	0
			1223	763	209	251				
1	F	325	Total	C	N	O	S	0	0	0
			2517	1577	426	511	3			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	87	ALA	PRO	conflict	UNP O69230
A	98	GLY	ARG	conflict	UNP O69230
A	177	ASP	TYR	conflict	UNP O69230
B	87	ALA	PRO	conflict	UNP O69230
B	98	GLY	ARG	conflict	UNP O69230
B	177	ASP	TYR	conflict	UNP O69230
C	87	ALA	PRO	conflict	UNP O69230
C	98	GLY	ARG	conflict	UNP O69230
C	177	ASP	TYR	conflict	UNP O69230
D	87	ALA	PRO	conflict	UNP O69230
D	98	GLY	ARG	conflict	UNP O69230
D	177	ASP	TYR	conflict	UNP O69230
E	87	ALA	PRO	conflict	UNP O69230
E	98	GLY	ARG	conflict	UNP O69230
E	177	ASP	TYR	conflict	UNP O69230
F	87	ALA	PRO	conflict	UNP O69230
F	98	GLY	ARG	conflict	UNP O69230

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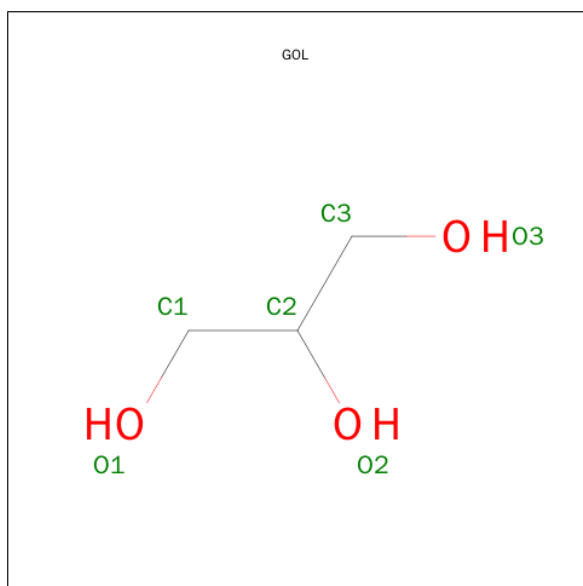
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Chain	Residue	Modelled	Actual	Comment	Reference
F	177	ASP	TYR	conflict	UNP O69230

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	D	2	Total 2 Ca 2	0	0
2	E	1	Total 1 Ca 1	0	0
2	B	2	Total 2 Ca 2	0	0
2	C	1	Total 1 Ca 1	0	0
2	A	2	Total 2 Ca 2	0	0
2	F	2	Total 2 Ca 2	0	0

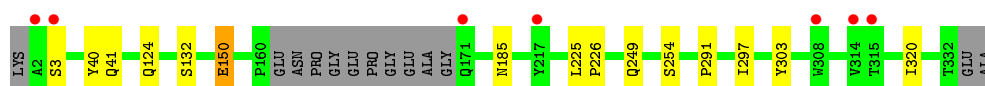
- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



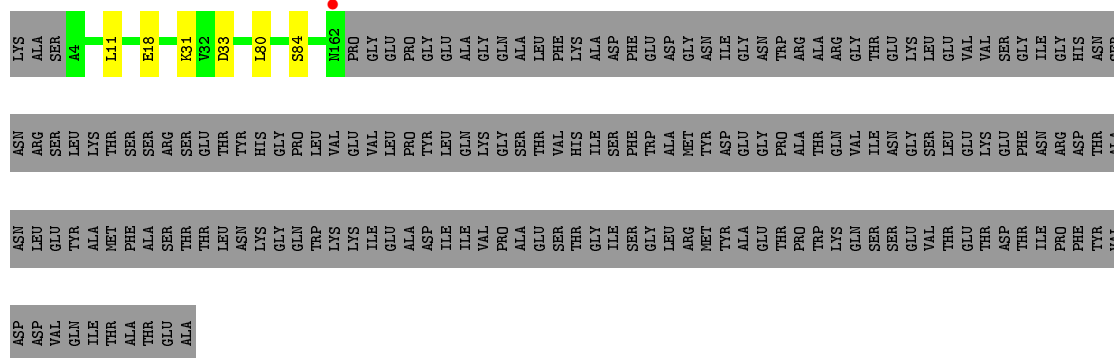
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total 6 C 3 O 3	0	0
3	D	1	Total 6 C 3 O 3	0	0

- Molecule 4 is water.

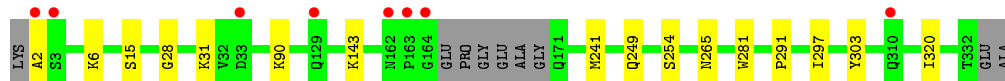
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	8	Total 8	O 8	0	0
4	B	6	Total 6	O 6	0	0
4	C	2	Total 2	O 2	0	0
4	D	8	Total 8	O 8	0	0
4	E	2	Total 2	O 2	0	0
4	F	4	Total 4	O 4	0	0



- Molecule 1: Endo-1,4-beta-xylanase C



- Molecule 1: Endo-1,4-beta-xylanase C



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	84.17Å 110.36Å 118.52Å 90.00° 90.63° 90.00°	Depositor
Resolution (Å)	118.52 – 2.43 84.16 – 2.43	Depositor EDS
% Data completeness (in resolution range)	99.3 (118.52-2.43) 99.3 (84.16-2.43)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.14 (at 2.42Å)	Xtriage
Refinement program	REFMAC 5.8.0103	Depositor
R, R_{free}	0.227 , 0.265 0.231 , 0.266	Depositor DCC
R_{free} test set	4072 reflections (5.29%)	DCC
Wilson B-factor (Å ²)	43.0	Xtriage
Anisotropy	0.390	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 25.1	EDS
Estimated twinning fraction	0.029 for h,-k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 81095 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	12540	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.02% 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: GOL, CA

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	0.72	2/2598 (0.1%)	0.84	0/3526
1	B	0.69	0/2537	0.84	0/3443
1	C	0.71	0/1241	0.89	2/1683 (0.1%)
1	D	0.70	1/2537 (0.0%)	0.85	0/3443
1	E	0.78	0/1243	0.86	0/1686
1	F	0.69	0/2566	0.83	0/3483
All	All	0.71	3/12722 (0.0%)	0.85	2/17264 (0.0%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	161	GLU	CD-OE1	5.39	1.31	1.25
1	A	161	GLU	CD-OE2	5.13	1.31	1.25
1	D	150	GLU	CD-OE2	-5.04	1.20	1.25

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	90	LYS	N-CA-CB	-5.17	101.29	110.60
1	C	90	LYS	CA-CB-CG	5.06	124.54	113.40

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	2549	0	2455	23	0
1	B	2489	0	2406	15	0
1	C	1221	0	1184	9	0
1	D	2489	0	2407	13	0
1	E	1223	0	1184	6	0
1	F	2517	0	2430	11	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	1	0	0	0	0
2	D	2	0	0	1	0
2	E	1	0	0	0	0
2	F	2	0	0	0	0
3	B	6	0	8	0	0
3	D	6	0	8	1	0
4	A	8	0	0	0	0
4	B	6	0	0	1	0
4	C	2	0	0	0	0
4	D	8	0	0	2	0
4	E	2	0	0	0	0
4	F	4	0	0	1	0
All	All	12540	0	12082	60	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 (60) 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:150:GLU:OE2	4:D:501:HOH:O	1.89	0.91
1:A:108:ASN:HD22	1:D:41:GLN:NE2	1.84	0.76
1:B:41:GLN:HB3	4:B:501:HOH:O	1.87	0.75
1:A:216:THR:HG21	1:A:308:TRP:HA	1.66	0.74
1:B:183:ILE:HG13	1:B:196:VAL:HG23	1.73	0.70
1:C:63:THR:HG22	1:C:132:SER:HA	1.76	0.66
1:E:18:GLU:HG3	1:F:31:LYS:HD3	1.79	0.64
2:D:401:CA:CA	4:D:501:HOH:O	1.74	0.64
1:A:170:GLY:O	1:A:331:ALA:HB3	1.97	0.63
1:B:277:ASN:HB2	1:B:280:GLN:HG2	1.82	0.62
1:A:132:SER:C	1:A:133:LEU:HD12	2.21	0.61
1:B:183:ILE:HG23	1:B:186:TRP:HB2	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:80:LEU:HD13	1:C:113:VAL:HG12	1.86	0.57
1:A:122:GLN:HG2	1:D:40:TYR:CZ	2.44	0.52
1:E:80:LEU:HB3	1:E:84:SER:OG	2.09	0.52
1:A:122:GLN:H	1:D:41:GLN:HG3	1.76	0.51
1:A:170:GLY:O	1:A:171:GLN:HB3	2.11	0.50
1:D:124:GLN:O	3:D:403:GOL:H2	2.11	0.50
1:B:63:THR:O	1:C:96:ARG:NH1	2.44	0.49
1:E:11:LEU:O	1:F:28:GLY:HA2	2.13	0.49
1:F:249:GLN:HG2	1:F:320:ILE:HD11	1.95	0.48
1:C:81:VAL:HG11	1:C:145:ALA:HB3	1.96	0.48
1:C:31:LYS:NZ	1:C:33:ASP:OD2	2.43	0.48
1:B:249:GLN:HG2	1:B:320:ILE:HD11	1.95	0.47
1:B:96:ARG:HG2	1:B:132:SER:OG	2.13	0.47
1:D:249:GLN:HG2	1:D:320:ILE:HD11	1.94	0.47
1:A:169:ALA:HB2	1:D:185:ASN:OD1	2.13	0.47
1:B:99:ASN:OD1	1:B:99:ASN:N	2.48	0.47
1:A:249:GLN:HG2	1:A:320:ILE:HD11	1.97	0.47
1:B:206:ARG:HH22	1:C:160:PRO:HB2	1.78	0.47
1:C:81:VAL:CG1	1:C:145:ALA:HB3	2.46	0.46
1:A:131:SER:O	1:B:96:ARG:NH2	2.45	0.45
1:A:244:GLU:HG2	1:A:245:GLY:N	2.31	0.45
1:A:170:GLY:O	1:A:171:GLN:CB	2.64	0.45
1:A:131:SER:O	1:B:96:ARG:NH1	2.49	0.45
1:F:15:SER:HB2	4:F:501:HOH:O	2.17	0.45
1:A:108:ASN:HD22	1:D:41:GLN:HE21	1.60	0.45
1:A:96:ARG:NH2	1:C:66:VAL:O	2.51	0.44
1:A:216:THR:CG2	1:A:308:TRP:HA	2.43	0.44
1:A:132:SER:O	1:A:133:LEU:HD12	2.17	0.44
1:A:216:THR:HG21	1:A:309:LYS:H	1.81	0.44
1:E:11:LEU:O	1:F:28:GLY:CA	2.66	0.43
1:E:31:LYS:NZ	1:E:33:ASP:OD2	2.46	0.43
1:A:122:GLN:N	1:D:41:GLN:HG3	2.34	0.43
1:F:254:SER:HB2	1:F:303:TYR:CZ	2.52	0.43
1:F:143:LYS:HE3	1:F:143:LYS:HB3	1.90	0.43
1:D:225:LEU:N	1:D:226:PRO:CD	2.81	0.43
1:B:291:PRO:HG2	1:B:297:ILE:HD13	2.01	0.43
1:A:108:ASN:HD22	1:D:41:GLN:HE22	1.63	0.43
1:B:254:SER:HB2	1:B:303:TYR:CZ	2.55	0.42
1:E:18:GLU:CG	1:F:31:LYS:HD3	2.48	0.42
1:D:291:PRO:HG2	1:D:297:ILE:HD13	2.00	0.42
1:F:291:PRO:HG2	1:F:297:ILE:HD13	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:64:ASP:HB3	1:C:134:LEU:HD22	2.01	0.42
1:B:225:LEU:N	1:B:226:PRO:CD	2.82	0.42
1:A:254:SER:HB2	1:A:303:TYR:CZ	2.55	0.41
1:D:254:SER:HB2	1:D:303:TYR:CZ	2.55	0.41
1:F:241:MET:HB2	1:F:281:TRP:CD2	2.56	0.41
1:A:291:PRO:HG2	1:A:297:ILE:HD13	2.02	0.40
1:A:261:ARG:HB2	1:F:2:ALA:HB3	2.03	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	326/334 (98%)	321 (98%)	5 (2%)	0	100	100
1	B	317/334 (95%)	310 (98%)	7 (2%)	0	100	100
1	C	157/334 (47%)	152 (97%)	5 (3%)	0	100	100
1	D	317/334 (95%)	312 (98%)	5 (2%)	0	100	100
1	E	157/334 (47%)	151 (96%)	6 (4%)	0	100	100
1	F	321/334 (96%)	315 (98%)	6 (2%)	0	100	100
All	All	1595/2004 (80%)	1561 (98%)	34 (2%)	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	270/273 (99%)	267 (99%)	3 (1%)	80	87
1	B	265/273 (97%)	264 (100%)	1 (0%)	93	96
1	C	128/273 (47%)	128 (100%)	0	100	100
1	D	265/273 (97%)	263 (99%)	2 (1%)	86	92
1	E	128/273 (47%)	128 (100%)	0	100	100
1	F	268/273 (98%)	265 (99%)	3 (1%)	80	87
All	All	1324/1638 (81%)	1315 (99%)	9 (1%)	88	93

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	129	GLN
1	A	216	THR
1	A	244	GLU
1	B	99	ASN
1	D	3	SER
1	D	132	SER
1	F	6	LYS
1	F	90	LYS
1	F	265	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	116	GLN
1	A	122	GLN
1	A	129	GLN
1	A	203	ASN
1	A	277	ASN
1	B	116	GLN
1	B	185	ASN
1	B	203	ASN
1	C	122	GLN
1	D	41	GLN
1	D	122	GLN
1	D	203	ASN
1	D	277	ASN
1	E	56	ASN

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Mol	Chain	Res	Type
1	E	106	GLN
1	F	116	GLN
1	F	122	GLN
1	F	129	GLN
1	F	185	ASN
1	F	203	ASN
1	F	265	ASN

5.3.3 RNA ⓘ

There are no RNA molecules 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 ⓘ

Of 12 ligands modelled in this entry, 10 are monoatomic - leaving 2 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	$\# Z > 2$
3	GOL	B	403	-	5,5,5	0.77	0	5,5,5	1.08	0
3	GOL	D	403	-	5,5,5	0.95	0	5,5,5	0.88	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	GOL	B	403	-	-	0/4/4/4	0/0/0/0
3	GOL	D	403	-	-	0/4/4/4	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	403	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 ⓘ

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	330/334 (98%)	0.20	8 (2%) 62 61	23, 47, 75, 84	0
1	B	321/334 (96%)	0.31	7 (2%) 65 65	26, 45, 70, 86	0
1	C	159/334 (47%)	0.18	2 (1%) 79 80	28, 47, 75, 84	0
1	D	321/334 (96%)	0.33	7 (2%) 65 65	23, 48, 69, 83	0
1	E	159/334 (47%)	0.05	1 (0%) 90 91	20, 38, 67, 77	0
1	F	325/334 (97%)	0.29	8 (2%) 61 59	31, 55, 74, 84	0
All	All	1615/2004 (80%)	0.25	33 (2%) 68 68	20, 48, 73, 86	0

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	314	VAL	5.4
1	E	162	ASN	4.2
1	D	171	GLN	4.1
1	F	164	GLY	4.0
1	F	163	PRO	3.9
1	A	171	GLN	3.9
1	A	169	ALA	3.7
1	F	162	ASN	3.7
1	D	315	THR	3.4
1	D	2	ALA	3.2
1	F	3	SER	3.2
1	B	2	ALA	3.0
1	A	2	ALA	2.9
1	A	168	GLU	2.9
1	A	131	SER	2.8
1	A	3	SER	2.8
1	A	170	GLY	2.8
1	D	308	TRP	2.7
1	C	3	SER	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	3	SER	2.5
1	D	217	TYR	2.5
1	D	3	SER	2.4
1	B	308	TRP	2.4
1	B	214	SER	2.3
1	F	129	GLN	2.2
1	F	33	ASP	2.2
1	C	109	ALA	2.2
1	B	279	GLY	2.1
1	F	2	ALA	2.1
1	F	310	GLN	2.0
1	B	106	GLN	2.0
1	B	161	GLU	2.0
1	A	253	GLY	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	CA	A	1002	1/1	0.98	0.14	-0.09	55,55,55,55	0
3	GOL	D	403	6/6	0.80	0.15	-0.24	44,50,51,52	0
2	CA	F	401	1/1	0.99	0.14	-0.88	36,36,36,36	0
2	CA	E	401	1/1	1.00	0.13	-1.39	26,26,26,26	0
2	CA	A	1001	1/1	0.99	0.12	-1.54	38,38,38,38	0
2	CA	B	401	1/1	0.98	0.14	-1.55	42,42,42,42	0
2	CA	D	401	1/1	0.92	0.13	-1.75	36,36,36,36	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	CA	C	401	1/1	0.98	0.09	-2.46	43,43,43,43	0
2	CA	D	402	1/1	0.98	0.08	-3.80	40,40,40,40	0
3	GOL	B	403	6/6	0.86	0.13	-3.92	48,53,56,57	0
2	CA	F	402	1/1	0.95	0.07	-3.94	50,50,50,50	0
2	CA	B	402	1/1	0.99	0.05	-6.24	42,42,42,42	0

6.5 Other polymers [i](#)

There are no such residues in this entry.