



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 03:42 PM GMT

PDB ID : 4GWM
Title : Crystal structure of human promeprin beta
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Stocker, W.; Becker-Pauly, C.; Gomis-Ruth, F.X.
Deposited on : 2012-09-03
Resolution : 1.85 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

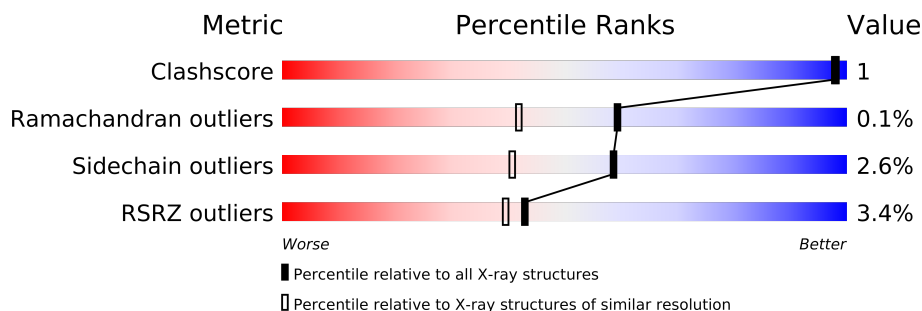
The following versions of software and data (see [references](#)) were used in the production of this report:

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 1.85 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))
Clashscore	79885	1470 (1.86-1.86)
Ramachandran outliers	78287	1451 (1.86-1.86)
Sidechain outliers	78261	1451 (1.86-1.86)
RSRZ outliers	66119	1269 (1.86-1.86)

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	592	
1	B	592	

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

Mol	Type	Chain	Res	Geometry	Electron density
8	NAG	A	727	-	X

2 Entry composition

There are 14 unique types of molecules in this entry. The entry contains 10380 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 Meprin A subunit beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	561	Total	C	N	O	S	0	0	0
			4508	2834	782	872	20			
1	B	554	Total	C	N	O	S	0	0	0
			4458	2810	774	854	20			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	23	PRO	THR	ENGINEERED MUTATION	UNP Q16820
A	24	TRP	PRO	ENGINEERED MUTATION	UNP Q16820
B	23	PRO	THR	ENGINEERED MUTATION	UNP Q16820
B	24	TRP	PRO	ENGINEERED MUTATION	UNP Q16820

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Zn	0	0
			1	1		
2	A	1	Total	Zn	0	0
			1	1		

- Molecule 3 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Na	0	0
			1	1		
3	A	1	Total	Na	0	0
			1	1		

- Molecule 4 is a polymer of unknown type called SUGAR (3-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	3	Total	C	N	O	0	0
			39	22	2	15		
4	A	3	Total	C	N	O	0	0
			39	22	2	15		

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	23	PRO	THR	ENGINEERED MUTATION	UNP Q16820
A	24	TRP	PRO	ENGINEERED MUTATION	UNP Q16820
A	23	PRO	THR	ENGINEERED MUTATION	UNP Q16820
A	24	TRP	PRO	ENGINEERED MUTATION	UNP Q16820

- Molecule 5 is a polymer of unknown type called SUGAR (7-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	7	Total	C	N	O	0	0
			81	46	2	33		

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	23	PRO	THR	ENGINEERED MUTATION	UNP Q16820
A	24	TRP	PRO	ENGINEERED MUTATION	UNP Q16820

- Molecule 6 is a polymer of unknown type called SUGAR (2-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	2	Total	C	N	O	0	0
			24	14	1	9		
6	B	2	Total	C	N	O	0	0
			24	14	1	9		

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	23	PRO	THR	ENGINEERED MUTATION	UNP Q16820
A	24	TRP	PRO	ENGINEERED MUTATION	UNP Q16820
B	23	PRO	THR	ENGINEERED MUTATION	UNP Q16820
B	24	TRP	PRO	ENGINEERED MUTATION	UNP Q16820

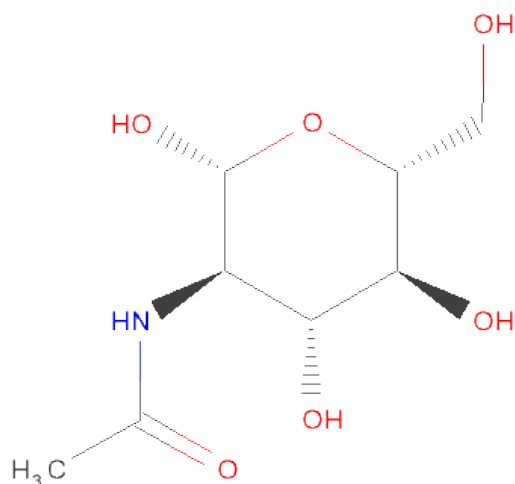
- Molecule 7 is a polymer of unknown type called SUGAR (9-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	A	9	Total	C	N	O	0	0
			105	58	2	45		
7	B	9	Total	C	N	O	0	0
			105	58	2	45		

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	23	PRO	THR	ENGINEERED MUTATION	UNP Q16820
A	24	TRP	PRO	ENGINEERED MUTATION	UNP Q16820
B	23	PRO	THR	ENGINEERED MUTATION	UNP Q16820
B	24	TRP	PRO	ENGINEERED MUTATION	UNP Q16820

- Molecule 8 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



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

- Molecule 9 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

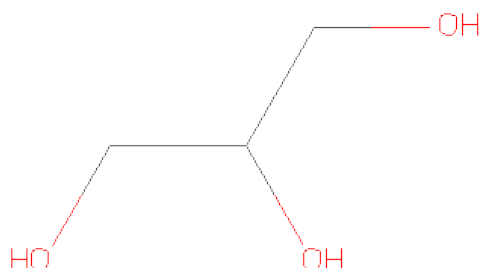
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	B	3	Total	Cl	0	0
			3	3		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	2	Total	Cl	0	0
			2	2		

- Molecule 10 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



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

- Molecule 11 is a polymer of unknown type called SUGAR (2-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
11	B	2	Total	C	N	O	0	0
			28	16	2	10		

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	23	PRO	THR	ENGINEERED MUTATION	UNP Q16820
B	24	TRP	PRO	ENGINEERED MUTATION	UNP Q16820

- Molecule 12 is a polymer of unknown type called SUGAR (6-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
12	B	6	Total	C	N	O	0	0
			70	40	2	28		

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	23	PRO	THR	ENGINEERED MUTATION	UNP Q16820
B	24	TRP	PRO	ENGINEERED MUTATION	UNP Q16820

- Molecule 13 is a polymer of unknown type called SUGAR (4-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
13	B	4	Total	C	N	O	0	0
			48	28	2	18		

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	23	PRO	THR	ENGINEERED MUTATION	UNP Q16820
B	24	TRP	PRO	ENGINEERED MUTATION	UNP Q16820

- Molecule 14 is water.

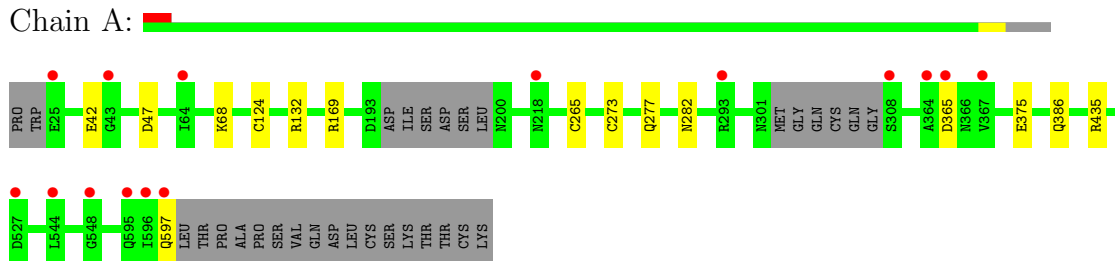
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
14	A	352	Total	O	0	0
			352	352		
14	B	456	Total	O	0	0
			456	456		

3 Residue-property plots

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.

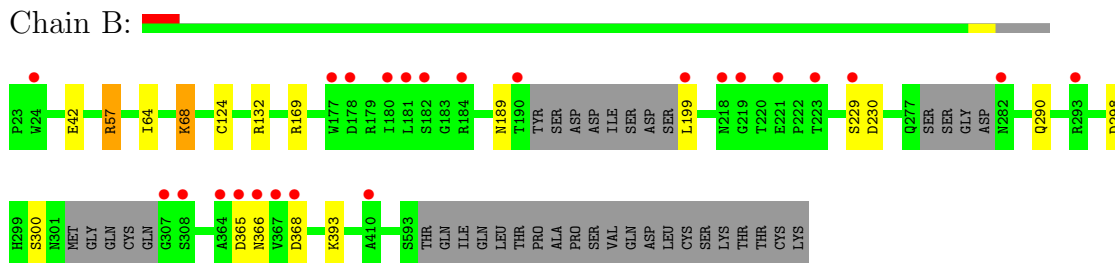
- Molecule 1: Meprin A subunit beta

Chain A:



- Molecule 1: Meprin A subunit beta

Chain B:



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	69.62Å 71.12Å 85.74Å 74.87° 80.08° 65.13°	Depositor
Resolution (Å)	48.45 – 1.85 48.44 – 1.85	Depositor EDS
% Data completeness (in resolution range)	96.6 (48.45-1.85) 96.6 (48.44-1.85)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.95 (at 1.86Å)	Xtriage
Refinement program	BUSTER 2.11.2	Depositor
R, R_{free}	0.168 , 0.188 0.173 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	25.4	Xtriage
Anisotropy	0.128	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 46.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 118474 reflections	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	10380	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.40% 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, ZN, BMA, NAG, CL, NA, FUC, MAN

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	0.51	0/4617	0.62	0/6252
1	B	0.58	0/4568	0.63	0/6185
All	All	0.54	0/9185	0.62	0/12437

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	4508	0	0	3	0
1	B	4458	0	0	1	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	78	0	0	0	0
5	A	81	0	0	0	0
6	A	24	0	0	0	0
6	B	24	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	A	105	0	0	0	0
7	B	105	0	0	0	0
8	A	14	0	0	0	0
8	B	14	0	0	0	0
9	A	2	0	0	0	0
9	B	3	0	0	0	0
10	A	6	0	0	0	0
11	B	28	0	0	0	0
12	B	70	0	0	0	0
13	B	48	0	0	0	0
14	A	352	0	0	0	0
14	B	456	0	0	0	0
All	All	10380	0	0	4	0

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 1.

All (4) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:265:CYS:SG	1:A:273:CYS:SG	3.05	0.55
1:B:68:LYS:O	1:B:132:ARG:NE	2.53	0.42
1:A:386:GLN:OE1	1:A:435:ARG:NH2	2.54	0.40
1:A:68:LYS:O	1:A:132:ARG:NE	2.55	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	555/592 (94%)	543 (98%)	12 (2%)	0	100	100
1	B	546/592 (92%)	534 (98%)	11 (2%)	1 (0%)	56	38
All	All	1101/1184 (93%)	1077 (98%)	23 (2%)	1 (0%)	59	42

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	57	ARG

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	497/525 (95%)	488 (98%)	9 (2%)	71	57
1	B	490/525 (93%)	473 (96%)	17 (4%)	48	26
All	All	987/1050 (94%)	961 (97%)	26 (3%)	59	41

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	42	GLU
1	A	47	ASP
1	A	124	CYS
1	A	169	ARG
1	A	277	GLN
1	A	282	ASN
1	A	365	ASP
1	A	375	GLU
1	A	597	GLN
1	B	42	GLU
1	B	57	ARG
1	B	64	ILE
1	B	68	LYS
1	B	124	CYS
1	B	169	ARG
1	B	189	ASN
1	B	199	LEU
1	B	229	SER
1	B	230	ASP
1	B	290	GLN
1	B	298	ASP
1	B	300	SER
1	B	365	ASP

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Mol	Chain	Res	Type
1	B	366	ASN
1	B	368	ASP
1	B	393	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 ⓘ

47 carbohydrates 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	$\# Z > 2$
4	NAG	A	703	1,4	12,14,15	0.27	0	15,19,21	0.50	0
4	NAG	A	704	4	12,14,15	0.28	0	15,19,21	0.63	0
4	BMA	A	705	4	10,11,12	0.37	0	11,15,17	0.42	0
5	NAG	A	706	1,5	12,14,15	0.34	0	15,19,21	0.52	0
5	FUC	A	707	5	9,10,11	0.35	0	10,14,16	0.33	0
5	FUC	A	708	5	9,10,11	0.43	0	10,14,16	0.62	0
5	NAG	A	709	5	12,14,15	0.32	0	15,19,21	0.47	0
5	BMA	A	710	5	10,11,12	0.37	0	11,15,17	0.37	0
5	MAN	A	711	5	10,11,12	0.32	0	11,15,17	0.66	0
5	MAN	A	712	5	10,11,12	0.38	0	11,15,17	0.48	0
6	NAG	A	713	1,6	12,14,15	0.32	0	15,19,21	0.51	0
6	FUC	A	714	6	9,10,11	0.36	0	10,14,16	0.41	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
7	NAG	A	715	1,7	12,14,15	0.28	0	15,19,21	0.48	0
7	NAG	A	716	7	12,14,15	0.33	0	15,19,21	0.54	0
7	BMA	A	717	7	10,11,12	0.39	0	11,15,17	0.70	0
7	MAN	A	718	7	10,11,12	0.36	0	11,15,17	0.49	0
7	MAN	A	719	7	10,11,12	0.37	0	11,15,17	0.51	0
7	MAN	A	720	7	10,11,12	0.35	0	11,15,17	0.43	0
7	MAN	A	721	7	10,11,12	0.37	0	11,15,17	0.37	0
7	MAN	A	722	7	10,11,12	0.36	0	11,15,17	0.35	0
7	MAN	A	723	7	10,11,12	0.36	0	11,15,17	0.46	0
4	NAG	A	724	1,4	12,14,15	0.28	0	15,19,21	0.50	0
4	NAG	A	725	4	12,14,15	0.27	0	15,19,21	0.45	0
4	BMA	A	726	4	10,11,12	0.37	0	11,15,17	0.37	0
11	NAG	B	703	11,1	12,14,15	0.32	0	15,19,21	0.53	0
11	NAG	B	704	11	12,14,15	0.26	0	15,19,21	0.40	0
12	NAG	B	705	1,12	12,14,15	0.30	0	15,19,21	0.56	0
12	FUC	B	706	12	9,10,11	0.36	0	10,14,16	0.40	0
12	FUC	B	707	12	9,10,11	0.37	0	10,14,16	0.48	0
12	NAG	B	708	12	12,14,15	0.29	0	15,19,21	0.44	0
12	BMA	B	709	12	10,11,12	0.38	0	11,15,17	0.43	0
12	MAN	B	710	12	10,11,12	0.32	0	11,15,17	0.78	1 (9%)
13	NAG	B	711	1,13	12,14,15	0.34	0	15,19,21	0.57	0
13	FUC	B	712	13	9,10,11	0.36	0	10,14,16	0.46	0
13	FUC	B	713	13	9,10,11	0.36	0	10,14,16	0.48	0
13	NAG	B	714	13	12,14,15	0.29	0	15,19,21	0.51	0
7	NAG	B	715	1,7	12,14,15	0.29	0	15,19,21	0.47	0
7	NAG	B	716	7	12,14,15	0.28	0	15,19,21	0.49	0
7	BMA	B	717	7	10,11,12	0.39	0	11,15,17	0.58	0
7	MAN	B	718	7	10,11,12	0.36	0	11,15,17	0.47	0
7	MAN	B	719	7	10,11,12	0.38	0	11,15,17	0.47	0
7	MAN	B	720	7	10,11,12	0.33	0	11,15,17	0.33	0
7	MAN	B	721	7	10,11,12	0.37	0	11,15,17	0.38	0
7	MAN	B	722	7	10,11,12	0.33	0	11,15,17	0.37	0
7	MAN	B	723	7	10,11,12	0.35	0	11,15,17	0.38	0
6	NAG	B	725	1,6	12,14,15	0.23	0	15,19,21	0.37	0
6	FUC	B	726	6	9,10,11	0.36	0	10,14,16	0.45	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
4	NAG	A	703	1,4	-	0/6/23/26	0/1/1/1
4	NAG	A	704	4	-	0/6/23/26	0/1/1/1
4	BMA	A	705	4	-	0/2/19/22	0/1/1/1
5	NAG	A	706	1,5	-	0/6/23/26	0/1/1/1
5	FUC	A	707	5	-	0/0/17/20	0/1/1/1
5	FUC	A	708	5	-	0/0/17/20	0/1/1/1
5	NAG	A	709	5	-	0/6/23/26	0/1/1/1
5	BMA	A	710	5	-	0/2/19/22	0/1/1/1
5	MAN	A	711	5	-	0/2/19/22	0/1/1/1
5	MAN	A	712	5	-	0/2/19/22	0/1/1/1
6	NAG	A	713	1,6	-	0/6/23/26	0/1/1/1
6	FUC	A	714	6	-	0/0/17/20	0/1/1/1
7	NAG	A	715	1,7	-	0/6/23/26	0/1/1/1
7	NAG	A	716	7	-	0/6/23/26	0/1/1/1
7	BMA	A	717	7	-	0/2/19/22	0/1/1/1
7	MAN	A	718	7	-	0/2/19/22	0/1/1/1
7	MAN	A	719	7	-	0/2/19/22	0/1/1/1
7	MAN	A	720	7	-	0/2/19/22	0/1/1/1
7	MAN	A	721	7	-	0/2/19/22	0/1/1/1
7	MAN	A	722	7	-	0/2/19/22	0/1/1/1
7	MAN	A	723	7	-	0/2/19/22	0/1/1/1
4	NAG	A	724	1,4	-	0/6/23/26	0/1/1/1
4	NAG	A	725	4	-	0/6/23/26	0/1/1/1
4	BMA	A	726	4	-	0/2/19/22	0/1/1/1
11	NAG	B	703	11,1	-	0/6/23/26	0/1/1/1
11	NAG	B	704	11	-	0/6/23/26	0/1/1/1
12	NAG	B	705	1,12	-	0/6/23/26	0/1/1/1
12	FUC	B	706	12	-	0/0/17/20	0/1/1/1
12	FUC	B	707	12	-	0/0/17/20	0/1/1/1
12	NAG	B	708	12	-	0/6/23/26	0/1/1/1
12	BMA	B	709	12	-	0/2/19/22	0/1/1/1
12	MAN	B	710	12	-	0/2/19/22	0/1/1/1
13	NAG	B	711	1,13	-	0/6/23/26	0/1/1/1
13	FUC	B	712	13	-	0/0/17/20	0/1/1/1
13	FUC	B	713	13	-	0/0/17/20	0/1/1/1
13	NAG	B	714	13	-	0/6/23/26	0/1/1/1
7	NAG	B	715	1,7	-	0/6/23/26	0/1/1/1
7	NAG	B	716	7	-	0/6/23/26	0/1/1/1
7	BMA	B	717	7	-	0/2/19/22	0/1/1/1
7	MAN	B	718	7	-	0/2/19/22	0/1/1/1
7	MAN	B	719	7	-	0/2/19/22	0/1/1/1
7	MAN	B	720	7	-	0/2/19/22	0/1/1/1
7	MAN	B	721	7	-	0/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
7	MAN	B	722	7	-	0/2/19/22	0/1/1/1
7	MAN	B	723	7	-	0/2/19/22	0/1/1/1
6	NAG	B	725	1,6	-	0/6/23/26	0/1/1/1
6	FUC	B	726	6	-	0/0/17/20	0/1/1/1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
12	B	710	MAN	O5-C5-C6	2.02	109.10	106.98

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.6 Ligand geometry ⓘ

Of 12 ligands modelled in this entry, 9 are monoatomic - leaving 3 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	$\# Z > 2$
8	NAG	A	727	1	12,14,15	0.24	0	15,19,21	0.37	0
10	GOL	A	730	-	5,5,5	0.14	0	5,5,5	0.32	0
8	NAG	B	724	1	12,14,15	0.25	0	15,19,21	0.38	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
8	NAG	A	727	1	-	0/6/23/26	0/1/1/1
10	GOL	A	730	-	-	0/4/4/4	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
8	NAG	B	724	1	-	0/6/23/26	0/1/1/1

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.

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	561/592 (94%)	-0.02	15 (2%)	52 48	21, 36, 66, 128	0
1	B	554/592 (93%)	0.03	24 (4%)	34 31	16, 30, 63, 90	0
All	All	1115/1184 (94%)	0.01	39 (3%)	43 39	16, 34, 65, 128	0

All (39) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	307	GLY	8.5
1	B	199	LEU	6.5
1	A	596	ILE	6.5
1	B	308	SER	5.6
1	A	597	GLN	5.5
1	B	364	ALA	5.5
1	B	178	ASP	5.2
1	B	365	ASP	4.3
1	B	181	LEU	4.3
1	A	43	GLY	3.7
1	A	293	ARG	3.7
1	B	366	ASN	3.7
1	A	527	ASP	3.5
1	B	368	ASP	3.5
1	B	190	THR	3.5
1	A	365	ASP	3.5
1	A	308	SER	3.4
1	B	24	TRP	3.3
1	B	177	TRP	3.2
1	B	367	VAL	3.1
1	A	364	ALA	3.1
1	B	180	ILE	3.1
1	B	182	SER	3.0
1	B	410	ALA	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	367	VAL	2.8
1	B	223	THR	2.6
1	B	184	ARG	2.5
1	B	219	GLY	2.5
1	A	218	ASN	2.5
1	B	293	ARG	2.4
1	A	544	LEU	2.3
1	A	64	ILE	2.3
1	A	548	GLY	2.3
1	B	229	SER	2.3
1	A	25	GLU	2.2
1	A	595	GLN	2.2
1	B	218	ASN	2.1
1	B	221	GLU	2.0
1	B	282	ASN	2.0

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 ⓘ

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
6	NAG	B	725	14/15	0.25	24.23	70,74,83,88	0
13	NAG	B	714	14/15	0.26	20.44	48,53,61,66	0
6	NAG	A	713	14/15	0.28	9.38	80,85,91,93	0
13	FUC	B	713	10/11	0.18	7.27	42,47,49,51	0
7	NAG	B	715	14/15	0.11	3.26	41,46,60,62	0
11	NAG	B	704	14/15	0.25	2.20	51,55,59,61	0
7	MAN	B	721	11/12	0.21	2.16	51,54,56,59	0
7	MAN	A	722	11/12	0.14	2.01	64,68,74,76	0
7	MAN	B	720	11/12	0.15	1.90	47,50,59,64	0
4	NAG	A	704	14/15	0.24	1.14	50,59,69,82	0
7	BMA	A	717	11/12	0.18	1.04	62,67,74,82	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
13	NAG	B	711	14/15	0.09	0.97	33,38,45,50	0
4	NAG	A	703	14/15	0.12	0.44	35,43,48,50	0
11	NAG	B	703	14/15	0.11	0.32	33,39,45,46	0
4	NAG	A	724	14/15	0.20	0.32	70,73,79,81	0
7	NAG	A	716	14/15	0.14	0.27	48,52,58,61	0
7	NAG	A	715	14/15	0.17	0.18	52,56,64,65	0
7	NAG	B	716	14/15	0.08	-0.10	30,41,53,55	0
12	FUC	B	707	10/11	0.09	-0.17	39,42,46,49	0
12	NAG	B	705	14/15	0.16	-0.39	43,57,64,66	0
7	MAN	A	723	11/12	0.12	-0.45	66,69,73,75	0
12	FUC	B	706	10/11	0.16	-0.46	59,62,69,70	0
5	NAG	A	706	14/15	0.07	-0.81	36,43,56,59	0
5	FUC	A	708	10/11	0.07	-0.95	36,38,41,48	0
7	MAN	A	719	11/12	0.09	-1.12	59,66,69,74	0
7	BMA	B	717	11/12	0.08	-3.57	47,55,62,70	0
7	MAN	B	719	11/12	0.06	-3.96	42,45,49,56	0
7	MAN	B	722	11/12	0.13	-	68,74,79,84	0
7	MAN	A	721	11/12	0.23	-	97,102,103,103	0
7	MAN	B	718	11/12	0.19	-	80,86,90,91	0
6	FUC	A	714	10/11	0.34	-	97,102,103,103	0
4	BMA	A	726	11/12	0.39	-	99,102,105,106	0
4	NAG	A	725	14/15	0.36	-	83,89,97,98	0
7	MAN	A	720	11/12	0.17	-	81,84,90,91	0
5	FUC	A	707	10/11	0.24	-	62,67,70,73	0
6	FUC	B	726	10/11	0.38	-	84,86,88,89	0
13	FUC	B	712	10/11	0.28	-	62,72,79,79	0
5	NAG	A	709	14/15	0.10	-	48,56,62,66	0
5	MAN	A	712	11/12	0.17	-	80,83,87,89	0
12	MAN	B	710	11/12	0.52	-	114,116,117,118	0
7	MAN	A	718	11/12	0.42	-	90,95,98,99	0
5	BMA	A	710	11/12	0.17	-	73,80,84,87	0
7	MAN	B	723	11/12	0.42	-	91,97,99,100	0
5	MAN	A	711	11/12	0.25	-	93,98,100,101	0
12	NAG	B	708	14/15	0.18	-	61,70,75,83	0
12	BMA	B	709	11/12	0.38	-	95,106,109,110	0
4	BMA	A	705	11/12	0.45	-	92,101,105,107	0

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
8	NAG	A	727	14/15	0.35	9.13	81,85,90,92	0
10	GOL	A	730	6/6	0.17	1.51	53,54,55,59	0
2	ZN	B	701	1/1	0.10	0.33	20,20,20,20	0
2	ZN	A	701	1/1	0.10	0.02	25,25,25,25	0
3	NA	B	702	1/1	0.09	-0.25	27,27,27,27	0
9	CL	A	729	1/1	0.10	-0.29	28,28,28,28	0
9	CL	B	728	1/1	0.07	-0.36	32,32,32,32	1
9	CL	B	729	1/1	0.08	-1.13	20,20,20,20	1
9	CL	B	727	1/1	0.11	-1.24	18,18,18,18	0
8	NAG	B	724	14/15	0.24	-1.40	84,89,92,94	0
9	CL	A	728	1/1	0.07	-2.05	28,28,28,28	1
3	NA	A	702	1/1	0.05	-2.30	31,31,31,31	0

6.5 Other polymers ⓘ

There are no such residues in this entry.