



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 31, 2016 – 05:08 PM EDT

PDB ID : 5AM8  
Title : Crystal structure of the Angiotensin-1 converting enzyme N-domain in complex with amyloid-beta 4-10  
Authors : Masuyer, G.; Larmuth, K.M.; Douglas, R.G.; Sturrock, E.D.; Acharya, K.R.  
Deposited on : 2015-03-10  
Resolution : 1.90 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](mailto: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.

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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-20027107  
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-20027107

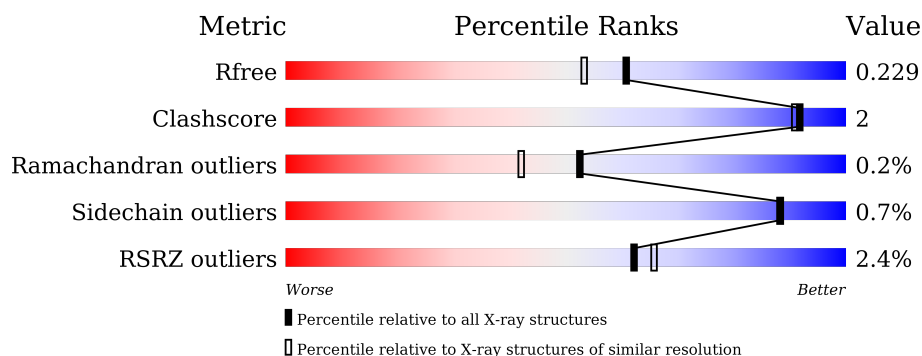
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.90 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	4755 (1.90-1.90)
Clashscore	102246	5398 (1.90-1.90)
Ramachandran outliers	100387	5338 (1.90-1.90)
Sidechain outliers	100360	5339 (1.90-1.90)
RSRZ outliers	91569	4766 (1.90-1.90)

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  $\geq 3$ , 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	629	<div> <div>2%</div> <div>91%</div> <div>5%</div> <div>.</div> </div>
1	B	629	<div> <div>3%</div> <div>91%</div> <div>.</div> <div>.</div> <div>.</div> </div>
1	C	629	<div> <div>%</div> <div>91%</div> <div>6%</div> <div>.</div> </div>
1	D	629	<div> <div>3%</div> <div>92%</div> <div>6%</div> <div>.</div> </div>
2	P	7	<div> <div>29%</div> <div>14%</div> <div>57%</div> </div>
2	Q	7	<div> <div>14%</div> <div>29%</div> <div>57%</div> </div>

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Mol	Chain	Length	Quality of chain
2	R	7	
2	S	7	

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	SO4	A	1003	-	-	-	X
5	SO4	C	1003	-	-	-	X
5	SO4	D	1003	-	-	-	X
6	NAG	A	1100	-	-	-	X
6	NAG	D	1101	-	-	-	X
7	NAG	A	1101	-	-	-	X
7	NAG	A	1103	-	-	-	X
7	NAG	B	1102	-	-	-	X
8	PEG	A	1200	-	-	-	X
8	PEG	A	1201	-	-	-	X
8	PEG	A	1203	-	-	-	X
8	PEG	B	1201	-	-	-	X
8	PEG	B	1202	-	-	-	X
8	PEG	B	1204	-	-	-	X
8	PEG	B	1205	-	-	-	X
8	PEG	C	1201	-	-	-	X
8	PEG	C	1202	-	-	-	X
8	PEG	C	1203	-	-	-	X
8	PEG	D	1203	-	-	-	X
8	PEG	D	1204	-	-	-	X
9	P6G	B	1203	-	-	-	X
9	P6G	D	1201	-	-	-	X

## 2 Entry composition

There are 14 unique types of molecules in this entry. The entry contains 22316 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 ANGIOTENSIN-CONVERTING ENZYME.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	604	Total	C	N	O	S	0	1	0
			4939	3173	849	898	19			
1	B	606	Total	C	N	O	S	0	1	0
			4950	3179	852	900	19			
1	C	606	Total	C	N	O	S	0	1	0
			4959	3186	852	902	19			
1	D	612	Total	C	N	O	S	0	5	0
			5029	3229	867	914	19			

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	9	GLN	ASN	ENGINEERED MUTATION	UNP P12821
A	25	GLN	ASN	ENGINEERED MUTATION	UNP P12821
A	82	GLN	ASN	ENGINEERED MUTATION	UNP P12821
A	117	GLN	ASN	ENGINEERED MUTATION	UNP P12821
A	289	GLN	ASN	ENGINEERED MUTATION	UNP P12821
A	545	ARG	GLN	ENGINEERED MUTATION	UNP P12821
A	576	LEU	PRO	ENGINEERED MUTATION	UNP P12821
A	629	LEU	ARG	ENGINEERED MUTATION	UNP P12821
B	9	GLN	ASN	ENGINEERED MUTATION	UNP P12821
B	25	GLN	ASN	ENGINEERED MUTATION	UNP P12821
B	82	GLN	ASN	ENGINEERED MUTATION	UNP P12821
B	117	GLN	ASN	ENGINEERED MUTATION	UNP P12821
B	289	GLN	ASN	ENGINEERED MUTATION	UNP P12821
B	545	ARG	GLN	ENGINEERED MUTATION	UNP P12821
B	576	LEU	PRO	ENGINEERED MUTATION	UNP P12821
B	629	LEU	ARG	ENGINEERED MUTATION	UNP P12821
C	9	GLN	ASN	ENGINEERED MUTATION	UNP P12821
C	25	GLN	ASN	ENGINEERED MUTATION	UNP P12821
C	82	GLN	ASN	ENGINEERED MUTATION	UNP P12821
C	117	GLN	ASN	ENGINEERED MUTATION	UNP P12821
C	289	GLN	ASN	ENGINEERED MUTATION	UNP P12821

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Chain	Residue	Modelled	Actual	Comment	Reference
C	545	ARG	GLN	ENGINEERED MUTATION	UNP P12821
C	576	LEU	PRO	ENGINEERED MUTATION	UNP P12821
C	629	LEU	ARG	ENGINEERED MUTATION	UNP P12821
D	9	GLN	ASN	ENGINEERED MUTATION	UNP P12821
D	25	GLN	ASN	ENGINEERED MUTATION	UNP P12821
D	82	GLN	ASN	ENGINEERED MUTATION	UNP P12821
D	117	GLN	ASN	ENGINEERED MUTATION	UNP P12821
D	289	GLN	ASN	ENGINEERED MUTATION	UNP P12821
D	545	ARG	GLN	ENGINEERED MUTATION	UNP P12821
D	576	LEU	PRO	ENGINEERED MUTATION	UNP P12821
D	629	LEU	ARG	ENGINEERED MUTATION	UNP P12821

- Molecule 2 is a protein called BETA-AMYLOID PROTEIN 42.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf	Trace
2	P	3	Total C N O 15 7 3 5	0	0	1
2	Q	3	Total C N O 15 7 3 5	0	0	1
2	R	3	Total C N O 15 7 3 5	0	0	1
2	S	3	Total C N O 15 7 3 5	0	0	1

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total Zn 1 1	0	0
3	A	1	Total Zn 1 1	0	0
3	D	1	Total Zn 1 1	0	0
3	C	1	Total Zn 1 1	0	0

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

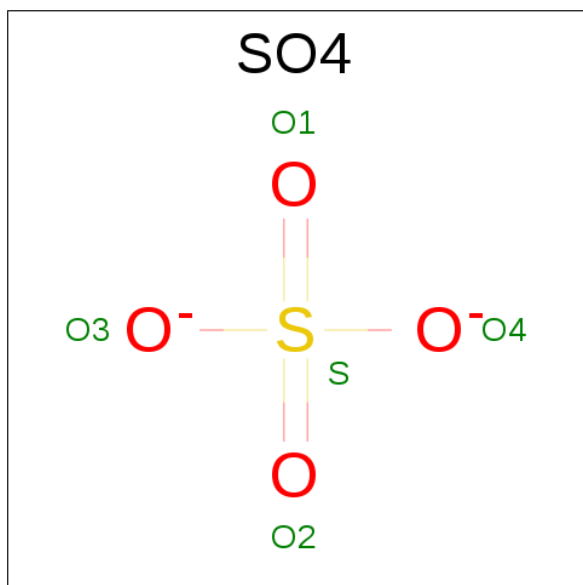
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total Cl 1 1	0	0

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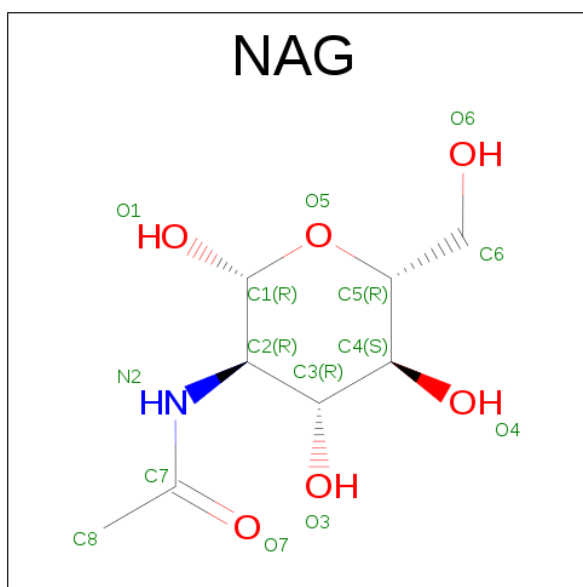
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Cl	0	0
			1	1		
4	D	1	Total	Cl	0	0
			1	1		
4	C	1	Total	Cl	0	0
			1	1		

- Molecule 5 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	O	S	0	0
			5	4	1		
5	B	1	Total	O	S	0	0
			5	4	1		
5	C	1	Total	O	S	0	0
			5	4	1		
5	D	1	Total	O	S	0	0
			5	4	1		

- Molecule 6 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: C<sub>8</sub>H<sub>15</sub>NO<sub>6</sub>).

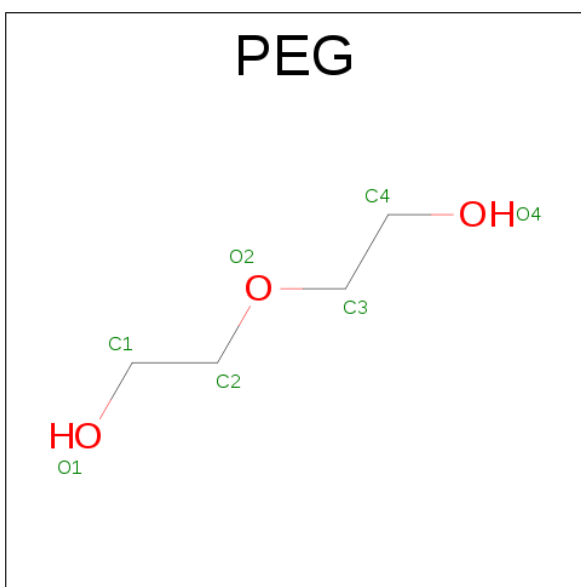


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total	C	N	O	0	0
			14	8	1	5		
6	D	1	Total	C	N	O	0	0
			14	8	1	5		
6	D	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 7 is a polymer of unknown type called SUGAR (2-MER).

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

- Molecule 8 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: C<sub>4</sub>H<sub>10</sub>O<sub>3</sub>).



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

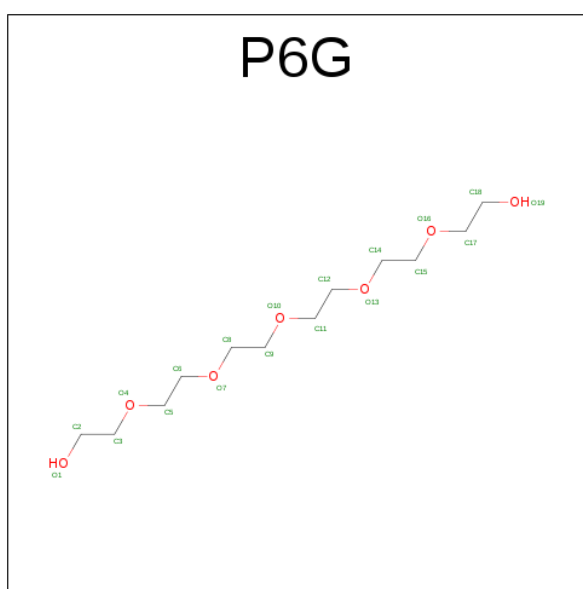
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	D	1	Total	C	O	0	0
			7	4	3		
8	D	1	Total	C	O	0	0
			7	4	3		
8	D	1	Total	C	O	0	0
			7	4	3		
8	D	1	Total	C	O	0	0
			7	4	3		

- Molecule 9 is HEXAETHYLENE GLYCOL (three-letter code: P6G) (formula:  $C_{12}H_{26}O_7$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	A	1	Total	C	O	0	0
			19	12	7		
9	B	1	Total	C	O	0	0
			19	12	7		
9	D	1	Total	C	O	0	0
			19	12	7		

- Molecule 10 is a polymer of unknown type called SUGAR (2-MER).

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

- Molecule 11 is a polymer of unknown type called SUGAR (3-MER).

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

- Molecule 12 is a polymer of unknown type called SUGAR (4-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
12	C	4	Total	C	N	O	0	0
			49	28	2	19		

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

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
13	D	4	Total	C	N	O	0	0
			49	28	2	19		

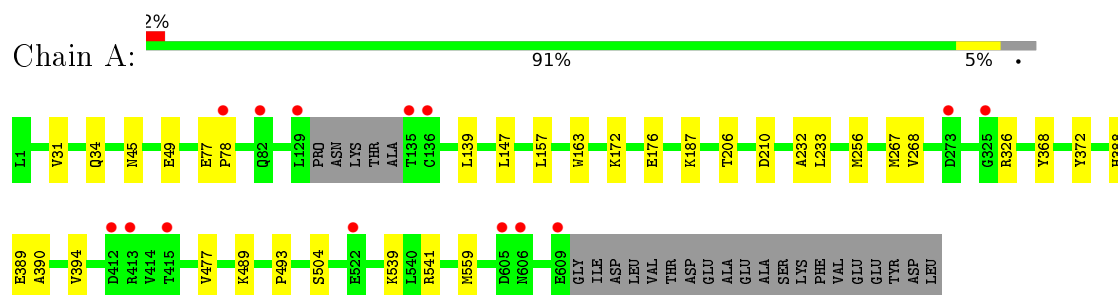
- Molecule 14 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
14	A	391	Total	O	0	0
			391	391		
14	B	458	Total	O	0	0
			458	458		
14	C	513	Total	O	0	0
			513	513		
14	D	464	Total	O	0	0
			464	464		
14	R	1	Total	O	0	0
			1	1		
14	S	2	Total	O	0	0
			2	2		

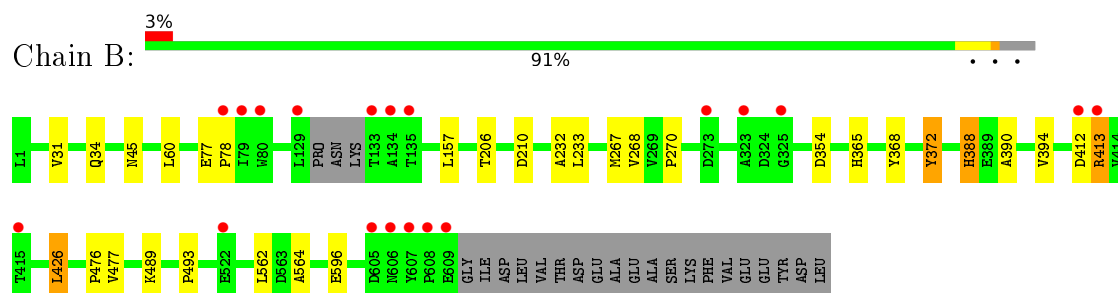
### 3 Residue-property plots [i](#)

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.

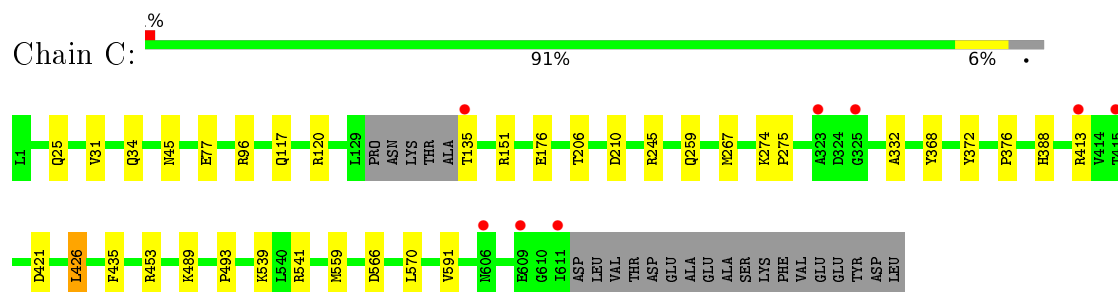
#### • Molecule 1: ANGIOTENSIN-CONVERTING ENZYME



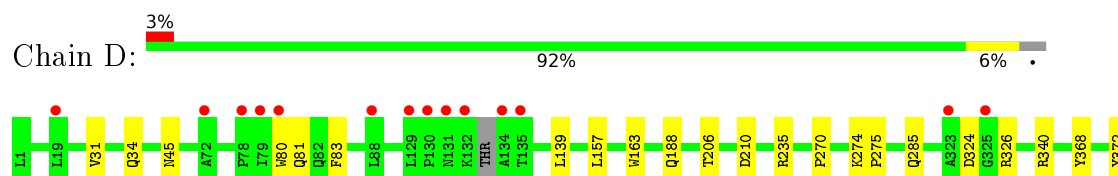
#### • Molecule 1: ANGIOTENSIN-CONVERTING ENZYME

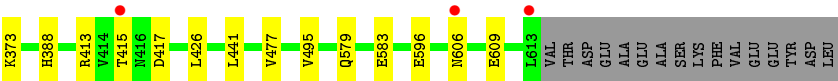


#### • Molecule 1: ANGIOTENSIN-CONVERTING ENZYME



#### • Molecule 1: ANGIOTENSIN-CONVERTING ENZYME

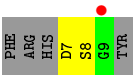




● Molecule 2: BETA-AMYLOID PROTEIN 42



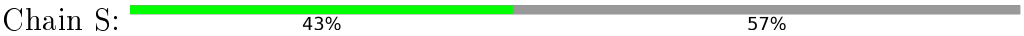
● Molecule 2: BETA-AMYLOID PROTEIN 42



● Molecule 2: BETA-AMYLOID PROTEIN 42



● Molecule 2: BETA-AMYLOID PROTEIN 42



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	73.45Å 101.76Å 114.36Å 85.23° 86.07° 81.45°	Depositor
Resolution (Å)	113.78 – 1.90 36.81 – 1.90	Depositor EDS
% Data completeness (in resolution range)	93.2 (113.78-1.90) 91.3 (36.81-1.90)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.63 (at 1.89Å)	Xtriage
Refinement program	REFMAC 5.8.0073	Depositor
R, $R_{free}$	0.185 , 0.224 0.194 , 0.229	Depositor DCC
$R_{free}$ test set	12020 reflections (5.29%)	DCC
Wilson B-factor (Å <sup>2</sup> )	20.0	Xtriage
Anisotropy	0.428	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 45.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 239414 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	22316	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 39.74 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 3.0333e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: ZN, BMA, NAG, CL, MAN, FUC, P6G, FUL, PEG, SO4

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	0.47	0/5098	0.64	1/6944 (0.0%)
1	B	0.50	0/5108	0.65	1/6957 (0.0%)
1	C	0.51	0/5117	0.67	2/6968 (0.0%)
1	D	0.51	0/5195	0.65	0/7075
2	P	1.41	0/14	1.12	0/18
2	Q	1.60	1/14 (7.1%)	0.97	0/18
2	R	1.33	0/14	1.25	0/18
2	S	1.43	0/14	0.92	0/18
All	All	0.50	1/20574 (0.0%)	0.65	4/28016 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	Q	8	SER	C-N	-5.48	1.23	1.33

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	541	ARG	NE-CZ-NH1	6.28	123.44	120.30
1	C	426	LEU	CA-CB-CG	-5.83	101.90	115.30
1	B	426	LEU	CA-CB-CG	-5.13	103.50	115.30
1	C	566	ASP	CB-CG-OD1	5.12	122.91	118.30

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	4939	0	4712	15	0
1	B	4950	0	4725	21	0
1	C	4959	0	4745	21	0
1	D	5029	0	4808	22	0
2	P	15	0	8	1	0
2	Q	15	0	8	1	0
2	R	15	0	8	1	0
2	S	15	0	8	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
5	A	5	0	0	0	0
5	B	5	0	0	0	0
5	C	5	0	0	0	0
5	D	5	0	0	0	0
6	A	14	0	13	0	0
6	D	28	0	26	1	0
7	A	56	0	50	1	0
7	B	28	0	25	0	0
7	C	28	0	25	0	0
8	A	28	0	40	3	0
8	B	35	0	50	0	0
8	C	28	0	40	0	0
8	D	35	0	50	0	0
9	A	19	0	26	1	0
9	B	19	0	26	0	0
9	D	19	0	26	3	0
10	B	24	0	22	1	0
10	C	24	0	22	1	0
11	B	39	0	34	0	0
12	C	49	0	43	0	0
13	D	49	0	43	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
14	A	391	0	0	4	0
14	B	458	0	0	3	0
14	C	513	0	0	5	0
14	D	464	0	0	2	0
14	R	1	0	0	0	0
14	S	2	0	0	0	0
All	All	22316	0	19583	85	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 (85) 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:B:270:PRO:HD3	1:B:426:LEU:HD22	1.67	0.75
1:A:147:LEU:HD22	1:A:256:MET:HA	1.68	0.74
1:C:206:THR:HG23	1:C:210:ASP:OD2	1.89	0.73
1:A:206:THR:HG23	1:A:210:ASP:OD2	1.91	0.71
1:B:365:HIS:HD1	1:B:388:HIS:CD2	2.09	0.70
1:D:206:THR:HG23	1:D:210:ASP:OD2	1.93	0.69
1:C:453:ARG:NH1	14:C:2414:HOH:O	2.23	0.69
1:C:413:ARG:NH2	14:C:2379:HOH:O	2.27	0.66
1:A:539:LYS:HE3	1:A:559:MET:O	1.96	0.65
1:C:176:GLU:OE1	14:C:2196:HOH:O	2.13	0.65
1:B:354:ASP:OD2	14:B:2315:HOH:O	2.15	0.63
1:B:206:THR:HG23	1:B:210:ASP:OD2	1.99	0.62
1:C:426:LEU:O	1:C:426:LEU:HG	1.99	0.61
8:A:1201:PEG:O4	9:A:1202:P6G:O19	2.19	0.60
1:B:365:HIS:HD1	1:B:388:HIS:HD2	1.49	0.60
1:C:245:ARG:HG2	1:C:591:VAL:HG11	1.83	0.59
1:B:596:GLU:OE2	6:D:1100:NAG:O7	2.19	0.59
10:C:1101:FUC:H4	14:C:2505:HOH:O	2.03	0.58
1:B:157:LEU:HD11	1:B:477:VAL:HG13	1.86	0.58
1:D:235:ARG:HH11	9:D:1201:P6G:H172	1.68	0.57
1:A:233:LEU:HD23	1:A:267:MET:HE1	1.88	0.56
1:D:157:LEU:HD11	1:D:477:VAL:HG13	1.88	0.55
1:C:274:LYS:HB3	1:C:275:PRO:CD	2.37	0.54
1:B:412:ASP:O	1:B:413:ARG:CB	2.56	0.53
1:D:188:GLN:HG2	14:D:2186:HOH:O	2.08	0.53
1:D:275:PRO:HG3	1:D:413:ARG:HG2	1.91	0.52
1:D:235:ARG:HH12	9:D:1201:P6G:H112	1.75	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:233:LEU:HD23	1:B:267:MET:HE1	1.93	0.51
1:A:172:LYS:O	1:A:176:GLU:HG3	2.11	0.51
10:B:1100:NAG:O7	1:D:596:GLU:OE2	2.29	0.50
1:A:187:LYS:HE3	14:A:2178:HOH:O	2.10	0.50
1:D:274:LYS:HB3	1:D:275:PRO:CD	2.42	0.50
1:B:232:ALA:CB	1:B:268:VAL:HG12	2.42	0.50
1:D:579:GLN:NE2	1:D:583:GLU:OE2	2.34	0.50
1:C:539:LYS:HE3	1:C:559:MET:O	2.12	0.49
1:A:389:GLU:HB2	1:A:504:SER:HB2	1.94	0.49
1:A:77:GLU:N	1:A:78:PRO:HD2	2.28	0.49
1:C:151:ARG:HD2	1:C:267:MET:SD	2.53	0.49
8:A:1201:PEG:H22	14:A:2389:HOH:O	2.11	0.49
1:B:412:ASP:O	1:B:413:ARG:HB3	2.13	0.49
1:B:489:LYS:O	1:B:493:PRO:HD2	2.14	0.48
1:D:340[B]:ARG:HG2	1:D:373:LYS:O	2.13	0.48
1:D:270:PRO:HD3	1:D:426:LEU:HD22	1.95	0.48
1:A:489:LYS:O	1:A:493:PRO:HD2	2.14	0.48
1:D:83:PHE:HA	14:D:2006:HOH:O	2.15	0.47
1:B:31:VAL:O	1:B:34:GLN:HG3	2.15	0.46
1:D:495:VAL:HG12	1:D:495:VAL:O	2.16	0.46
1:D:31:VAL:O	1:D:34:GLN:HG3	2.15	0.45
1:C:77:GLU:OE2	1:C:96:ARG:NH2	2.49	0.45
1:D:235:ARG:NH1	9:D:1201:P6G:H172	2.31	0.45
1:C:135:THR:O	1:C:135:THR:HG23	2.17	0.45
1:B:372:TYR:OH	1:B:388:HIS:HE1	1.99	0.44
1:D:80:TRP:CE2	1:D:81:GLN:HG3	2.53	0.44
1:A:232:ALA:CB	1:A:268:VAL:HG12	2.48	0.44
1:C:245:ARG:HG2	1:C:591:VAL:CG1	2.47	0.44
1:C:25:GLN:OE1	1:C:376:PRO:HA	2.17	0.44
1:C:570:LEU:C	1:C:570:LEU:HD23	2.37	0.44
1:A:31:VAL:O	1:A:34:GLN:HG3	2.18	0.44
1:B:426:LEU:HG	1:B:426:LEU:O	2.15	0.43
1:C:541[B]:ARG:NH1	14:C:2316:HOH:O	2.51	0.43
1:D:139:LEU:HD22	1:D:163:TRP:CZ2	2.53	0.43
1:A:49:GLU:HG2	7:A:1102:NAG:H82	2.00	0.43
1:C:259:GLN:O	1:C:435:PHE:HA	2.18	0.43
1:B:233:LEU:HD23	1:B:267:MET:CE	2.49	0.43
1:B:157:LEU:HD13	1:B:476:PRO:HB2	2.00	0.43
1:B:60:LEU:HD11	14:B:2001:HOH:O	2.18	0.42
1:D:606:ASN:O	1:D:609:GLU:N	2.52	0.42
1:C:332:ALA:HB3	2:R:7:ASP:HB2	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:117:GLN:HE22	1:C:120:ARG:HE	1.66	0.42
1:D:270:PRO:HD3	1:D:426:LEU:CD2	2.49	0.42
14:B:2323:HOH:O	2:Q:7:ASP:N	2.52	0.42
14:A:2288:HOH:O	2:P:7:ASP:N	2.53	0.42
1:A:157:LEU:HD11	1:A:477:VAL:HG13	2.02	0.41
1:B:77:GLU:HB3	1:B:78:PRO:HD3	2.02	0.41
1:D:324:ASP:OD1	1:D:326:ARG:HB2	2.21	0.41
1:C:489:LYS:O	1:C:493:PRO:HD2	2.20	0.41
1:D:415:THR:OG1	1:D:417:ASP:OD2	2.38	0.41
1:C:31:VAL:O	1:C:34:GLN:HG3	2.20	0.41
1:A:390:ALA:O	1:A:394:VAL:HG23	2.21	0.41
1:D:441:LEU:C	1:D:441:LEU:HD12	2.40	0.41
1:A:139:LEU:HD22	1:A:163:TRP:CZ2	2.56	0.41
1:C:77:GLU:OE1	1:C:77:GLU:HA	2.20	0.41
1:B:562:LEU:HD23	1:B:564:ALA:O	2.21	0.40
8:A:1203:PEG:H41	14:A:2127:HOH:O	2.20	0.40
1:B:390:ALA:O	1:B:394:VAL:HG23	2.22	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	601/629 (96%)	590 (98%)	9 (2%)	2 (0%)	46	35
1	B	603/629 (96%)	593 (98%)	8 (1%)	2 (0%)	46	35
1	C	603/629 (96%)	593 (98%)	9 (2%)	1 (0%)	52	42
1	D	613/629 (98%)	601 (98%)	11 (2%)	1 (0%)	52	42
2	P	1/7 (14%)	1 (100%)	0	0	100	100
2	Q	1/7 (14%)	1 (100%)	0	0	100	100
2	R	1/7 (14%)	1 (100%)	0	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	S	1/7 (14%)	1 (100%)	0	0	100	100
All	All	2424/2544 (95%)	2381 (98%)	37 (2%)	6 (0%)	52	42

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	413	ARG
1	A	326	ARG
1	D	45	ASN
1	B	45	ASN
1	A	45	ASN
1	C	45	ASN

### 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	519/541 (96%)	516 (99%)	3 (1%)	90	90
1	B	519/541 (96%)	516 (99%)	3 (1%)	90	90
1	C	522/541 (96%)	518 (99%)	4 (1%)	86	86
1	D	529/541 (98%)	525 (99%)	4 (1%)	86	86
2	P	2/6 (33%)	2 (100%)	0	100	100
2	Q	2/6 (33%)	2 (100%)	0	100	100
2	R	2/6 (33%)	2 (100%)	0	100	100
2	S	2/6 (33%)	2 (100%)	0	100	100
All	All	2097/2188 (96%)	2083 (99%)	14 (1%)	88	88

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	368	TYR
1	A	372	TYR

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Mol	Chain	Res	Type
1	A	388	HIS
1	B	368	TYR
1	B	372	TYR
1	B	388	HIS
1	C	368	TYR
1	C	372	TYR
1	C	388	HIS
1	C	421	ASP
1	D	285	GLN
1	D	368	TYR
1	D	372	TYR
1	D	388	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	203	ASN
1	A	289	GLN
1	B	388	HIS
1	C	117	GLN
1	C	203	ASN
1	C	598	GLN
1	D	18	GLN
1	D	188	GLN

### 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 ⓘ

23 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
7	NAG	A	1101	1,7	14,14,15	0.54	0	15,19,21	1.19	1 (6%)
7	NAG	A	1102	7	14,14,15	0.56	0	15,19,21	0.77	0
7	NAG	A	1103	1,7	14,14,15	0.69	0	15,19,21	0.82	0
7	NAG	A	1104	7	14,14,15	0.53	0	15,19,21	1.09	1 (6%)
10	NAG	B	1100	1,10	14,14,15	0.75	0	15,19,21	2.33	3 (20%)
10	FUC	B	1101	10	10,10,11	0.73	0	13,14,16	1.59	2 (15%)
7	NAG	B	1102	1,7	14,14,15	0.69	0	15,19,21	1.10	1 (6%)
7	NAG	B	1103	7	14,14,15	0.49	0	15,19,21	0.79	0
11	NAG	B	1104	1,11	14,14,15	0.59	0	15,19,21	0.83	0
11	NAG	B	1105	11	14,14,15	0.65	0	15,19,21	0.76	0
11	BMA	B	1106	11	11,11,12	0.36	0	15,15,17	0.86	0
10	NAG	C	1100	1,10	14,14,15	0.66	0	15,19,21	1.56	4 (26%)
10	FUC	C	1101	10	10,10,11	0.66	0	13,14,16	2.15	5 (38%)
7	NAG	C	1102	1,7	14,14,15	0.58	0	15,19,21	1.20	1 (6%)
7	NAG	C	1103	7	14,14,15	0.55	0	15,19,21	1.25	3 (20%)
12	NAG	C	1104	1,12	14,14,15	0.87	1 (7%)	15,19,21	0.87	0
12	NAG	C	1105	12	14,14,15	0.69	0	15,19,21	0.90	0
12	BMA	C	1106	12	11,11,12	0.65	0	15,15,17	2.39	3 (20%)
12	FUL	C	1107	12	10,10,11	0.86	0	13,14,16	1.66	4 (30%)
13	NAG	D	1102	1,13	14,14,15	0.85	1 (7%)	15,19,21	0.74	0
13	NAG	D	1103	13	14,14,15	0.63	0	15,19,21	0.72	0
13	MAN	D	1104	13	11,11,12	0.63	0	15,15,17	2.09	4 (26%)
13	FUC	D	1105	13	10,10,11	0.57	0	13,14,16	1.31	1 (7%)

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
7	NAG	A	1101	1,7	-	0/6/23/26	0/1/1/1
7	NAG	A	1102	7	-	0/6/23/26	0/1/1/1
7	NAG	A	1103	1,7	-	0/6/23/26	0/1/1/1
7	NAG	A	1104	7	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
10	NAG	B	1100	1,10	-	0/6/23/26	0/1/1/1
10	FUC	B	1101	10	-	0/0/17/20	0/1/1/1
7	NAG	B	1102	1,7	-	0/6/23/26	0/1/1/1
7	NAG	B	1103	7	-	0/6/23/26	0/1/1/1
11	NAG	B	1104	1,11	-	0/6/23/26	0/1/1/1
11	NAG	B	1105	11	-	0/6/23/26	0/1/1/1
11	BMA	B	1106	11	-	0/2/19/22	0/1/1/1
10	NAG	C	1100	1,10	-	0/6/23/26	0/1/1/1
10	FUC	C	1101	10	-	0/0/17/20	0/1/1/1
7	NAG	C	1102	1,7	-	0/6/23/26	0/1/1/1
7	NAG	C	1103	7	-	0/6/23/26	0/1/1/1
12	NAG	C	1104	1,12	-	0/6/23/26	0/1/1/1
12	NAG	C	1105	12	-	0/6/23/26	0/1/1/1
12	BMA	C	1106	12	-	0/2/19/22	0/1/1/1
12	FUL	C	1107	12	-	0/0/17/20	0/1/1/1
13	NAG	D	1102	1,13	-	0/6/23/26	0/1/1/1
13	NAG	D	1103	13	-	0/6/23/26	0/1/1/1
13	MAN	D	1104	13	-	0/2/19/22	0/1/1/1
13	FUC	D	1105	13	-	0/0/17/20	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
12	C	1104	NAG	O5-C1	-2.18	1.40	1.43
13	D	1102	NAG	O5-C1	-2.08	1.40	1.43

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
10	B	1100	NAG	C2-N2-C7	-6.43	114.74	123.11
10	B	1101	FUC	O5-C1-C2	-3.67	105.03	110.89
10	B	1100	NAG	O7-C7-N2	-3.56	114.58	121.84
13	D	1105	FUC	C1-C2-C3	-2.73	106.24	109.55
7	A	1104	NAG	O7-C7-C8	-2.42	117.62	122.07
10	C	1101	FUC	O5-C5-C4	-2.29	105.61	109.58
7	C	1103	NAG	C3-C4-C5	-2.28	106.17	110.23
12	C	1107	FUL	O3-C3-C2	-2.22	105.94	110.01
10	C	1100	NAG	O3-C3-C4	-2.13	105.57	110.36
7	C	1103	NAG	O5-C5-C4	-2.09	106.67	110.13
12	C	1107	FUL	C3-C4-C5	2.11	112.80	109.66
10	C	1100	NAG	O5-C5-C6	2.12	111.88	107.34
10	C	1100	NAG	C2-N2-C7	2.21	125.98	123.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	C	1103	NAG	O5-C5-C6	2.24	112.14	107.34
7	B	1102	NAG	C1-O5-C5	2.31	115.53	112.14
10	C	1101	FUC	C2-C3-C4	2.39	115.22	111.05
12	C	1107	FUL	C1-C2-C3	2.50	112.58	109.55
10	C	1101	FUC	O5-C1-C2	2.58	115.02	110.89
7	C	1102	NAG	C1-O5-C5	2.86	116.34	112.14
13	D	1104	MAN	C2-C3-C4	2.95	116.20	111.05
10	B	1101	FUC	O5-C5-C6	3.03	111.71	106.28
10	C	1100	NAG	C4-C3-C2	3.45	116.70	111.34
12	C	1107	FUL	O5-C5-C6	3.54	112.63	106.28
13	D	1104	MAN	O5-C5-C6	3.67	115.20	107.34
7	A	1101	NAG	C1-O5-C5	3.76	117.67	112.14
13	D	1104	MAN	C1-C2-C3	3.77	114.12	109.55
10	B	1100	NAG	C8-C7-N2	3.80	123.38	116.10
12	C	1106	BMA	O5-C5-C4	3.91	116.61	110.13
10	C	1101	FUC	C3-C4-C5	4.02	115.66	109.66
12	C	1106	BMA	C3-C4-C5	4.22	117.76	110.23
13	D	1104	MAN	C3-C4-C5	4.29	117.87	110.23
10	C	1101	FUC	C1-C2-C3	4.67	115.21	109.55
12	C	1106	BMA	C1-O5-C5	6.67	121.95	112.14

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	A	1102	NAG	1	0
10	B	1100	NAG	1	0
10	C	1101	FUC	1	0

## 5.6 Ligand geometry ⓘ

Of 36 ligands modelled in this entry, 8 are monoatomic - leaving 28 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
5	SO4	A	1003	-	4,4,4	0.49	0	6,6,6	0.16	0
6	NAG	A	1100	1	14,14,15	0.44	0	15,19,21	1.21	1 (6%)
8	PEG	A	1200	-	6,6,6	0.44	0	5,5,5	0.40	0
8	PEG	A	1201	-	6,6,6	0.40	0	5,5,5	0.43	0
9	P6G	A	1202	-	18,18,18	0.49	0	17,17,17	0.59	0
8	PEG	A	1203	-	6,6,6	0.51	0	5,5,5	0.31	0
8	PEG	A	1204	-	6,6,6	0.36	0	5,5,5	0.44	0
5	SO4	B	1003	-	4,4,4	0.46	0	6,6,6	0.34	0
8	PEG	B	1200	-	6,6,6	0.52	0	5,5,5	0.32	0
8	PEG	B	1201	-	6,6,6	0.61	0	5,5,5	0.51	0
8	PEG	B	1202	-	6,6,6	0.49	0	5,5,5	0.41	0
9	P6G	B	1203	-	18,18,18	0.64	0	17,17,17	0.79	0
8	PEG	B	1204	-	6,6,6	0.63	0	5,5,5	0.23	0
8	PEG	B	1205	-	6,6,6	0.54	0	5,5,5	0.28	0
5	SO4	C	1003	-	4,4,4	0.74	0	6,6,6	0.34	0
8	PEG	C	1200	-	6,6,6	0.44	0	5,5,5	0.35	0
8	PEG	C	1201	-	6,6,6	0.45	0	5,5,5	0.34	0
8	PEG	C	1202	-	6,6,6	0.36	0	5,5,5	0.61	0
8	PEG	C	1203	-	6,6,6	0.38	0	5,5,5	0.48	0
5	SO4	D	1003	-	4,4,4	0.61	0	6,6,6	0.19	0
6	NAG	D	1100	1	14,14,15	0.29	0	15,19,21	0.53	0
6	NAG	D	1101	1	14,14,15	0.49	0	15,19,21	1.03	1 (6%)
8	PEG	D	1200	-	6,6,6	0.39	0	5,5,5	0.43	0
9	P6G	D	1201	-	18,18,18	0.59	0	17,17,17	0.66	0
8	PEG	D	1202	-	6,6,6	0.52	0	5,5,5	0.40	0
8	PEG	D	1203	-	6,6,6	0.52	0	5,5,5	0.19	0
8	PEG	D	1204	-	6,6,6	0.45	0	5,5,5	0.38	0
8	PEG	D	1205	-	6,6,6	0.63	0	5,5,5	0.50	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
5	SO4	A	1003	-	-	0/0/0/0	0/0/0/0
6	NAG	A	1100	1	-	0/6/23/26	0/1/1/1
8	PEG	A	1200	-	-	0/4/4/4	0/0/0/0
8	PEG	A	1201	-	-	0/4/4/4	0/0/0/0
9	P6G	A	1202	-	-	0/16/16/16	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
8	PEG	A	1203	-	-	0/4/4/4	0/0/0/0
8	PEG	A	1204	-	-	0/4/4/4	0/0/0/0
5	SO4	B	1003	-	-	0/0/0/0	0/0/0/0
8	PEG	B	1200	-	-	0/4/4/4	0/0/0/0
8	PEG	B	1201	-	-	0/4/4/4	0/0/0/0
8	PEG	B	1202	-	-	0/4/4/4	0/0/0/0
9	P6G	B	1203	-	-	0/16/16/16	0/0/0/0
8	PEG	B	1204	-	-	0/4/4/4	0/0/0/0
8	PEG	B	1205	-	-	0/4/4/4	0/0/0/0
5	SO4	C	1003	-	-	0/0/0/0	0/0/0/0
8	PEG	C	1200	-	-	0/4/4/4	0/0/0/0
8	PEG	C	1201	-	-	0/4/4/4	0/0/0/0
8	PEG	C	1202	-	-	0/4/4/4	0/0/0/0
8	PEG	C	1203	-	-	0/4/4/4	0/0/0/0
5	SO4	D	1003	-	-	0/0/0/0	0/0/0/0
6	NAG	D	1100	1	-	0/6/23/26	0/1/1/1
6	NAG	D	1101	1	-	0/6/23/26	0/1/1/1
8	PEG	D	1200	-	-	0/4/4/4	0/0/0/0
9	P6G	D	1201	-	-	0/16/16/16	0/0/0/0
8	PEG	D	1202	-	-	0/4/4/4	0/0/0/0
8	PEG	D	1203	-	-	0/4/4/4	0/0/0/0
8	PEG	D	1204	-	-	0/4/4/4	0/0/0/0
8	PEG	D	1205	-	-	0/4/4/4	0/0/0/0

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	D	1101	NAG	C1-O5-C5	2.45	115.75	112.14
6	A	1100	NAG	C1-O5-C5	3.09	116.69	112.14

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	A	1201	PEG	2	0
9	A	1202	P6G	1	0
8	A	1203	PEG	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	D	1100	NAG	1	0
9	D	1201	P6G	3	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	604/629 (96%)	-0.13	14 (2%) 64 67	15, 25, 46, 73	0
1	B	606/629 (96%)	-0.14	19 (3%) 52 56	13, 23, 43, 68	0
1	C	606/629 (96%)	-0.29	8 (1%) 79 82	12, 20, 36, 69	0
1	D	612/629 (97%)	-0.16	17 (2%) 56 60	11, 20, 42, 81	0
2	P	3/7 (42%)	1.12	0 100 100	23, 23, 23, 29	0
2	Q	3/7 (42%)	1.52	1 (33%) 0 0	19, 19, 19, 28	0
2	R	3/7 (42%)	0.35	0 100 100	20, 20, 22, 29	0
2	S	3/7 (42%)	0.40	0 100 100	18, 18, 19, 27	0
All	All	2440/2544 (95%)	-0.18	59 (2%) 62 66	11, 22, 42, 81	0

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	325	GLY	8.7
1	B	325	GLY	7.2
1	B	134	ALA	6.5
1	D	130	PRO	5.8
1	B	135	THR	5.2
1	A	415	THR	5.0
1	D	79	ILE	4.9
1	D	613	LEU	4.8
1	D	134	ALA	4.5
1	A	606	ASN	4.4
1	C	135	THR	4.4
1	D	131	ASN	4.3
1	A	413	ARG	4.2
1	A	135	THR	4.1
1	C	611	ILE	4.0
1	A	325	GLY	3.7

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Mol	Chain	Res	Type	RSRZ
1	B	80	TRP	3.6
1	A	78	PRO	3.6
1	C	413	ARG	3.5
1	B	129	LEU	3.5
1	C	606	ASN	3.4
1	B	606	ASN	3.3
1	D	132	LYS	3.3
1	B	609	GLU	3.3
1	C	325	GLY	3.2
1	A	412	ASP	3.2
1	B	415	THR	3.0
1	B	605	ASP	3.0
1	D	135	THR	3.0
1	B	607	TYR	2.9
1	B	79	ILE	2.8
1	C	609	GLU	2.7
1	D	19	LEU	2.7
2	Q	9	GLY	2.6
1	A	605	ASP	2.5
1	D	323	ALA	2.5
1	B	412	ASP	2.5
1	D	80	TRP	2.5
1	A	82	GLN	2.4
1	B	78	PRO	2.4
1	B	323	ALA	2.3
1	D	129	LEU	2.3
1	A	273	ASP	2.3
1	D	78	PRO	2.3
1	C	415	THR	2.3
1	A	609	GLU	2.3
1	B	273	ASP	2.3
1	B	522	GLU	2.2
1	D	72	ALA	2.2
1	A	522	GLU	2.2
1	A	136	CYS	2.1
1	B	133	THR	2.1
1	B	413	ARG	2.1
1	D	606	ASN	2.1
1	C	323	ALA	2.1
1	D	415	THR	2.0
1	B	608	PRO	2.0
1	A	129	LEU	2.0

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Mol	Chain	Res	Type	RSRZ
1	D	88	LEU	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](#)

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
7	NAG	A	1103	14/15	0.88	0.24	3.87	62,64,68,70	0
7	NAG	A	1101	14/15	0.91	0.26	2.62	36,38,44,51	0
7	NAG	B	1102	14/15	0.83	0.22	2.56	38,42,49,58	0
10	NAG	C	1100	14/15	0.85	0.15	1.89	34,40,48,50	0
7	NAG	C	1102	14/15	0.93	0.15	1.74	27,31,35,38	0
10	NAG	B	1100	14/15	0.80	0.13	1.49	33,36,43,44	0
11	NAG	B	1104	14/15	0.93	0.14	0.88	46,48,51,52	0
13	NAG	D	1102	14/15	0.95	0.11	0.17	30,33,35,36	0
12	NAG	C	1104	14/15	0.96	0.10	0.08	29,30,34,38	0
11	BMA	B	1106	11/12	0.72	0.25	-	64,67,69,70	0
10	FUC	B	1101	10/11	0.78	0.20	-	46,50,52,53	0
7	NAG	B	1103	14/15	0.67	0.39	-	62,69,75,79	0
13	FUC	D	1105	10/11	0.91	0.15	-	37,40,44,44	0
13	MAN	D	1104	11/12	0.69	0.23	-	55,59,60,61	0
12	NAG	C	1105	14/15	0.92	0.16	-	36,40,46,46	0
13	NAG	D	1103	14/15	0.90	0.13	-	39,45,49,54	0
11	NAG	B	1105	14/15	0.91	0.23	-	51,54,57,62	0
7	NAG	C	1103	14/15	0.86	0.30	-	43,47,54,57	0
12	FUL	C	1107	10/11	0.82	0.19	-	43,47,49,51	0
12	BMA	C	1106	11/12	0.77	0.21	-	45,48,53,54	0
7	NAG	A	1102	14/15	0.79	0.38	-	57,62,64,65	0
10	FUC	C	1101	10/11	0.62	0.32	-	54,61,65,69	0
7	NAG	A	1104	14/15	0.87	0.31	-	66,67,70,75	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
8	PEG	C	1202	7/7	0.79	0.26	24.55	46,47,50,52	0
8	PEG	D	1203	7/7	0.83	0.20	17.64	46,49,53,53	0
5	SO4	D	1003	5/5	0.63	0.29	8.41	69,69,79,84	0
8	PEG	D	1204	7/7	0.89	0.24	7.74	43,45,53,55	0
8	PEG	C	1203	7/7	0.93	0.20	7.70	41,43,44,46	0
8	PEG	B	1205	7/7	0.79	0.32	5.72	50,51,54,58	0
8	PEG	B	1201	7/7	0.72	0.25	5.57	41,43,46,47	0
8	PEG	B	1204	7/7	0.79	0.24	5.41	49,51,53,53	0
9	P6G	B	1203	19/19	0.76	0.16	4.52	37,39,51,51	0
8	PEG	C	1201	7/7	0.93	0.22	4.50	36,38,43,47	0
5	SO4	C	1003	5/5	0.88	0.18	3.59	60,62,65,74	0
8	PEG	A	1203	7/7	0.89	0.27	3.55	49,50,52,52	0
6	NAG	A	1100	14/15	0.71	0.19	3.48	42,47,52,60	0
5	SO4	A	1003	5/5	0.80	0.25	3.42	74,80,84,88	0
8	PEG	A	1201	7/7	0.85	0.13	2.67	40,43,45,45	0
8	PEG	B	1202	7/7	0.88	0.12	2.46	33,33,34,34	0
8	PEG	A	1200	7/7	0.84	0.22	2.31	47,48,52,54	0
6	NAG	D	1101	14/15	0.82	0.23	2.27	37,41,46,46	0
9	P6G	D	1201	19/19	0.87	0.14	2.24	31,36,50,51	0
6	NAG	D	1100	14/15	0.78	0.14	1.91	30,36,40,42	0
9	P6G	A	1202	19/19	0.90	0.11	0.69	36,39,40,41	0
5	SO4	B	1003	5/5	0.88	0.18	0.54	71,73,75,77	0
8	PEG	C	1200	7/7	0.86	0.13	0.12	43,44,45,47	0
4	CL	A	1002	1/1	1.00	0.09	-1.02	18,18,18,18	0
4	CL	B	1002	1/1	0.99	0.09	-1.43	16,16,16,16	0
4	CL	C	1002	1/1	1.00	0.08	-1.94	15,15,15,15	0
3	ZN	B	1001	1/1	1.00	0.08	-2.41	16,16,16,16	0
4	CL	D	1002	1/1	1.00	0.07	-2.45	14,14,14,14	0
3	ZN	C	1001	1/1	1.00	0.06	-2.59	15,15,15,15	0
3	ZN	D	1001	1/1	1.00	0.07	-3.77	15,15,15,15	0
3	ZN	A	1001	1/1	1.00	0.06	-3.99	19,19,19,19	0
8	PEG	D	1200	7/7	0.81	0.19	-	52,53,55,60	0
8	PEG	D	1205	7/7	0.83	0.18	-	37,40,42,43	0
8	PEG	A	1204	7/7	0.83	0.24	-	50,52,54,55	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
8	PEG	D	1202	7/7	0.82	0.16	-	52,55,58,58	0
8	PEG	B	1200	7/7	0.75	0.26	-	59,61,62,63	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.