



# wwPDB X-ray Structure Validation Summary Report ⓘ

Aug 8, 2020 – 01:12 AM BST

PDB ID : 5LG6  
Title : Structure of the deglycosylated porcine aminopeptidase N ectodomain  
Authors : Santiago, C.; Reguera, J.; Mudgal, G.; Casasnovas, J.M.  
Deposited on : 2016-07-06  
Resolution : 2.50 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

<https://www.wwpdb.org/validation/2017/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 : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.13.1  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.13.1

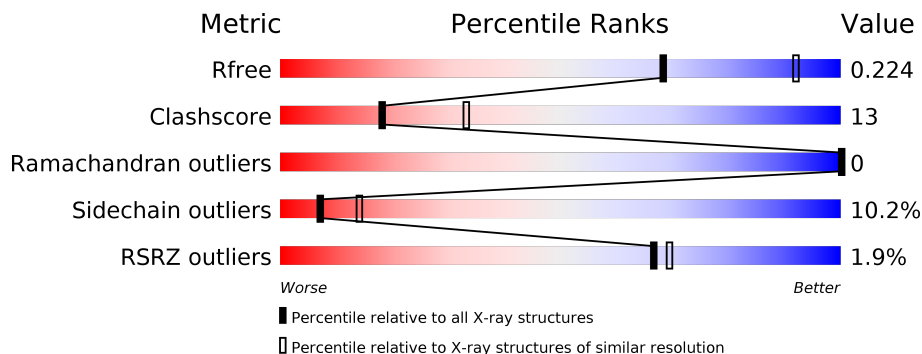
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.50 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))
$R_{free}$	130704	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

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 respectively. 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	968	 2% 67% 20% 8%
1	B	968	 2% 67% 20% 8%

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 14927 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 Aminopeptidase N.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	A	889	7154	4578	1190	1356	6	24	0	0	0
1	B	887	7138	4564	1188	1356	6	24	0	0	0

There are 84 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	19	TYR	-	expression tag	UNP P15145
A	20	PRO	-	expression tag	UNP P15145
A	21	TYR	-	expression tag	UNP P15145
A	22	ASP	-	expression tag	UNP P15145
A	23	VAL	-	expression tag	UNP P15145
A	24	PRO	-	expression tag	UNP P15145
A	25	ASP	-	expression tag	UNP P15145
A	26	TYR	-	expression tag	UNP P15145
A	27	ALA	-	expression tag	UNP P15145
A	28	GLY	-	expression tag	UNP P15145
A	29	ALA	-	expression tag	UNP P15145
A	30	GLN	-	expression tag	UNP P15145
A	31	PRO	-	expression tag	UNP P15145
A	32	ALA	-	expression tag	UNP P15145
A	33	ARG	-	expression tag	UNP P15145
A	34	SER	-	expression tag	UNP P15145
A	35	PRO	-	expression tag	UNP P15145
A	107	PHE	LEU	conflict	UNP P15145
A	108	ILE	LEU	conflict	UNP P15145
A	964	VAL	-	expression tag	UNP P15145
A	965	LEU	-	expression tag	UNP P15145
A	966	ASN	-	expression tag	UNP P15145
A	967	TRP	-	expression tag	UNP P15145
A	968	PHE	-	expression tag	UNP P15145
A	969	ILE	-	expression tag	UNP P15145

*Continued on next page...*

*Continued from previous page...*

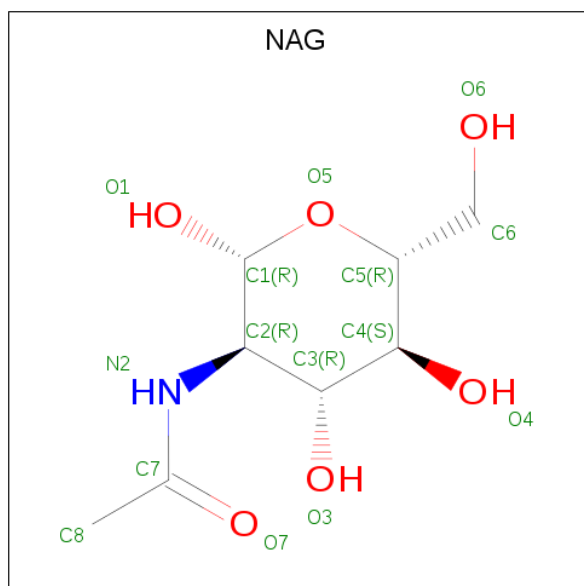
Chain	Residue	Modelled	Actual	Comment	Reference
A	970	GLU	-	expression tag	UNP P15145
A	971	HIS	-	expression tag	UNP P15145
A	972	SER	-	expression tag	UNP P15145
A	973	LEU	-	expression tag	UNP P15145
A	974	VAL	-	expression tag	UNP P15145
A	975	PRO	-	expression tag	UNP P15145
A	976	ARG	-	expression tag	UNP P15145
A	977	GLY	-	expression tag	UNP P15145
A	978	SER	-	expression tag	UNP P15145
A	979	ASP	-	expression tag	UNP P15145
A	980	TYR	-	expression tag	UNP P15145
A	981	LYS	-	expression tag	UNP P15145
A	982	ASP	-	expression tag	UNP P15145
A	983	ASP	-	expression tag	UNP P15145
A	984	ASP	-	expression tag	UNP P15145
A	985	ASP	-	expression tag	UNP P15145
A	986	LYS	-	expression tag	UNP P15145
B	19	TYR	-	expression tag	UNP P15145
B	20	PRO	-	expression tag	UNP P15145
B	21	TYR	-	expression tag	UNP P15145
B	22	ASP	-	expression tag	UNP P15145
B	23	VAL	-	expression tag	UNP P15145
B	24	PRO	-	expression tag	UNP P15145
B	25	ASP	-	expression tag	UNP P15145
B	26	TYR	-	expression tag	UNP P15145
B	27	ALA	-	expression tag	UNP P15145
B	28	GLY	-	expression tag	UNP P15145
B	29	ALA	-	expression tag	UNP P15145
B	30	GLN	-	expression tag	UNP P15145
B	31	PRO	-	expression tag	UNP P15145
B	32	ALA	-	expression tag	UNP P15145
B	33	ARG	-	expression tag	UNP P15145
B	34	SER	-	expression tag	UNP P15145
B	35	PRO	-	expression tag	UNP P15145
B	107	PHE	LEU	conflict	UNP P15145
B	108	ILE	LEU	conflict	UNP P15145
B	964	VAL	-	expression tag	UNP P15145
B	965	LEU	-	expression tag	UNP P15145
B	966	ASN	-	expression tag	UNP P15145
B	967	TRP	-	expression tag	UNP P15145
B	968	PHE	-	expression tag	UNP P15145
B	969	ILE	-	expression tag	UNP P15145

*Continued on next page...*

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
B	970	GLU	-	expression tag	UNP P15145
B	971	HIS	-	expression tag	UNP P15145
B	972	SER	-	expression tag	UNP P15145
B	973	LEU	-	expression tag	UNP P15145
B	974	VAL	-	expression tag	UNP P15145
B	975	PRO	-	expression tag	UNP P15145
B	976	ARG	-	expression tag	UNP P15145
B	977	GLY	-	expression tag	UNP P15145
B	978	SER	-	expression tag	UNP P15145
B	979	ASP	-	expression tag	UNP P15145
B	980	TYR	-	expression tag	UNP P15145
B	981	LYS	-	expression tag	UNP P15145
B	982	ASP	-	expression tag	UNP P15145
B	983	ASP	-	expression tag	UNP P15145
B	984	ASP	-	expression tag	UNP P15145
B	985	ASP	-	expression tag	UNP P15145
B	986	LYS	-	expression tag	UNP P15145

- Molecule 2 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula:  $C_8H_{15}NO_6$ ).



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

Continued on next page...

*Continued from previous page...*

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

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

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

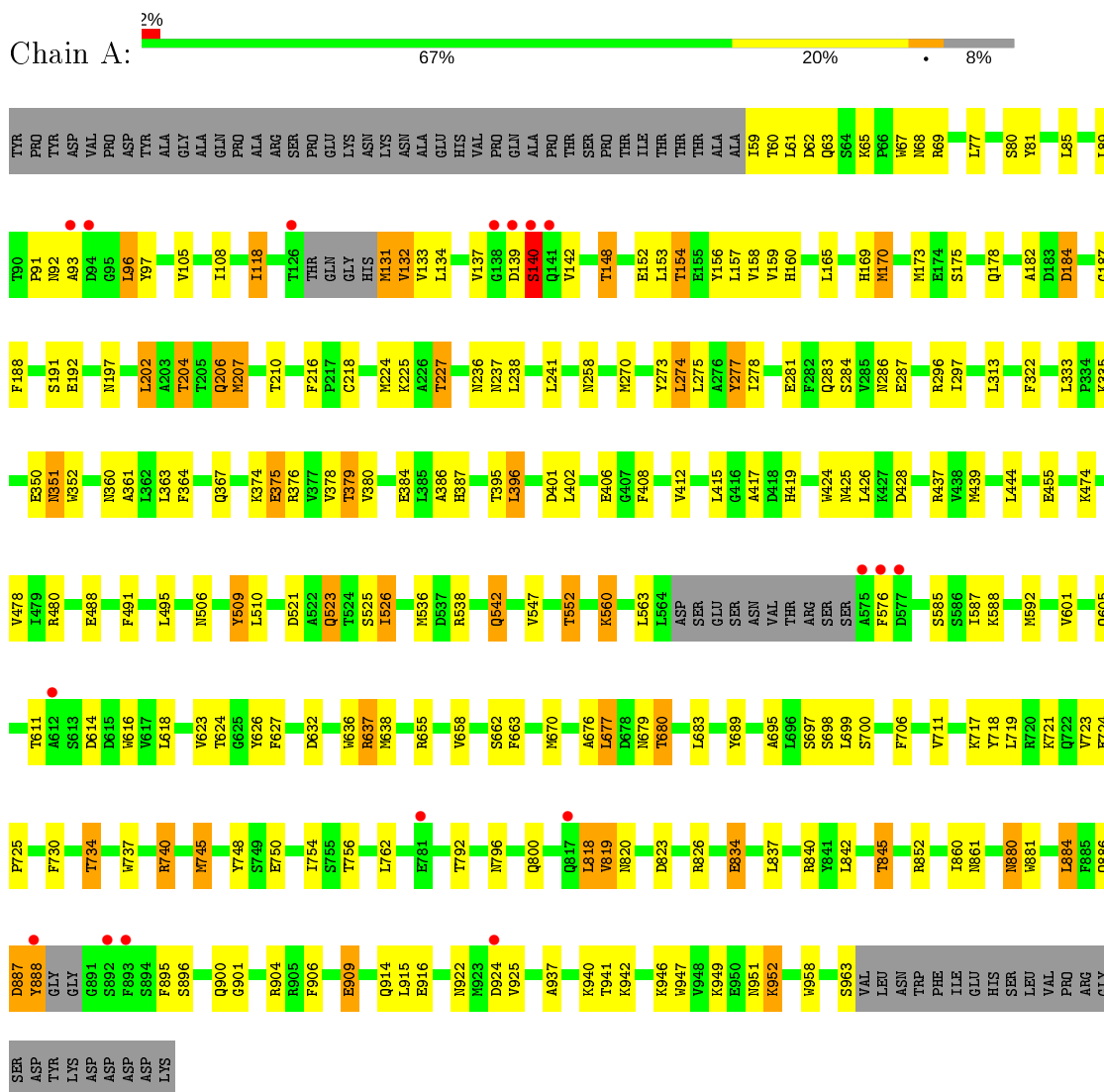
- Molecule 4 is water.

<b>Mol</b>	<b>Chain</b>	<b>Residues</b>	<b>Atoms</b>		<b>ZeroOcc</b>	<b>AltConf</b>
4	A	216	Total 216	O 216	0	0
4	B	165	Total 165	O 165	0	0

### 3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes 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.

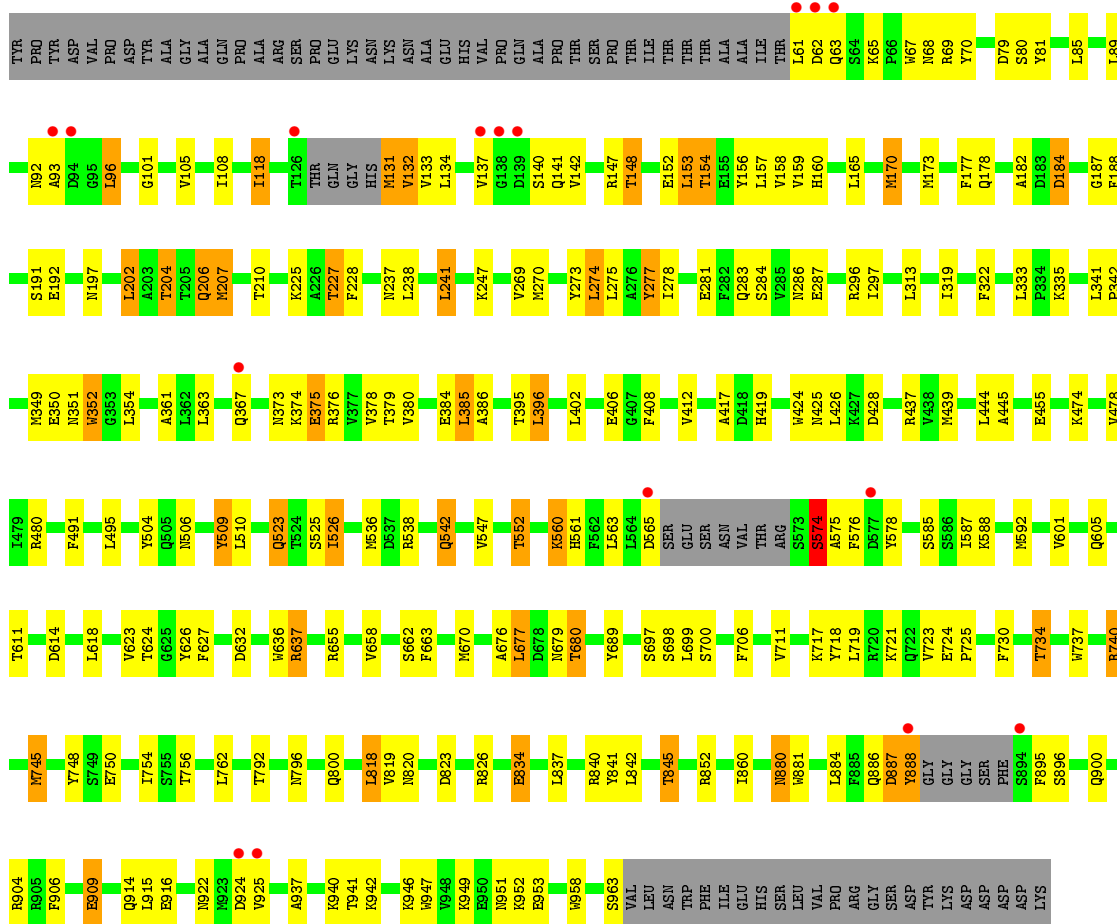
- Molecule 1: Aminopeptidase N



- Molecule 1: Aminopeptidase N







## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	66.47Å 215.69Å 78.63Å 90.00° 91.90° 90.00°	Depositor
Resolution (Å)	24.98 – 2.50 24.98 – 2.50	Depositor EDS
% Data completeness (in resolution range)	97.0 (24.98-2.50) 93.8 (24.98-2.50)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.10 (at 2.50Å)	Xtrriage
Refinement program	PHENIX 1.7.1_743	Depositor
R, $R_{free}$	0.191 , 0.229 0.185 , 0.224	Depositor DCC
$R_{free}$ test set	1989 reflections (2.69%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	15.0	Xtrriage
Anisotropy	0.786	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.29 , 29.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	0.047 for h,-k,-l	Xtrriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	14927	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	19.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.22% of the height of the origin peak. No significant pseudotranslation is detected.*

<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, 0.333 respectively for untwinned datasets, and 0.375, 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, NAG

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	0.45	2/7315 (0.0%)	0.62	2/9929 (0.0%)
1	B	0.46	4/7298 (0.1%)	0.62	1/9906 (0.0%)
All	All	0.46	6/14613 (0.0%)	0.62	3/19835 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2
1	B	0	2
All	All	0	4

The worst 5 of 6 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	350	GLU	CG-CD	-8.99	1.38	1.51
1	B	909	GLU	CD-OE2	-6.56	1.18	1.25
1	A	909	GLU	CD-OE2	-6.25	1.18	1.25
1	B	909	GLU	CD-OE1	-5.75	1.19	1.25
1	A	909	GLU	CD-OE1	-5.74	1.19	1.25

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	909	GLU	CA-CB-CG	6.34	127.35	113.40
1	A	909	GLU	CA-CB-CG	6.13	126.88	113.40
1	A	140	SER	N-CA-C	-5.50	96.16	111.00

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	139	ASP	Peptide
1	A	93	ALA	Peptide
1	B	574	SER	Peptide
1	B	93	ALA	Peptide

## 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	7154	0	6924	184	1
1	B	7138	0	6904	193	1
2	A	126	0	117	5	0
2	B	126	0	117	7	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	216	0	0	3	0
4	B	165	0	0	10	0
All	All	14927	0	14062	383	1

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 13.

The worst 5 of 383 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:542:GLN:NE2	1:A:576:PHE:HB2	1.58	1.17
1:B:947:TRP:O	1:B:951:ASN:ND2	1.81	1.14
1:A:947:TRP:O	1:A:951:ASN:ND2	1.84	1.08
1:B:187:GLY:HA2	1:B:207:MSE:HE1	1.34	1.06
1:A:187:GLY:HA2	1:A:207:MSE:HE1	1.35	1.05

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:521:ASP:O	1:B:141:GLN:OE1[1_656]	2.02	0.18

### 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	881/968 (91%)	850 (96%)	31 (4%)	0	100	100
1	B	879/968 (91%)	848 (96%)	31 (4%)	0	100	100
All	All	1760/1936 (91%)	1698 (96%)	62 (4%)	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	787/830 (95%)	708 (90%)	79 (10%)	7	15
1	B	786/830 (95%)	705 (90%)	81 (10%)	7	14
All	All	1573/1660 (95%)	1413 (90%)	160 (10%)	7	14

5 of 160 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	909	GLU
1	B	152	GLU
1	B	880	ASN

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	A	915	LEU
1	B	96	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 6 such sidechains are listed below:

Mol	Chain	Res	Type
1	A	542	GLN
1	B	880	ASN
1	A	880	ASN
1	A	351	ASN
1	B	351	ASN

### 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 monosaccharides in this entry.

### 5.6 Ligand geometry [i](#)

Of 20 ligands modelled in this entry, 2 are monoatomic - leaving 18 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	$\# Z  > 2$
2	NAG	B	1003	1	14,14,15	0.58	0	17,19,21	0.74	1 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	NAG	B	1008	1	14,14,15	0.68	0	17,19,21	1.48	3 (17%)
2	NAG	A	1009	1	14,14,15	0.53	0	17,19,21	1.42	1 (5%)
2	NAG	B	1007	1	14,14,15	0.48	0	17,19,21	1.36	1 (5%)
2	NAG	A	1002	1	14,14,15	0.57	0	17,19,21	0.92	0
2	NAG	B	1006	1	14,14,15	0.41	0	17,19,21	1.59	3 (17%)
2	NAG	A	1006	1	14,14,15	0.46	0	17,19,21	1.28	1 (5%)
2	NAG	A	1001	1	14,14,15	0.67	0	17,19,21	1.09	2 (11%)
2	NAG	B	1001	1	14,14,15	0.63	0	17,19,21	0.79	0
2	NAG	A	1008	1	14,14,15	0.56	0	17,19,21	0.85	0
2	NAG	A	1003	1	14,14,15	0.53	0	17,19,21	0.89	1 (5%)
2	NAG	B	1005	1	14,14,15	0.48	0	17,19,21	0.89	0
2	NAG	B	1002	1	14,14,15	0.62	0	17,19,21	1.21	3 (17%)
2	NAG	A	1007	1	14,14,15	0.52	0	17,19,21	1.43	1 (5%)
2	NAG	B	1009	1	14,14,15	0.57	0	17,19,21	1.24	1 (5%)
2	NAG	A	1004	1	14,14,15	0.50	0	17,19,21	1.31	3 (17%)
2	NAG	B	1004	1	14,14,15	0.50	0	17,19,21	0.63	0
2	NAG	A	1005	1	14,14,15	0.60	0	17,19,21	1.33	3 (17%)

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
2	NAG	B	1003	1	-	4/6/23/26	0/1/1/1
2	NAG	B	1008	1	-	0/6/23/26	0/1/1/1
2	NAG	A	1009	1	-	2/6/23/26	0/1/1/1
2	NAG	B	1007	1	-	0/6/23/26	0/1/1/1
2	NAG	A	1002	1	-	2/6/23/26	0/1/1/1
2	NAG	B	1006	1	-	0/6/23/26	0/1/1/1
2	NAG	A	1006	1	-	0/6/23/26	0/1/1/1
2	NAG	A	1001	1	-	4/6/23/26	0/1/1/1
2	NAG	B	1001	1	-	2/6/23/26	0/1/1/1
2	NAG	A	1008	1	-	0/6/23/26	0/1/1/1
2	NAG	A	1003	1	-	2/6/23/26	0/1/1/1
2	NAG	B	1005	1	-	4/6/23/26	0/1/1/1
2	NAG	B	1002	1	-	4/6/23/26	0/1/1/1
2	NAG	A	1007	1	-	0/6/23/26	0/1/1/1

Continued on next page...

*Continued from previous page...*

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	B	1009	1	-	2/6/23/26	0/1/1/1
2	NAG	A	1004	1	-	2/6/23/26	0/1/1/1
2	NAG	B	1004	1	-	3/6/23/26	0/1/1/1
2	NAG	A	1005	1	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

The worst 5 of 24 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1009	NAG	C1-O5-C5	4.95	118.90	112.19
2	A	1007	NAG	C1-O5-C5	4.67	118.52	112.19
2	B	1007	NAG	C1-O5-C5	4.26	117.96	112.19
2	B	1006	NAG	C2-N2-C7	-4.12	117.03	122.90
2	B	1009	NAG	C1-O5-C5	4.01	117.63	112.19

There are no chirality outliers.

5 of 31 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	1003	NAG	C8-C7-N2-C2
2	B	1003	NAG	O7-C7-N2-C2
2	A	1009	NAG	C8-C7-N2-C2
2	A	1009	NAG	O7-C7-N2-C2
2	A	1002	NAG	C8-C7-N2-C2

There are no ring outliers.

10 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1003	NAG	1	0
2	B	1008	NAG	2	0
2	A	1009	NAG	1	0
2	A	1002	NAG	1	0
2	A	1001	NAG	1	0
2	B	1001	NAG	1	0
2	A	1003	NAG	1	0
2	B	1002	NAG	3	0
2	B	1009	NAG	3	0
2	A	1005	NAG	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, 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	865/968 (89%)	-0.35	17 (1%) 65 68	4, 16, 43, 84	1 (0%)
1	B	863/968 (89%)	-0.32	16 (1%) 66 69	4, 16, 44, 83	1 (0%)
All	All	1728/1936 (89%)	-0.34	33 (1%) 66 69	4, 16, 44, 84	2 (0%)

The worst 5 of 33 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	61	LEU	7.2
1	A	93	ALA	5.7
1	B	925	VAL	5.2
1	B	137	VAL	5.0
1	A	94	ASP	4.5

### 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 monosaccharides 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. 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	B-factors( $\text{\AA}^2$ )	Q<0.9
2	NAG	B	1004	14/15	0.82	0.27	30,56,62,62	0
2	NAG	B	1005	14/15	0.87	0.28	44,62,72,74	0
2	NAG	B	1003	14/15	0.87	0.23	38,49,59,61	0
2	NAG	A	1008	14/15	0.88	0.22	37,48,57,67	0
2	NAG	A	1003	14/15	0.88	0.20	32,43,52,55	0
2	NAG	A	1001	14/15	0.88	0.18	38,46,54,57	0
2	NAG	B	1001	14/15	0.88	0.28	30,50,62,64	0
2	NAG	A	1004	14/15	0.89	0.19	26,41,50,53	0
2	NAG	B	1008	14/15	0.90	0.20	28,38,48,51	0
2	NAG	B	1009	14/15	0.92	0.23	27,40,56,59	0
2	NAG	A	1009	14/15	0.92	0.23	25,34,43,54	0
2	NAG	B	1002	14/15	0.92	0.18	12,18,26,26	0
2	NAG	A	1005	14/15	0.92	0.19	13,27,44,47	0
2	NAG	B	1007	14/15	0.93	0.16	24,29,40,42	0
2	NAG	B	1006	14/15	0.94	0.12	18,24,29,49	0
2	NAG	A	1006	14/15	0.95	0.10	14,18,24,26	0
2	NAG	A	1007	14/15	0.95	0.18	19,26,35,37	0
2	NAG	A	1002	14/15	0.96	0.11	9,14,19,24	0
3	ZN	B	1010	1/1	0.99	0.04	18,18,18,18	0
3	ZN	A	1010	1/1	1.00	0.03	17,17,17,17	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.