



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 14, 2026 – 11:54 AM UTC

PDB ID : 6OIU / pdb_00006oiu
Title : X-ray crystal structure of the ectodomain of the Toxoplasma gondii ME49 Aminopeptidase N (TGME49_224350)
Authors : McGowan, S.; Drinkwater, N.
Deposited on : 2019-04-09
Resolution : 2.20 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

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

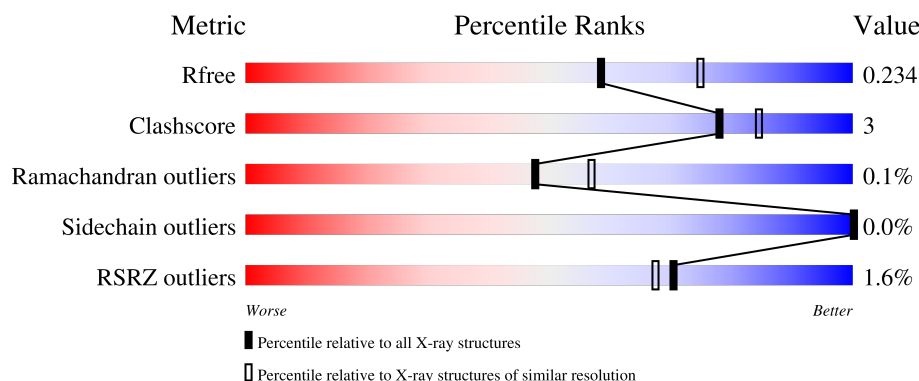
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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}	180053	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 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	907	90%8%
1	B	907	92%7%
1	C	907	94%6%
1	D	907	3% 89%8%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 30397 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
1	A	890	Total	C	N	O	S	0	0	0
			6951	4405	1212	1305	29			
1	B	899	Total	C	N	O	S	0	0	0
			6944	4403	1201	1311	29			
1	C	900	Total	C	N	O	S	0	0	0
			6998	4438	1212	1319	29			
1	D	886	Total	C	N	O	S	0	0	0
			6696	4244	1161	1262	29			

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	519	MET	-	initiating methionine	UNP S8G5K8
A	840	GLN	ASN	conflict	UNP S8G5K8
A	1303	GLN	ASN	conflict	UNP S8G5K8
A	1420	HIS	-	expression tag	UNP S8G5K8
A	1421	HIS	-	expression tag	UNP S8G5K8
A	1422	HIS	-	expression tag	UNP S8G5K8
A	1423	HIS	-	expression tag	UNP S8G5K8
A	1424	HIS	-	expression tag	UNP S8G5K8
A	1425	HIS	-	expression tag	UNP S8G5K8
B	519	MET	-	initiating methionine	UNP S8G5K8
B	840	GLN	ASN	conflict	UNP S8G5K8
B	1303	GLN	ASN	conflict	UNP S8G5K8
B	1420	HIS	-	expression tag	UNP S8G5K8
B	1421	HIS	-	expression tag	UNP S8G5K8
B	1422	HIS	-	expression tag	UNP S8G5K8
B	1423	HIS	-	expression tag	UNP S8G5K8
B	1424	HIS	-	expression tag	UNP S8G5K8
B	1425	HIS	-	expression tag	UNP S8G5K8
C	519	MET	-	initiating methionine	UNP S8G5K8
C	840	GLN	ASN	conflict	UNP S8G5K8
C	1303	GLN	ASN	conflict	UNP S8G5K8

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Chain	Residue	Modelled	Actual	Comment	Reference
C	1420	HIS	-	expression tag	UNP S8G5K8
C	1421	HIS	-	expression tag	UNP S8G5K8
C	1422	HIS	-	expression tag	UNP S8G5K8
C	1423	HIS	-	expression tag	UNP S8G5K8
C	1424	HIS	-	expression tag	UNP S8G5K8
C	1425	HIS	-	expression tag	UNP S8G5K8
D	519	MET	-	initiating methionine	UNP S8G5K8
D	840	GLN	ASN	conflict	UNP S8G5K8
D	1303	GLN	ASN	conflict	UNP S8G5K8
D	1420	HIS	-	expression tag	UNP S8G5K8
D	1421	HIS	-	expression tag	UNP S8G5K8
D	1422	HIS	-	expression tag	UNP S8G5K8
D	1423	HIS	-	expression tag	UNP S8G5K8
D	1424	HIS	-	expression tag	UNP S8G5K8
D	1425	HIS	-	expression tag	UNP S8G5K8

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

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

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



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

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	717	Total	O	0	0
			717	717		
4	B	750	Total	O	0	0
			750	750		
4	C	872	Total	O	0	0
			872	872		

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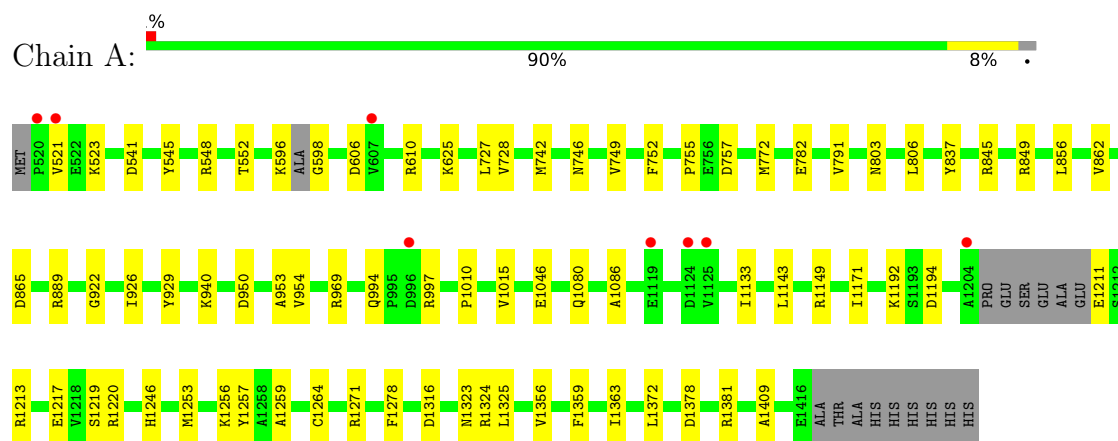
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	407	Total 407	O 407	0	0

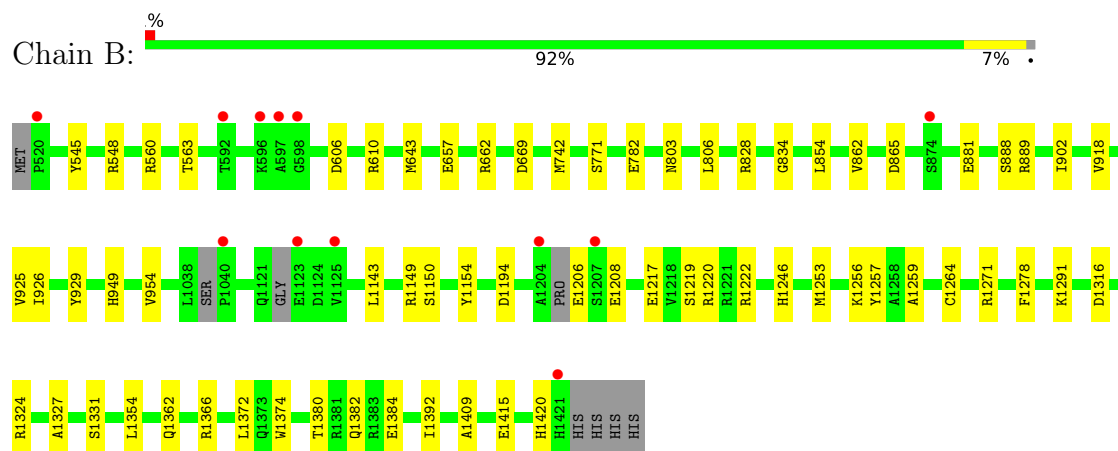
3 Residue-property plots [i](#)

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.

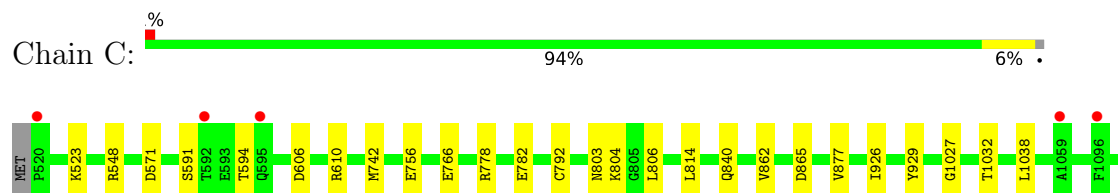
• Molecule 1: Aminopeptidase N



• Molecule 1: Aminopeptidase N

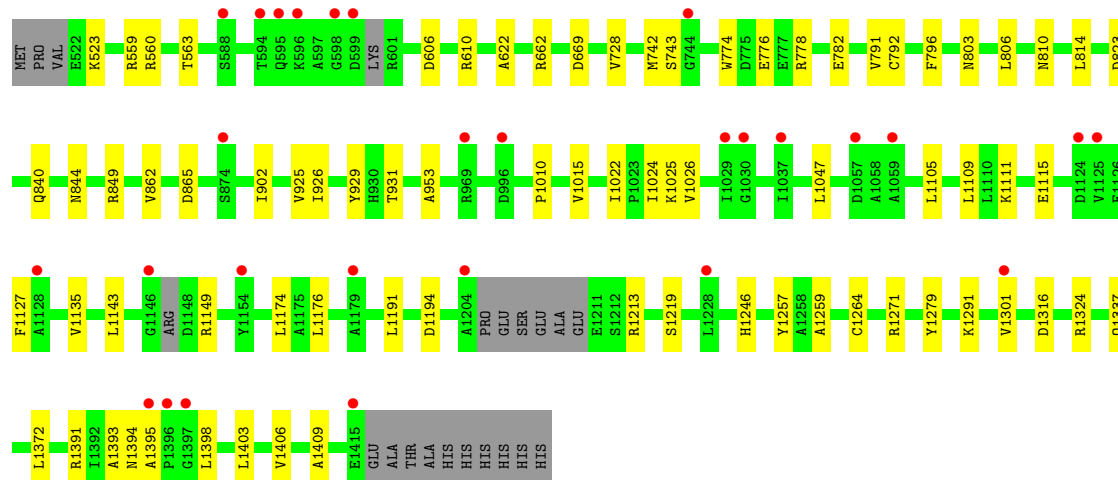
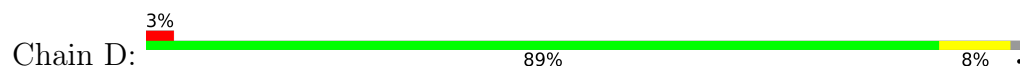


• Molecule 1: Aminopeptidase N





• Molecule 1: Aminopeptidase N



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	92.93Å 207.87Å 102.87Å 90.00° 94.34° 90.00°	Depositor
Resolution (Å)	38.54 – 2.20 38.54 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.6 (38.54-2.20) 91.8 (38.54-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.41 (at 2.20Å)	Xtriage
Refinement program	PHENIX 1.14_3260	Depositor
R, R_{free}	0.190 , 0.236 0.191 , 0.234	Depositor DCC
R_{free} test set	9637 reflections (4.90%)	wwPDB-VP
Wilson B-factor (Å ²)	19.0	Xtriage
Anisotropy	0.717	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 44.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	30397	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.50% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, GOL

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	0.22	3/7091 (0.0%)	0.29	0/9602
1	B	0.08	0/7085	0.26	0/9607
1	C	0.20	3/7142 (0.0%)	0.28	0/9682
1	D	0.24	3/6833 (0.0%)	0.30	0/9289
All	All	0.20	9/28151 (0.0%)	0.28	0/38180

The worst 5 of 9 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	1393	ALA	C-N	-8.96	1.22	1.33
1	D	1394	ASN	C-N	8.75	1.49	1.33
1	C	756	GLU	C-O	-7.45	1.14	1.24
1	A	755	PRO	C-O	-6.48	1.15	1.24
1	D	1127	PHE	C-O	-5.87	1.17	1.23

There are no bond angle outliers.

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	6951	0	6763	48	0
1	B	6944	0	6668	36	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	6998	0	6776	31	0
1	D	6696	0	6273	38	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0
3	A	18	0	24	0	0
3	B	18	0	24	0	0
3	C	18	0	24	3	0
4	A	717	0	0	9	0
4	B	750	0	0	2	0
4	C	872	0	0	9	0
4	D	407	0	0	2	0
All	All	30397	0	26552	154	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 3.

The worst 5 of 154 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:C:1032:THR:HG23	4:C:2296:HOH:O	1.68	0.92
1:D:1149:ARG:HB2	1:D:1219:SER:HB3	1.67	0.74
1:C:742:MET:HE2	1:C:782:GLU:HB2	1.71	0.72
1:A:596:LYS:HA	1:A:598:GLY:N	2.06	0.70
1:A:889:ARG:NH1	4:A:2102:HOH:O	2.25	0.70

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	884/907 (98%)	864 (98%)	20 (2%)	0	100	100
1	B	891/907 (98%)	869 (98%)	22 (2%)	0	100	100
1	C	896/907 (99%)	875 (98%)	19 (2%)	2 (0%)	43	51
1	D	878/907 (97%)	851 (97%)	27 (3%)	0	100	100
All	All	3549/3628 (98%)	3459 (98%)	88 (2%)	2 (0%)	48	57

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	1123	GLU
1	C	1205	PRO

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	717/775 (92%)	717 (100%)	0	100	100
1	B	705/775 (91%)	705 (100%)	0	100	100
1	C	719/775 (93%)	719 (100%)	0	100	100
1	D	653/775 (84%)	652 (100%)	1 (0%)	87	94
All	All	2794/3100 (90%)	2793 (100%)	1 (0%)	100	100

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	1174	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 22 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	1121	GLN
1	D	1078	HIS

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Mol	Chain	Res	Type
1	C	1296	GLN
1	D	1112	HIS
1	B	1080	GLN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 17 ligands modelled in this entry, 8 are monoatomic - leaving 9 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	A	2004	-	5,5,5	0.94	0	5,5,5	1.07	0
3	GOL	A	2003	-	5,5,5	0.93	0	5,5,5	1.07	0
3	GOL	B	1504	-	5,5,5	0.93	0	5,5,5	1.07	0
3	GOL	C	1504	-	5,5,5	0.92	0	5,5,5	1.09	0
3	GOL	C	1505	-	5,5,5	0.94	0	5,5,5	1.05	0
3	GOL	A	2005	-	5,5,5	0.92	0	5,5,5	1.10	0
3	GOL	B	1505	-	5,5,5	0.92	0	5,5,5	1.10	0
3	GOL	C	1503	-	5,5,5	0.94	0	5,5,5	1.09	0
3	GOL	B	1503	-	5,5,5	0.95	0	5,5,5	1.13	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	A	2004	-	-	2/4/4/4	-
3	GOL	A	2003	-	-	2/4/4/4	-
3	GOL	B	1504	-	-	2/4/4/4	-
3	GOL	C	1504	-	-	2/4/4/4	-
3	GOL	C	1505	-	-	1/4/4/4	-
3	GOL	A	2005	-	-	0/4/4/4	-
3	GOL	B	1505	-	-	0/4/4/4	-
3	GOL	C	1503	-	-	0/4/4/4	-
3	GOL	B	1503	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

5 of 11 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	2003	GOL	C1-C2-C3-O3
3	B	1503	GOL	C1-C2-C3-O3
3	B	1503	GOL	O2-C2-C3-O3
3	B	1504	GOL	C1-C2-C3-O3
3	A	2004	GOL	O1-C1-C2-C3

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	1505	GOL	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	890/907 (98%)	-0.05	8 (0%) 81 79	12, 25, 42, 68	1 (0%)
1	B	899/907 (99%)	-0.05	12 (1%) 75 73	14, 26, 42, 70	2 (0%)
1	C	900/907 (99%)	-0.20	10 (1%) 78 76	11, 21, 41, 76	0
1	D	886/907 (97%)	0.60	28 (3%) 50 47	22, 39, 59, 76	0
All	All	3575/3628 (98%)	0.07	58 (1%) 70 67	11, 27, 50, 76	3 (0%)

The worst 5 of 58 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	520	PRO	4.3
1	D	1146	GLY	4.0
1	D	1204	ALA	3.9
1	C	1210	GLU	3.9
1	D	596	LYS	3.7

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 oligosaccharides 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, 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	B-factors($\text{\AA}^2$)	Q<0.9
3	GOL	B	1503	6/6	0.77	0.24	37,38,43,45	0
3	GOL	C	1503	6/6	0.81	0.16	31,32,38,45	0
3	GOL	B	1504	6/6	0.82	0.15	38,44,51,54	0
3	GOL	A	2004	6/6	0.82	0.14	34,36,42,44	0
3	GOL	A	2003	6/6	0.83	0.12	34,35,36,38	0
3	GOL	A	2005	6/6	0.84	0.15	37,39,41,48	0
3	GOL	C	1505	6/6	0.84	0.13	32,33,39,44	0
3	GOL	C	1504	6/6	0.86	0.11	25,34,39,42	0
3	GOL	B	1505	6/6	0.88	0.12	36,39,44,49	0
2	ZN	B	1502	1/1	0.93	0.12	62,62,62,62	0
2	ZN	A	2002	1/1	0.93	0.10	56,56,56,56	0
2	ZN	C	1502	1/1	0.95	0.08	51,51,51,51	0
2	ZN	D	1502	1/1	0.96	0.12	78,78,78,78	0
2	ZN	D	1501	1/1	0.97	0.04	32,32,32,32	0
2	ZN	B	1501	1/1	0.99	0.03	18,18,18,18	0
2	ZN	C	1501	1/1	0.99	0.01	16,16,16,16	0
2	ZN	A	2001	1/1	1.00	0.01	16,16,16,16	0

6.5 Other polymers ⓘ

There are no such residues in this entry.