



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:40 PM UTC

PDB ID : 7K9O / pdb_00007k9o
Title : Co-crystal structure of alpha glucosidase with compound 3
Authors : Karade, S.S.; Mariuzza, R.A.
Deposited on : 2020-09-29
Resolution : 2.30 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<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
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
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

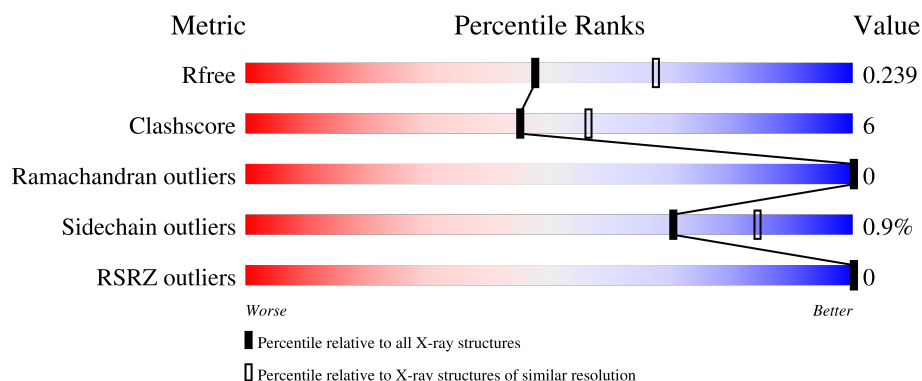
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	977	
1	C	977	
2	B	547	
2	D	547	

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 15936 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 Alpha glucosidase 2 alpha neutral subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	854	Total	C	N	O	S	0	10	0
			6925	4441	1194	1260	30			
1	C	856	Total	C	N	O	S	0	11	0
			6939	4450	1194	1265	30			

There are 88 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	2	MET	-	initiating methionine	UNP A1A4T2
A	3	GLY	-	expression tag	UNP A1A4T2
A	4	ILE	-	expression tag	UNP A1A4T2
A	5	LEU	-	expression tag	UNP A1A4T2
A	6	PRO	-	expression tag	UNP A1A4T2
A	7	SER	-	expression tag	UNP A1A4T2
A	8	PRO	-	expression tag	UNP A1A4T2
A	9	GLY	-	expression tag	UNP A1A4T2
A	10	MET	-	expression tag	UNP A1A4T2
A	11	PRO	-	expression tag	UNP A1A4T2
A	12	ALA	-	expression tag	UNP A1A4T2
A	13	LEU	-	expression tag	UNP A1A4T2
A	14	LEU	-	expression tag	UNP A1A4T2
A	15	SER	-	expression tag	UNP A1A4T2
A	16	LEU	-	expression tag	UNP A1A4T2
A	17	VAL	-	expression tag	UNP A1A4T2
A	18	SER	-	expression tag	UNP A1A4T2
A	19	LEU	-	expression tag	UNP A1A4T2
A	20	LEU	-	expression tag	UNP A1A4T2
A	21	SER	-	expression tag	UNP A1A4T2
A	22	VAL	-	expression tag	UNP A1A4T2
A	23	LEU	-	expression tag	UNP A1A4T2
A	24	LEU	-	expression tag	UNP A1A4T2
A	25	MET	-	expression tag	UNP A1A4T2
A	26	GLY	-	expression tag	UNP A1A4T2

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Chain	Residue	Modelled	Actual	Comment	Reference
A	27	CYS	-	expression tag	UNP A1A4T2
A	28	VAL	-	expression tag	UNP A1A4T2
A	29	ALA	-	expression tag	UNP A1A4T2
A	30	GLU	-	expression tag	UNP A1A4T2
A	31	THR	-	expression tag	UNP A1A4T2
A	32	GLY	-	expression tag	UNP A1A4T2
A	97	ASP	ASN	engineered mutation	UNP A1A4T2
A	967	SER	-	expression tag	UNP A1A4T2
A	968	ALA	-	expression tag	UNP A1A4T2
A	969	TRP	-	expression tag	UNP A1A4T2
A	970	SER	-	expression tag	UNP A1A4T2
A	971	HIS	-	expression tag	UNP A1A4T2
A	972	PRO	-	expression tag	UNP A1A4T2
A	973	GLN	-	expression tag	UNP A1A4T2
A	974	PHE	-	expression tag	UNP A1A4T2
A	975	GLU	-	expression tag	UNP A1A4T2
A	976	LYS	-	expression tag	UNP A1A4T2
A	977	LEU	-	expression tag	UNP A1A4T2
A	978	GLU	-	expression tag	UNP A1A4T2
C	2	MET	-	initiating methionine	UNP A1A4T2
C	3	GLY	-	expression tag	UNP A1A4T2
C	4	ILE	-	expression tag	UNP A1A4T2
C	5	LEU	-	expression tag	UNP A1A4T2
C	6	PRO	-	expression tag	UNP A1A4T2
C	7	SER	-	expression tag	UNP A1A4T2
C	8	PRO	-	expression tag	UNP A1A4T2
C	9	GLY	-	expression tag	UNP A1A4T2
C	10	MET	-	expression tag	UNP A1A4T2
C	11	PRO	-	expression tag	UNP A1A4T2
C	12	ALA	-	expression tag	UNP A1A4T2
C	13	LEU	-	expression tag	UNP A1A4T2
C	14	LEU	-	expression tag	UNP A1A4T2
C	15	SER	-	expression tag	UNP A1A4T2
C	16	LEU	-	expression tag	UNP A1A4T2
C	17	VAL	-	expression tag	UNP A1A4T2
C	18	SER	-	expression tag	UNP A1A4T2
C	19	LEU	-	expression tag	UNP A1A4T2
C	20	LEU	-	expression tag	UNP A1A4T2
C	21	SER	-	expression tag	UNP A1A4T2
C	22	VAL	-	expression tag	UNP A1A4T2
C	23	LEU	-	expression tag	UNP A1A4T2
C	24	LEU	-	expression tag	UNP A1A4T2

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Chain	Residue	Modelled	Actual	Comment	Reference
C	25	MET	-	expression tag	UNP A1A4T2
C	26	GLY	-	expression tag	UNP A1A4T2
C	27	CYS	-	expression tag	UNP A1A4T2
C	28	VAL	-	expression tag	UNP A1A4T2
C	29	ALA	-	expression tag	UNP A1A4T2
C	30	GLU	-	expression tag	UNP A1A4T2
C	31	THR	-	expression tag	UNP A1A4T2
C	32	GLY	-	expression tag	UNP A1A4T2
C	97	ASP	ASN	engineered mutation	UNP A1A4T2
C	967	SER	-	expression tag	UNP A1A4T2
C	968	ALA	-	expression tag	UNP A1A4T2
C	969	TRP	-	expression tag	UNP A1A4T2
C	970	SER	-	expression tag	UNP A1A4T2
C	971	HIS	-	expression tag	UNP A1A4T2
C	972	PRO	-	expression tag	UNP A1A4T2
C	973	GLN	-	expression tag	UNP A1A4T2
C	974	PHE	-	expression tag	UNP A1A4T2
C	975	GLU	-	expression tag	UNP A1A4T2
C	976	LYS	-	expression tag	UNP A1A4T2
C	977	LEU	-	expression tag	UNP A1A4T2
C	978	GLU	-	expression tag	UNP A1A4T2

- Molecule 2 is a protein called Glucosidase 2 subunit beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	83	Total	C	N	O	S	0	0	0
			603	359	98	136	10			
2	D	85	Total	C	N	O	S	0	0	0
			607	362	97	138	10			

There are 88 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-16	MET	-	initiating methionine	UNP O08795
B	-15	GLY	-	expression tag	UNP O08795
B	-14	ILE	-	expression tag	UNP O08795
B	-13	LEU	-	expression tag	UNP O08795
B	-12	PRO	-	expression tag	UNP O08795
B	-11	SER	-	expression tag	UNP O08795
B	-10	PRO	-	expression tag	UNP O08795
B	-9	GLY	-	expression tag	UNP O08795
B	-8	MET	-	expression tag	UNP O08795

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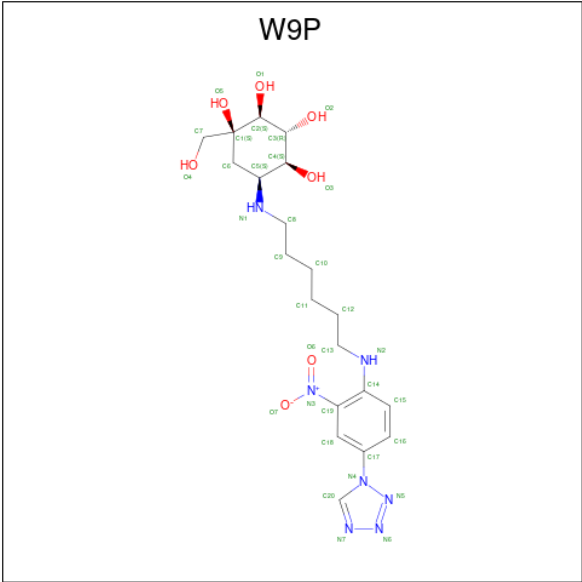
Chain	Residue	Modelled	Actual	Comment	Reference
B	-7	PRO	-	expression tag	UNP O08795
B	-6	ALA	-	expression tag	UNP O08795
B	-5	LEU	-	expression tag	UNP O08795
B	-4	LEU	-	expression tag	UNP O08795
B	-3	SER	-	expression tag	UNP O08795
B	-2	LEU	-	expression tag	UNP O08795
B	-1	VAL	-	expression tag	UNP O08795
B	0	SER	-	expression tag	UNP O08795
B	1	LEU	-	expression tag	UNP O08795
B	2	LEU	-	expression tag	UNP O08795
B	3	SER	-	expression tag	UNP O08795
B	4	VAL	-	expression tag	UNP O08795
B	5	LEU	-	expression tag	UNP O08795
B	6	LEU	-	expression tag	UNP O08795
B	7	MET	-	expression tag	UNP O08795
B	8	GLY	-	expression tag	UNP O08795
B	9	CYS	-	expression tag	UNP O08795
B	10	VAL	-	expression tag	UNP O08795
B	11	ALA	-	expression tag	UNP O08795
B	12	GLU	-	expression tag	UNP O08795
B	13	THR	-	expression tag	UNP O08795
B	14	GLY	-	expression tag	UNP O08795
B	518	SER	-	expression tag	UNP O08795
B	519	ALA	-	expression tag	UNP O08795
B	520	TRP	-	expression tag	UNP O08795
B	521	LEU	-	expression tag	UNP O08795
B	522	GLU	-	expression tag	UNP O08795
B	523	THR	-	expression tag	UNP O08795
B	524	LYS	-	expression tag	UNP O08795
B	525	HIS	-	expression tag	UNP O08795
B	526	HIS	-	expression tag	UNP O08795
B	527	HIS	-	expression tag	UNP O08795
B	528	HIS	-	expression tag	UNP O08795
B	529	HIS	-	expression tag	UNP O08795
B	530	HIS	-	expression tag	UNP O08795
D	-16	MET	-	initiating methionine	UNP O08795
D	-15	GLY	-	expression tag	UNP O08795
D	-14	ILE	-	expression tag	UNP O08795
D	-13	LEU	-	expression tag	UNP O08795
D	-12	PRO	-	expression tag	UNP O08795
D	-11	SER	-	expression tag	UNP O08795
D	-10	PRO	-	expression tag	UNP O08795

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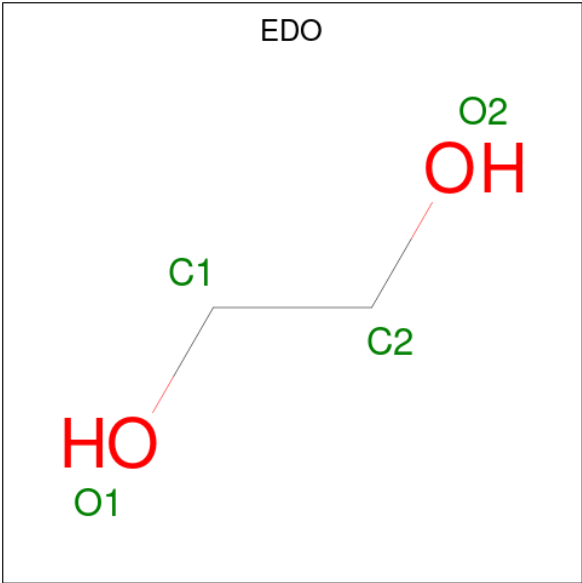
Chain	Residue	Modelled	Actual	Comment	Reference
D	-9	GLY	-	expression tag	UNP O08795
D	-8	MET	-	expression tag	UNP O08795
D	-7	PRO	-	expression tag	UNP O08795
D	-6	ALA	-	expression tag	UNP O08795
D	-5	LEU	-	expression tag	UNP O08795
D	-4	LEU	-	expression tag	UNP O08795
D	-3	SER	-	expression tag	UNP O08795
D	-2	LEU	-	expression tag	UNP O08795
D	-1	VAL	-	expression tag	UNP O08795
D	0	SER	-	expression tag	UNP O08795
D	1	LEU	-	expression tag	UNP O08795
D	2	LEU	-	expression tag	UNP O08795
D	3	SER	-	expression tag	UNP O08795
D	4	VAL	-	expression tag	UNP O08795
D	5	LEU	-	expression tag	UNP O08795
D	6	LEU	-	expression tag	UNP O08795
D	7	MET	-	expression tag	UNP O08795
D	8	GLY	-	expression tag	UNP O08795
D	9	CYS	-	expression tag	UNP O08795
D	10	VAL	-	expression tag	UNP O08795
D	11	ALA	-	expression tag	UNP O08795
D	12	GLU	-	expression tag	UNP O08795
D	13	THR	-	expression tag	UNP O08795
D	14	GLY	-	expression tag	UNP O08795
D	518	SER	-	expression tag	UNP O08795
D	519	ALA	-	expression tag	UNP O08795
D	520	TRP	-	expression tag	UNP O08795
D	521	LEU	-	expression tag	UNP O08795
D	522	GLU	-	expression tag	UNP O08795
D	523	THR	-	expression tag	UNP O08795
D	524	LYS	-	expression tag	UNP O08795
D	525	HIS	-	expression tag	UNP O08795
D	526	HIS	-	expression tag	UNP O08795
D	527	HIS	-	expression tag	UNP O08795
D	528	HIS	-	expression tag	UNP O08795
D	529	HIS	-	expression tag	UNP O08795
D	530	HIS	-	expression tag	UNP O08795

- Molecule 3 is (1 {S},2 {S},3 {R},4 {S},5 {S})-1-(hydroxymethyl)-5-[6-[[2-[oxidanyl(oxid anylidene)-\$l^{4}\$-azanyl]-4-(1,2,3,4-tetrazol-1-yl)phenyl]amino]hexylamino]cyclohexane -1,2,3,4-tetrol (CCD ID: W9P) (formula: C₂₀H₃₁N₇O₇) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			34	20	7	7		
3	C	1	Total	C	N	O	0	0
			34	20	7	7		

- Molecule 4 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			4	2	2		
4	A	1	Total	C	O	0	0
			4	2	2		

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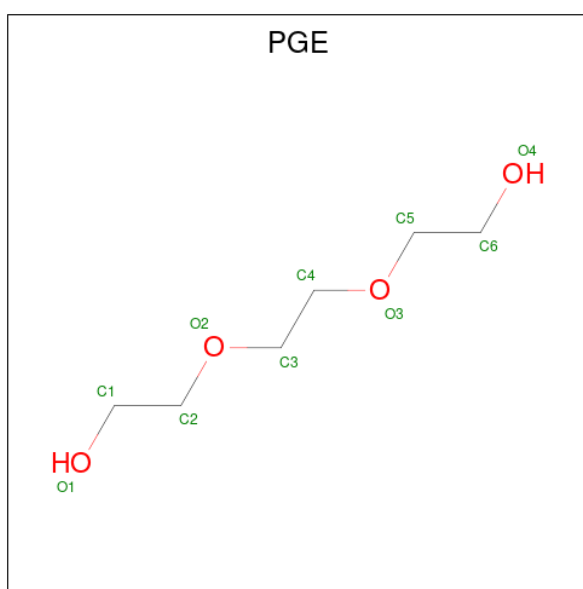
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total 4	C 2	O 2	0	0
4	A	1	Total 4	C 2	O 2	0	0
4	A	1	Total 4	C 2	O 2	0	0
4	A	1	Total 4	C 2	O 2	0	0
4	A	1	Total 4	C 2	O 2	0	0
4	A	1	Total 4	C 2	O 2	0	0
4	A	1	Total 4	C 2	O 2	0	0
4	A	1	Total 4	C 2	O 2	0	0
4	A	1	Total 4	C 2	O 2	0	0
4	A	1	Total 4	C 2	O 2	0	0
4	A	1	Total 4	C 2	O 2	0	0
4	A	1	Total 4	C 2	O 2	0	0
4	A	1	Total 4	C 2	O 2	0	0
4	B	1	Total 4	C 2	O 2	0	0
4	C	1	Total 4	C 2	O 2	0	0
4	C	1	Total 4	C 2	O 2	0	0
4	C	1	Total 4	C 2	O 2	0	0
4	C	1	Total 4	C 2	O 2	0	0
4	C	1	Total 4	C 2	O 2	0	0
4	C	1	Total 4	C 2	O 2	0	0
4	C	1	Total 4	C 2	O 2	0	0
4	C	1	Total 4	C 2	O 2	0	0

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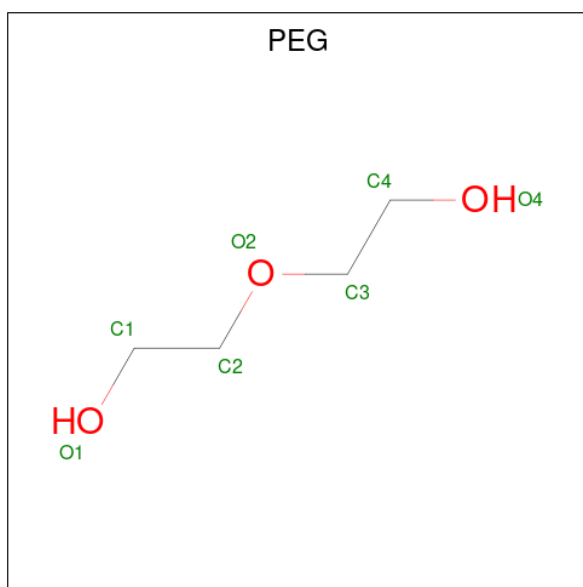
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	C	1	Total	C	O	0	0
			4	2	2		
4	C	1	Total	C	O	0	0
			4	2	2		
4	C	1	Total	C	O	0	0
			4	2	2		
4	C	1	Total	C	O	0	0
			4	2	2		

- Molecule 5 is TRIETHYLENE GLYCOL (CCD ID: PGE) (formula: $C_6H_{14}O_4$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			10	6	4		
5	A	1	Total	C	O	0	0
			10	6	4		
5	C	1	Total	C	O	0	0
			10	6	4		

- Molecule 6 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			7	4	3		
6	A	1	Total	C	O	0	0
			7	4	3		
6	A	1	Total	C	O	0	0
			7	4	3		
6	C	1	Total	C	O	0	0
			7	4	3		

- Molecule 7 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	B	2	Total	Ca	0	0
			2	2		
7	D	2	Total	Ca	0	0
			2	2		

- Molecule 8 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	C	1	Total	O	S	0	0
			5	4	1		

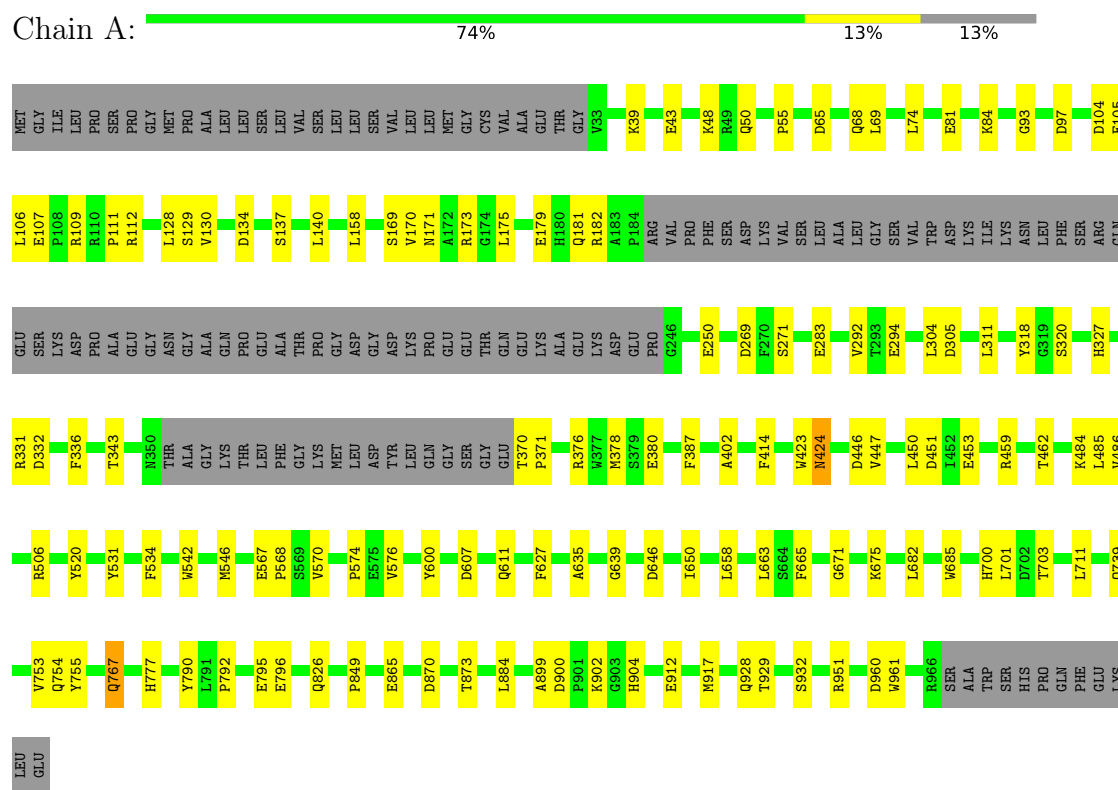
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	332	Total	O	0	0
			332	332		
9	B	14	Total	O	0	0
			14	14		
9	C	261	Total	O	0	0
			261	261		
9	D	12	Total	O	0	0
			12	12		

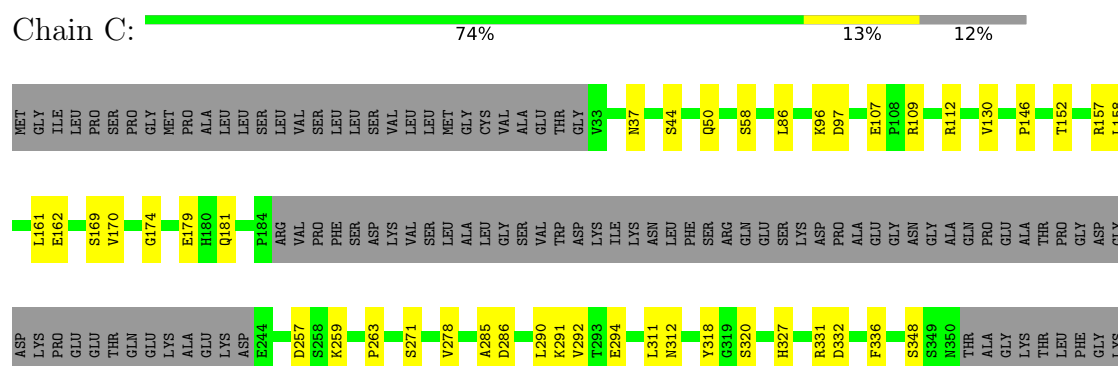
3 Residue-property plots

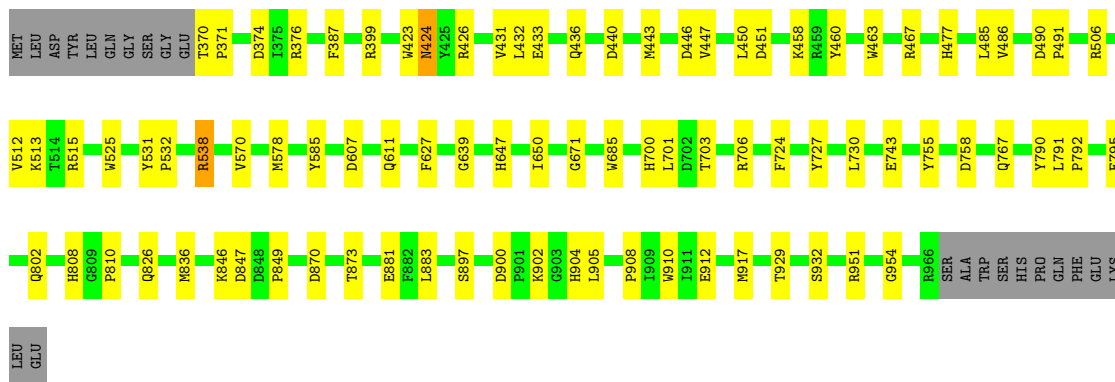
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Alpha glucosidase 2 alpha neutral subunit



- Molecule 1: Alpha glucosidase 2 alpha neutral subunit

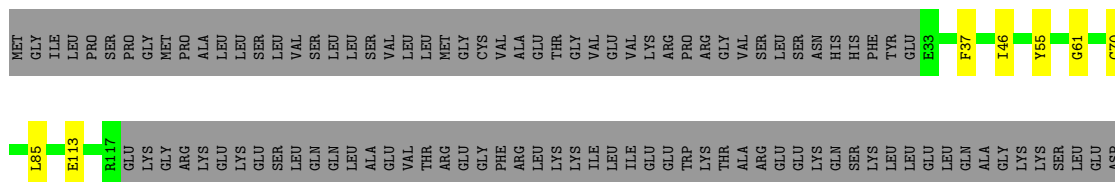




- Molecule 2: Glucosidase 2 subunit beta



- Molecule 2: Glucosidase 2 subunit beta



[illegible]

4 Data and refinement statistics

Property	Value	Source
Space group	P 32	Depositor
Cell constants a, b, c, α , β , γ	102.72Å 102.72Å 239.62Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	39.94 – 2.30 39.94 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.9 (39.94-2.30) 97.0 (39.94-2.30)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.59 (at 2.29Å)	Xtriage
Refinement program	PHENIX 1.14_3228	Depositor
R, R_{free}	0.210 , 0.238 0.211 , 0.239	Depositor DCC
R_{free} test set	2012 reflections (1.61%)	wwPDB-VP
Wilson B-factor (Å ²)	40.3	Xtriage
Anisotropy	0.044	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 26.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.39$, $\langle L^2 \rangle = 0.21$	Xtriage
Estimated twinning fraction	0.085 for -h,-k,l 0.247 for h,-h-k,-l 0.085 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	15936	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.48% 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: CA, PGE, EDO, W9P, 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	0.18	0/7165	0.41	0/9751
1	C	0.17	0/7183	0.40	0/9777
2	B	0.18	0/615	0.43	0/842
2	D	0.13	0/619	0.37	0/849
All	All	0.17	0/15582	0.41	0/21219

There are no bond length outliers.

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	6925	0	6689	90	1
1	C	6939	0	6692	79	1
2	B	603	0	500	5	0
2	D	607	0	493	5	0
3	A	34	0	0	0	0
3	C	34	0	0	0	0
4	A	56	0	84	4	0
4	B	4	0	6	0	0
4	C	48	0	72	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	20	0	28	5	0
5	C	10	0	14	1	0
6	A	21	0	29	1	0
6	C	7	0	10	1	0
7	B	2	0	0	0	0
7	D	2	0	0	0	0
8	C	5	0	0	0	0
9	A	332	0	0	25	0
9	B	14	0	0	1	0
9	C	261	0	0	22	0
9	D	12	0	0	1	0
All	All	15936	0	14617	177	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 6.

All (177) 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:463:TRP:O	9:C:1102:HOH:O	1.81	0.99
1:C:611:GLN:OE1	9:C:1101:HOH:O	1.80	0.97
1:C:802:GLN:NE2	9:C:1105:HOH:O	1.99	0.94
1:A:574:PRO:O	9:A:1101:HOH:O	1.90	0.89
1:C:263:PRO:O	9:C:1103:HOH:O	1.93	0.85
1:C:257:ASP:OD2	9:C:1104:HOH:O	1.94	0.84
1:A:65:ASP:O	9:A:1102:HOH:O	1.96	0.82
1:C:86:LEU:O	9:C:1106:HOH:O	1.99	0.79
1:C:37:ASN:O	9:C:1107:HOH:O	2.00	0.78
1:C:515:ARG:NE	9:C:1108:HOH:O	2.04	0.75
2:B:99:CYS:O	9:B:701:HOH:O	2.04	0.75
1:A:68:GLN:NE2	9:A:1116:HOH:O	2.19	0.74
1:A:675:LYS:HA	5:A:1007:PGE:H42	1.69	0.74
1:A:928:GLN:OE1	9:A:1104:HOH:O	2.06	0.73
1:A:646[A]:ASP:OD1	9:A:1105:HOH:O	2.07	0.71
1:A:43:GLU:OE2	9:A:1106:HOH:O	2.08	0.71
1:A:795:GLU:OE1	9:A:1107:HOH:O	2.08	0.71
1:C:900:ASP:OD1	9:C:1109:HOH:O	2.07	0.70
1:C:44:SER:OG	9:C:1110:HOH:O	2.10	0.70
1:C:506:ARG:HH12	4:C:1014:EDO:H11	1.56	0.70
1:A:753:VAL:O	9:A:1110:HOH:O	2.10	0.70
1:A:520:TYR:OH	9:A:1108:HOH:O	2.08	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:112:ARG:NH2	1:A:179:GLU:O	2.25	0.69
1:C:112:ARG:NH2	1:C:179:GLU:O	2.26	0.69
1:A:305:ASP:OD1	9:A:1109:HOH:O	2.10	0.69
1:C:897:SER:OG	9:C:1111:HOH:O	2.10	0.68
1:C:585:TYR:O	9:C:1112:HOH:O	2.11	0.68
1:A:796:GLU:OE1	9:A:1111:HOH:O	2.11	0.68
1:A:932:SER:O	9:A:1104:HOH:O	2.13	0.67
1:C:515:ARG:NH2	9:C:1108:HOH:O	2.25	0.65
1:A:960:ASP:HB3	5:A:1006:PGE:H5	1.79	0.64
1:A:576:VAL:O	9:A:1112:HOH:O	2.14	0.64
1:C:169:SER:HB2	1:C:271:SER:HB2	1.78	0.63
1:C:443:MET:O	9:C:1114:HOH:O	2.15	0.63
1:A:682:LEU:HD23	1:A:711:LEU:HD11	1.82	0.61
1:A:700:HIS:HB3	1:A:703:THR:HG23	1.83	0.60
1:A:506:ARG:NH1	9:A:1131:HOH:O	2.34	0.59
2:B:82:TYR:HB2	2:B:116:CYS:HB3	1.84	0.59
1:A:158:LEU:HB2	1:A:170:VAL:HB	1.84	0.59
1:C:846:LYS:HB2	4:C:1006:EDO:H11	1.85	0.59
1:A:106:LEU:HG	1:A:107:GLU:HG3	1.84	0.59
1:C:423:TRP:O	1:C:701:LEU:HA	2.03	0.59
1:C:700:HIS:HB3	1:C:703:THR:HG23	1.86	0.58
1:A:423:TRP:O	1:A:701:LEU:HA	2.03	0.57
1:A:607[B]:ASP:OD2	1:A:611:GLN:NE2	2.37	0.57
1:A:929:THR:O	9:A:1104:HOH:O	2.17	0.57
1:C:327:HIS:ND1	1:C:332:ASP:OD1	2.30	0.57
1:A:414:PHE:CD1	1:A:484:LYS:HG3	2.40	0.57
1:A:39:LYS:NZ	9:A:1138:HOH:O	2.32	0.57
1:A:380:GLU:OE1	9:A:1113:HOH:O	2.17	0.56
1:A:960:ASP:HB3	5:A:1006:PGE:H42	1.87	0.56
1:C:836:MET:HE2	6:C:1003:PEG:H22	1.87	0.56
1:C:50:GLN:O	1:C:376:ARG:NH2	2.40	0.55
1:C:513:LYS:NZ	9:C:1137:HOH:O	2.38	0.55
1:A:292:VAL:HG12	1:A:294:GLU:H	1.72	0.55
1:C:399:ARG:NH1	1:C:743:GLU:OE2	2.36	0.54
1:A:283:GLU:OE2	9:A:1114:HOH:O	2.18	0.53
1:C:900:ASP:OD1	1:C:902:LYS:HG2	2.08	0.53
1:C:336:PHE:HB3	1:C:387:PHE:HB2	1.90	0.52
1:A:294:GLU:OE1	9:A:1115:HOH:O	2.19	0.52
2:D:113:GLU:H	2:D:113:GLU:CD	2.17	0.52
1:A:870:ASP:OD2	1:A:873:THR:OG1	2.23	0.52
1:A:327:HIS:ND1	1:A:332:ASP:OD1	2.36	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:161:LEU:HD22	9:C:1125:HOH:O	2.09	0.52
1:A:951:ARG:HG3	2:B:55:TYR:CE1	2.45	0.52
1:A:169:SER:HB2	1:A:271:SER:HB2	1.92	0.51
1:C:436:GLN:NE2	1:C:440:ASP:OD1	2.35	0.51
1:C:460:TYR:CE2	1:C:490:ASP:HB2	2.46	0.51
1:C:458:LYS:HG2	1:C:525:TRP:HB3	1.92	0.51
1:C:607[B]:ASP:OD2	1:C:611:GLN:NE2	2.43	0.51
1:C:158:LEU:HB2	1:C:170:VAL:HB	1.93	0.51
1:A:81:GLU:O	1:A:84:LYS:NZ	2.44	0.51
1:A:675:LYS:H	5:A:1007:PGE:H6	1.75	0.51
1:A:790:TYR:CE2	1:A:792:PRO:HB3	2.45	0.51
1:A:69:LEU:HD11	1:A:128:LEU:HB2	1.93	0.50
1:A:534:PHE:HB3	1:A:600:TYR:HB3	1.92	0.50
1:A:370:THR:N	1:A:371:PRO:HD2	2.27	0.50
1:A:55:PRO:HB3	1:A:173:ARG:HB3	1.94	0.50
1:C:538:ARG:HD3	5:C:1008:PGE:H4	1.93	0.49
1:C:292:VAL:HG12	1:C:294:GLU:H	1.78	0.49
2:B:61:GLY:HA2	2:B:70:CYS:SG	2.53	0.49
1:C:370:THR:N	1:C:371:PRO:HD2	2.27	0.48
1:C:491:PRO:O	1:C:532:PRO:HD2	2.13	0.48
1:C:286:ASP:OD1	1:C:291:LYS:NZ	2.42	0.48
1:C:755:TYR:CE2	1:C:792:PRO:HG2	2.48	0.48
1:A:311:LEU:HD22	1:A:650:ILE:HD13	1.96	0.48
1:A:453:GLU:OE1	1:C:467:ARG:NE	2.46	0.48
1:A:48:LYS:NZ	9:A:1149:HOH:O	2.41	0.48
1:C:107:GLU:O	9:C:1115:HOH:O	2.20	0.48
1:A:111:PRO:HA	9:A:1202:HOH:O	2.14	0.48
1:A:447:VAL:HG11	1:A:486:VAL:HG23	1.96	0.48
1:A:69:LEU:HD13	1:A:74:LEU:HD12	1.95	0.48
1:A:658:LEU:HD22	1:A:663:LEU:HD22	1.95	0.47
1:A:900:ASP:OD1	1:A:902:LYS:HG2	2.14	0.47
1:A:542:TRP:CH2	1:A:546:MET:HE3	2.48	0.47
1:C:424:ASN:OD1	1:C:451:ASP:HB3	2.14	0.47
1:C:849:PRO:HB2	1:C:912:GLU:HB3	1.97	0.47
1:A:546:MET:HE2	4:A:1015:EDO:H21	1.96	0.47
1:C:447:VAL:HG11	1:C:486:VAL:HG23	1.97	0.47
1:A:171:ASN:HA	1:A:269:ASP:OD1	2.15	0.47
2:D:61:GLY:HA2	2:D:70:CYS:SG	2.55	0.47
1:A:175:LEU:HD13	1:A:376:ARG:HD2	1.96	0.47
1:C:58:SER:HB2	1:C:174:GLY:HA2	1.96	0.47
1:C:450:LEU:HG	1:C:485:LEU:HD21	1.95	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:104:ASP:OD1	1:A:105:GLU:N	2.47	0.46
1:A:134:ASP:OD1	1:A:137:SER:N	2.48	0.46
1:A:331:ARG:HG2	1:A:332:ASP:H	1.80	0.46
1:C:331:ARG:HG2	1:C:332:ASP:H	1.81	0.46
1:C:706:ARG:NH1	9:C:1133:HOH:O	2.35	0.46
1:C:320:SER:O	1:C:627:PHE:HA	2.16	0.46
1:A:109:ARG:NH1	1:A:181:GLN:O	2.48	0.46
1:C:847:ASP:HB3	1:C:908:PRO:HG2	1.98	0.46
1:C:259:LYS:HD2	9:C:1301:HOH:O	2.14	0.45
1:A:250:GLU:OE2	9:A:1117:HOH:O	2.20	0.45
1:A:318:TYR:CE2	1:A:639:GLY:HA3	2.51	0.45
1:A:50:GLN:O	1:A:376:ARG:NH2	2.50	0.45
1:C:278:VAL:HG23	1:C:290:LEU:HB2	1.99	0.45
1:A:675:LYS:HB3	5:A:1007:PGE:H2	1.99	0.45
1:A:849:PRO:HB2	1:A:912:GLU:HB3	1.99	0.45
1:C:311:LEU:HD22	1:C:650:ILE:HD13	1.99	0.45
1:A:459:ARG:O	1:A:462:THR:OG1	2.35	0.45
1:A:531:TYR:CZ	1:A:570:VAL:HG22	2.51	0.45
1:A:424:ASN:OD1	1:A:451:ASP:HB3	2.17	0.44
2:D:37:PHE:N	2:D:46:ILE:O	2.48	0.44
1:A:336:PHE:HB3	1:A:387:PHE:HB2	1.99	0.44
1:A:739:GLN:HG2	4:A:1002:EDO:H21	1.98	0.44
1:A:904:HIS:HB2	6:A:1020:PEG:O2	2.18	0.44
1:A:767:GLN:HG3	1:A:777:HIS:ND1	2.33	0.44
1:C:318:TYR:CE2	1:C:639:GLY:HA3	2.52	0.44
1:A:884:LEU:HG	1:A:899:ALA:HB3	2.00	0.44
1:A:671:GLY:H	1:A:685:TRP:CG	2.35	0.43
1:A:865:GLU:H	4:A:1003:EDO:H12	1.84	0.43
1:A:929:THR:HB	1:A:961:TRP:HB3	2.00	0.43
1:C:348:SER:OG	1:C:374:ASP:HB2	2.19	0.43
2:B:59:LYS:O	2:B:72:ASN:ND2	2.42	0.43
1:C:791:LEU:O	1:C:810:PRO:HA	2.19	0.43
1:C:881:GLU:HA	1:C:904:HIS:O	2.18	0.43
1:C:285:ALA:O	1:C:312:ASN:N	2.48	0.43
1:C:432:LEU:HD22	1:C:477[A]:HIS:ND1	2.34	0.43
1:A:402:ALA:O	4:A:1004:EDO:H11	2.19	0.42
1:A:450:LEU:HG	1:A:485:LEU:HD21	2.00	0.42
1:C:647:HIS:ND1	9:C:1134:HOH:O	2.36	0.42
1:A:84:LYS:NZ	9:A:1159:HOH:O	2.48	0.42
1:C:531:TYR:CZ	1:C:570:VAL:HG22	2.55	0.42
1:C:433:GLU:OE2	9:C:1116:HOH:O	2.22	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:320:SER:O	1:A:627:PHE:HA	2.20	0.42
1:C:910:TRP:CE3	1:C:954:GLY:HA2	2.55	0.42
1:A:635:ALA:HA	1:A:665:PHE:HB3	2.02	0.42
1:C:870:ASP:OD2	1:C:873:THR:OG1	2.24	0.42
1:A:93:GLY:H	1:A:128:LEU:HD21	1.85	0.42
1:A:826:GLN:OE1	1:A:917:MET:HE1	2.20	0.42
1:A:69:LEU:HD22	1:A:130:VAL:HG23	2.02	0.42
4:C:1009:EDO:O2	9:C:1117:HOH:O	2.22	0.42
1:A:304:LEU:O	9:A:1118:HOH:O	2.22	0.41
1:A:567:GLU:N	1:A:568:PRO:HA	2.35	0.41
1:C:512:VAL:HG11	1:C:578[B]:MET:SD	2.60	0.41
1:C:951:ARG:HG3	2:D:55:TYR:CE1	2.55	0.41
1:C:109:ARG:NH1	1:C:181:GLN:O	2.53	0.41
1:A:754:GLN:HB3	9:A:1184:HOH:O	2.20	0.41
1:C:146:PRO:HG2	1:C:162:GLU:HG3	2.03	0.41
1:A:129:SER:O	1:A:140:LEU:HA	2.20	0.41
1:A:635:ALA:HB2	1:A:665:PHE:CD2	2.56	0.41
1:C:152:THR:HB	1:C:157:ARG:HB3	2.03	0.41
1:C:431:VAL:HG21	1:C:450:LEU:HD21	2.01	0.41
1:A:343:THR:HA	1:A:378:MET:O	2.21	0.41
1:C:671:GLY:H	1:C:685:TRP:CG	2.39	0.41
1:C:758:ASP:OD2	1:C:790:TYR:OH	2.33	0.41
1:A:755:TYR:CE2	1:A:792:PRO:HG2	2.56	0.41
1:C:883:LEU:HG	1:C:905:LEU:HB3	2.03	0.41
2:D:70:CYS:HB2	9:D:707:HOH:O	2.20	0.41
1:C:790:TYR:CE2	1:C:792:PRO:HB3	2.56	0.40
1:C:929:THR:HG23	1:C:932:SER:HB2	2.04	0.40
1:C:826:GLN:OE1	1:C:917:MET:HE1	2.20	0.40
1:C:727:TYR:HA	1:C:730:LEU:HG	2.03	0.40
1:C:846:LYS:HB2	4:C:1006:EDO:C1	2.52	0.40
1:A:900:ASP:OD2	1:A:902:LYS:HE2	2.21	0.40
1:C:724:PHE:CE2	4:C:1011:EDO:H11	2.56	0.40

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:182:ARG:NH2	1:C:795:GLU:OE1[3_665]	2.14	0.06

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	858/977 (88%)	825 (96%)	33 (4%)	0	100	100
1	C	861/977 (88%)	831 (96%)	30 (4%)	0	100	100
2	B	81/547 (15%)	80 (99%)	1 (1%)	0	100	100
2	D	83/547 (15%)	81 (98%)	2 (2%)	0	100	100
All	All	1883/3048 (62%)	1817 (96%)	66 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	751/846 (89%)	747 (100%)	4 (0%)	81	90
1	C	752/846 (89%)	743 (99%)	9 (1%)	63	79
2	B	69/478 (14%)	68 (99%)	1 (1%)	59	76
2	D	68/478 (14%)	67 (98%)	1 (2%)	57	75
All	All	1640/2648 (62%)	1625 (99%)	15 (1%)	70	84

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	97	ASP
1	A	424	ASN

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Mol	Chain	Res	Type
1	A	446	ASP
1	A	767	GLN
2	B	117	ARG
1	C	96	LYS
1	C	97	ASP
1	C	130	VAL
1	C	424	ASN
1	C	426	ARG
1	C	446	ASP
1	C	538	ARG
1	C	767	GLN
1	C	808	HIS
2	D	85	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	77	HIS
1	A	277	HIS
1	A	372	GLN
1	A	545	ASN
1	A	805	GLN
1	A	807	HIS
1	C	77	HIS
1	C	154	GLN
1	C	408	GLN
1	C	563	ASN
1	C	808	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 41 ligands modelled in this entry, 4 are monoatomic - leaving 37 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	$\# Z > 2$
4	EDO	C	1015	-	3,3,3	0.39	0	2,2,2	0.59	0
4	EDO	A	1011	-	3,3,3	0.43	0	2,2,2	0.46	0
4	EDO	A	1018	-	3,3,3	0.39	0	2,2,2	0.51	0
5	PGE	A	1006	-	9,9,9	0.29	0	8,8,8	0.39	0
4	EDO	A	1003	-	3,3,3	0.43	0	2,2,2	0.62	0
4	EDO	A	1002	-	3,3,3	0.39	0	2,2,2	0.54	0
4	EDO	C	1013	-	3,3,3	0.46	0	2,2,2	0.41	0
4	EDO	C	1007	-	3,3,3	0.43	0	2,2,2	0.41	0
4	EDO	A	1015	-	3,3,3	0.41	0	2,2,2	0.46	0
5	PGE	C	1008	-	9,9,9	0.31	0	8,8,8	0.43	0
4	EDO	C	1006	-	3,3,3	0.42	0	2,2,2	0.45	0
4	EDO	A	1016	-	3,3,3	0.39	0	2,2,2	0.52	0
4	EDO	A	1019	-	3,3,3	0.38	0	2,2,2	0.50	0
6	PEG	A	1020	-	6,6,6	0.49	0	5,5,5	0.24	0
4	EDO	A	1017	-	3,3,3	0.41	0	2,2,2	0.47	0
4	EDO	A	1004	-	3,3,3	0.44	0	2,2,2	0.34	0
4	EDO	C	1009	-	3,3,3	0.42	0	2,2,2	0.36	0
4	EDO	A	1013	-	3,3,3	0.45	0	2,2,2	0.37	0
4	EDO	C	1011	-	3,3,3	0.38	0	2,2,2	0.53	0
4	EDO	A	1005	-	3,3,3	0.41	0	2,2,2	0.46	0
4	EDO	B	603	-	3,3,3	0.43	0	2,2,2	0.37	0
4	EDO	C	1012	-	3,3,3	0.41	0	2,2,2	0.47	0
4	EDO	C	1004	-	3,3,3	0.41	0	2,2,2	0.45	0
4	EDO	C	1010	-	3,3,3	0.44	0	2,2,2	0.48	0
4	EDO	A	1010	-	3,3,3	0.41	0	2,2,2	0.38	0
4	EDO	A	1012	-	3,3,3	0.44	0	2,2,2	0.39	0
4	EDO	C	1016	-	3,3,3	0.45	0	2,2,2	0.35	0
4	EDO	A	1008	-	3,3,3	0.44	0	2,2,2	0.40	0
4	EDO	C	1005	-	3,3,3	0.40	0	2,2,2	0.36	0
8	SO4	C	1002	-	4,4,4	0.24	0	6,6,6	0.13	0
4	EDO	C	1014	-	3,3,3	0.41	0	2,2,2	0.49	0
6	PEG	A	1009	-	6,6,6	0.49	0	5,5,5	0.30	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	PGE	A	1007	-	9,9,9	0.31	0	8,8,8	0.33	0
6	PEG	A	1014	-	6,6,6	0.50	0	5,5,5	0.34	0
3	W9P	C	1001	-	36,36,36	2.07	3 (8%)	43,50,50	1.40	5 (11%)
6	PEG	C	1003	-	6,6,6	0.49	0	5,5,5	0.36	0
3	W9P	A	1001	-	36,36,36	2.10	4 (11%)	43,50,50	1.18	4 (9%)

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	EDO	C	1015	-	-	1/1/1/1	-
4	EDO	A	1011	-	-	1/1/1/1	-
4	EDO	A	1018	-	-	1/1/1/1	-
5	PGE	A	1006	-	-	3/7/7/7	-
4	EDO	A	1003	-	-	1/1/1/1	-
4	EDO	A	1002	-	-	0/1/1/1	-
4	EDO	C	1013	-	-	0/1/1/1	-
4	EDO	C	1007	-	-	1/1/1/1	-
4	EDO	A	1015	-	-	0/1/1/1	-
5	PGE	C	1008	-	-	3/7/7/7	-
4	EDO	C	1006	-	-	1/1/1/1	-
4	EDO	A	1016	-	-	1/1/1/1	-
4	EDO	A	1019	-	-	1/1/1/1	-
6	PEG	A	1020	-	-	1/4/4/4	-
4	EDO	A	1017	-	-	1/1/1/1	-
4	EDO	A	1004	-	-	1/1/1/1	-
4	EDO	C	1009	-	-	1/1/1/1	-
4	EDO	A	1013	-	-	0/1/1/1	-
4	EDO	C	1011	-	-	1/1/1/1	-
4	EDO	A	1005	-	-	0/1/1/1	-
4	EDO	B	603	-	-	1/1/1/1	-
4	EDO	C	1012	-	-	0/1/1/1	-
4	EDO	C	1004	-	-	1/1/1/1	-
4	EDO	C	1010	-	-	0/1/1/1	-
4	EDO	A	1010	-	-	0/1/1/1	-
4	EDO	A	1012	-	-	0/1/1/1	-
4	EDO	C	1016	-	-	1/1/1/1	-
4	EDO	A	1008	-	-	1/1/1/1	-
4	EDO	C	1005	-	-	0/1/1/1	-
4	EDO	C	1014	-	-	0/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	PEG	A	1009	-	-	1/4/4/4	-
5	PGE	A	1007	-	-	5/7/7/7	-
6	PEG	A	1014	-	-	3/4/4/4	-
3	W9P	C	1001	-	-	12/20/45/45	0/3/3/3
6	PEG	C	1003	-	-	2/4/4/4	-
3	W9P	A	1001	-	-	6/20/45/45	0/3/3/3

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1001	W9P	O6-N3	10.40	1.40	1.22
3	C	1001	W9P	O6-N3	10.31	1.40	1.22
3	A	1001	W9P	C14-N2	4.36	1.48	1.37
3	C	1001	W9P	C14-N2	4.19	1.48	1.37
3	A	1001	W9P	C20-N7	2.75	1.37	1.31
3	C	1001	W9P	C20-N7	2.71	1.37	1.31
3	A	1001	W9P	C17-N4	2.01	1.47	1.43

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	1001	W9P	C8-N1-C5	-5.56	105.94	114.05
3	A	1001	W9P	C17-N4-N5	3.71	126.10	120.28
3	C	1001	W9P	C17-N4-N5	3.34	125.52	120.28
3	C	1001	W9P	N4-C20-N7	-3.04	105.97	109.60
3	A	1001	W9P	N4-C20-N7	-2.88	106.16	109.60
3	A	1001	W9P	C18-C19-C14	-2.48	119.29	121.55
3	A	1001	W9P	C15-C14-N2	-2.12	118.25	121.75
3	C	1001	W9P	C6-C5-N1	-2.07	107.46	112.03
3	C	1001	W9P	C19-C14-N2	-2.02	119.92	123.38

There are no chirality outliers.

All (52) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1001	W9P	C19-C14-N2-C13
3	C	1001	W9P	C19-C14-N2-C13
5	C	1008	PGE	C1-C2-O2-C3
5	C	1008	PGE	O3-C5-C6-O4
3	C	1001	W9P	C9-C10-C11-C12

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Mol	Chain	Res	Type	Atoms
3	C	1001	W9P	C15-C14-N2-C13
3	A	1001	W9P	C15-C14-N2-C13
5	A	1006	PGE	O2-C3-C4-O3
3	C	1001	W9P	C11-C12-C13-N2
3	A	1001	W9P	C11-C12-C13-N2
6	A	1014	PEG	O1-C1-C2-O2
3	A	1001	W9P	C9-C10-C11-C12
4	A	1004	EDO	O1-C1-C2-O2
4	C	1009	EDO	O1-C1-C2-O2
3	A	1001	W9P	C10-C11-C12-C13
5	A	1006	PGE	O1-C1-C2-O2
3	C	1001	W9P	C10-C11-C12-C13
4	A	1017	EDO	O1-C1-C2-O2
4	C	1006	EDO	O1-C1-C2-O2
5	A	1006	PGE	C4-C3-O2-C2
3	C	1001	W9P	C18-C17-N4-C20
5	A	1007	PGE	O3-C5-C6-O4
3	C	1001	W9P	C16-C17-N4-C20
4	C	1004	EDO	O1-C1-C2-O2
3	A	1001	W9P	C12-C13-N2-C14
6	C	1003	PEG	O2-C3-C4-O4
6	A	1014	PEG	O2-C3-C4-O4
6	A	1009	PEG	C4-C3-O2-C2
6	C	1003	PEG	C1-C2-O2-C3
5	A	1007	PGE	C1-C2-O2-C3
3	C	1001	W9P	C16-C17-N4-N5
3	C	1001	W9P	C18-C19-N3-O6
5	A	1007	PGE	C4-C3-O2-C2
3	C	1001	W9P	C18-C17-N4-N5
3	C	1001	W9P	C14-C19-N3-O6
4	A	1011	EDO	O1-C1-C2-O2
4	C	1011	EDO	O1-C1-C2-O2
4	C	1016	EDO	O1-C1-C2-O2
6	A	1014	PEG	C1-C2-O2-C3
3	C	1001	W9P	C12-C13-N2-C14
4	A	1018	EDO	O1-C1-C2-O2
4	B	603	EDO	O1-C1-C2-O2
5	C	1008	PGE	C6-C5-O3-C4
5	A	1007	PGE	C6-C5-O3-C4
4	A	1008	EDO	O1-C1-C2-O2
4	A	1019	EDO	O1-C1-C2-O2
4	C	1015	EDO	O1-C1-C2-O2

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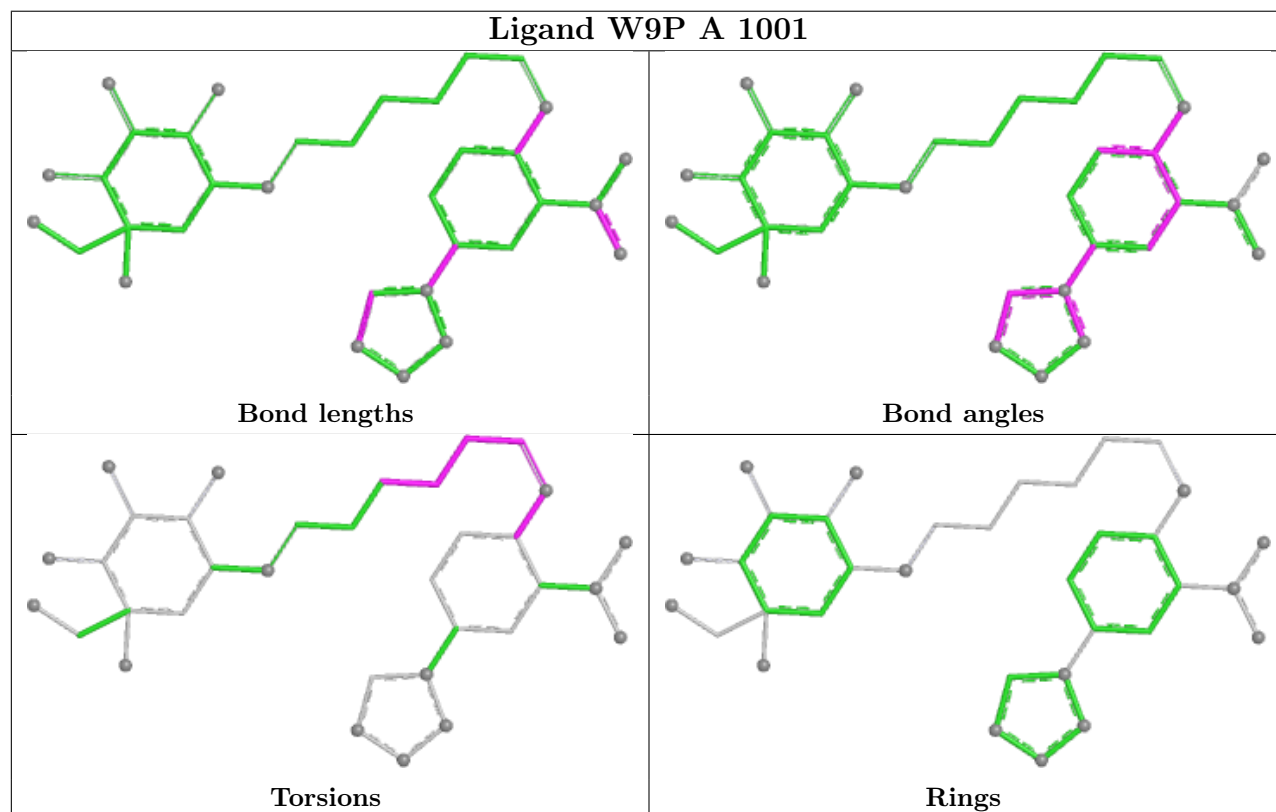
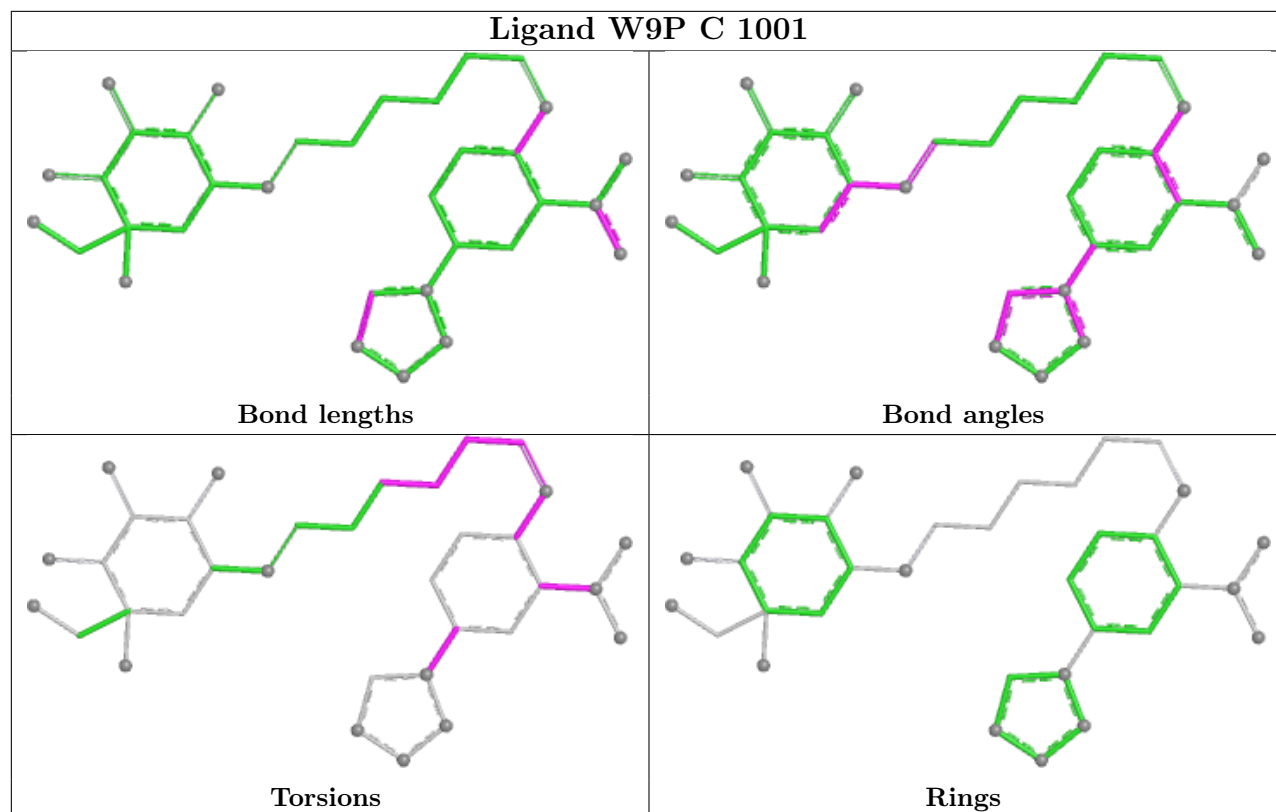
Mol	Chain	Res	Type	Atoms
4	A	1003	EDO	O1-C1-C2-O2
4	A	1016	EDO	O1-C1-C2-O2
4	C	1007	EDO	O1-C1-C2-O2
6	A	1020	PEG	O1-C1-C2-O2
5	A	1007	PGE	O2-C3-C4-O3

There are no ring outliers.

13 monomers are involved in 17 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	1006	PGE	2	0
4	A	1003	EDO	1	0
4	A	1002	EDO	1	0
4	A	1015	EDO	1	0
5	C	1008	PGE	1	0
4	C	1006	EDO	2	0
6	A	1020	PEG	1	0
4	A	1004	EDO	1	0
4	C	1009	EDO	1	0
4	C	1011	EDO	1	0
4	C	1014	EDO	1	0
5	A	1007	PGE	3	0
6	C	1003	PEG	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



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 [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	854/977 (87%)	-1.38	0 100 100	20, 42, 63, 92	10 (1%)
1	C	856/977 (87%)	-1.29	0 100 100	21, 47, 82, 100	11 (1%)
2	B	83/547 (15%)	-1.18	0 100 100	39, 55, 89, 102	0
2	D	85/547 (15%)	-1.06	0 100 100	40, 56, 93, 102	0
All	All	1878/3048 (61%)	-1.32	0 100 100	20, 45, 80, 102	21 (1%)

There are no RSRZ outliers to report.

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(Å ²)	Q<0.9
4	EDO	A	1008	4/4	0.94	0.07	58,60,63,64	0
4	EDO	A	1017	4/4	0.95	0.07	59,62,64,68	0
4	EDO	B	603	4/4	0.95	0.09	57,60,64,68	0
6	PEG	A	1014	7/7	0.95	0.08	57,63,69,69	0

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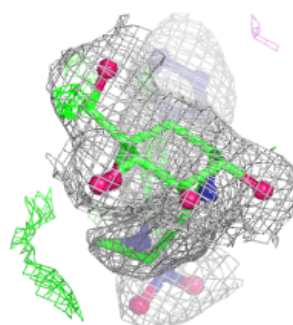
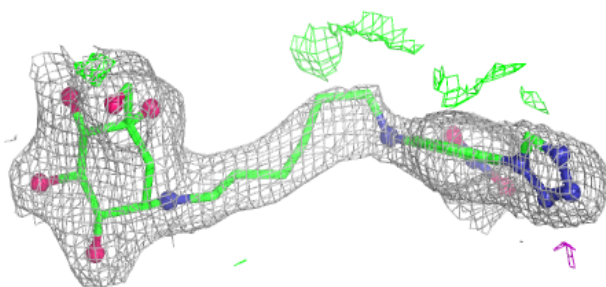
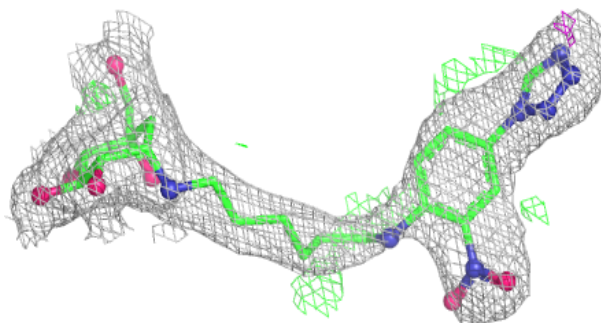
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	EDO	C	1005	4/4	0.96	0.07	51,53,55,65	0
4	EDO	C	1014	4/4	0.96	0.09	58,59,59,64	0
4	EDO	C	1004	4/4	0.96	0.10	54,58,61,67	0
6	PEG	A	1020	7/7	0.96	0.07	49,55,56,57	0
4	EDO	C	1009	4/4	0.97	0.08	54,55,56,57	0
4	EDO	A	1004	4/4	0.97	0.07	48,49,49,55	0
5	PGE	C	1008	10/10	0.97	0.07	52,56,64,70	0
4	EDO	A	1013	4/4	0.97	0.08	63,63,67,67	0
4	EDO	C	1007	4/4	0.97	0.10	47,50,52,55	0
6	PEG	C	1003	7/7	0.97	0.07	48,48,55,59	0
4	EDO	A	1005	4/4	0.98	0.05	42,48,50,54	0
4	EDO	A	1003	4/4	0.98	0.05	40,43,45,49	0
4	EDO	C	1012	4/4	0.98	0.06	41,48,49,53	0
4	EDO	C	1013	4/4	0.98	0.05	51,52,52,55	0
4	EDO	A	1018	4/4	0.98	0.08	41,42,46,47	0
4	EDO	C	1015	4/4	0.98	0.09	48,50,52,54	0
5	PGE	A	1006	10/10	0.98	0.06	52,60,68,76	0
5	PGE	A	1007	10/10	0.98	0.05	44,49,56,58	0
4	EDO	A	1010	4/4	0.98	0.06	48,50,55,58	0
6	PEG	A	1009	7/7	0.98	0.05	49,52,57,57	0
4	EDO	A	1011	4/4	0.98	0.05	54,57,59,60	0
4	EDO	A	1012	4/4	0.98	0.06	48,49,57,57	0
4	EDO	C	1006	4/4	0.98	0.08	44,46,50,52	0
8	SO4	C	1002	5/5	0.98	0.09	58,60,64,77	0
4	EDO	C	1016	4/4	0.99	0.05	51,56,58,61	0
3	W9P	A	1001	34/34	0.99	0.04	32,54,71,72	0
4	EDO	A	1019	4/4	0.99	0.06	42,44,47,53	0
4	EDO	C	1010	4/4	0.99	0.06	42,43,46,50	0
4	EDO	C	1011	4/4	0.99	0.06	47,49,53,55	0
3	W9P	C	1001	34/34	0.99	0.04	35,58,74,80	0
4	EDO	A	1015	4/4	0.99	0.06	47,49,50,61	0
4	EDO	A	1016	4/4	0.99	0.07	45,46,46,52	0
4	EDO	A	1002	4/4	0.99	0.06	40,41,42,49	0
7	CA	B	602	1/1	1.00	0.01	42,42,42,42	0
7	CA	D	601	1/1	1.00	0.01	49,49,49,49	0
7	CA	D	602	1/1	1.00	0.01	45,45,45,45	0
7	CA	B	601	1/1	1.00	0.01	51,51,51,51	0

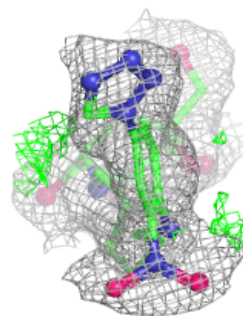
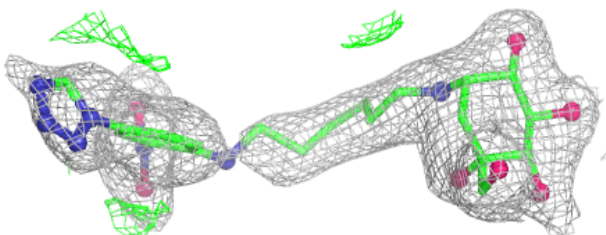
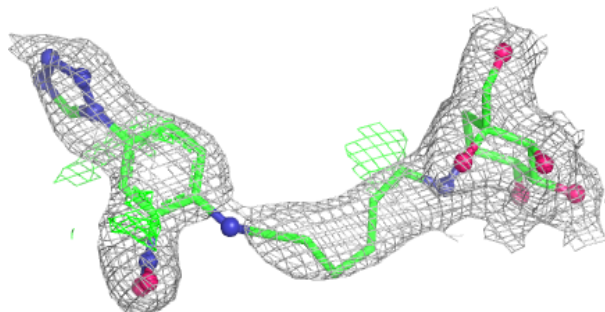
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around W9P A 1001:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around W9P C 1001:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.