



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 8, 2026 – 05:21 PM UTC

PDB ID : 3ZUK / pdb_00003zuk
Title : CRYSTAL STRUCTURE OF MYCOBACTERIUM TUBERCULOSIS ZINC METALLOPROTEASE ZMP1 IN COMPLEX WITH INHIBITOR
Authors : Ferraris, D.M.; Sbardella, D.; Petrera, A.; Marini, S.; Amstutz, B.; Coletta, M.; Sander, P.; Rizzi, M.
Deposited on : 2011-07-19
Resolution : 2.60 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
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

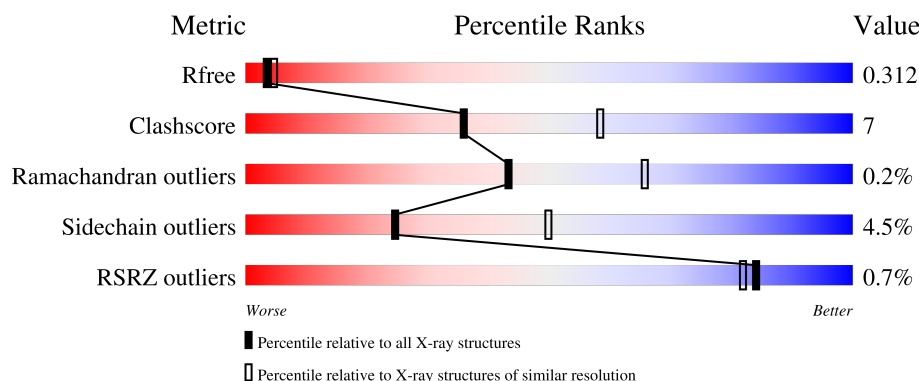
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	699	
1	B	699	

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	PEG	A	1673	-	X	-	-
5	PEG	A	1676	-	X	X	-
5	PEG	B	1666	-	X	-	-
5	PEG	B	1672	-	X	-	-
7	PG4	B	1684	-	X	X	-
8	ACT	B	1692	-	-	X	-

2 Entry composition

There are 12 unique types of molecules in this entry. The entry contains 11220 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 ENDOPEPTIDASE, PEPTIDASE FAMILY M13.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	657	Total	C	N	O	S	0	0	0
			5187	3277	925	976	9			
1	B	657	Total	C	N	O	S	0	0	0
			5187	3277	925	976	9			

There are 72 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-35	MET	-	expression tag	UNP O53649
A	-34	ARG	-	expression tag	UNP O53649
A	-33	GLY	-	expression tag	UNP O53649
A	-32	SER	-	expression tag	UNP O53649
A	-31	HIS	-	expression tag	UNP O53649
A	-30	HIS	-	expression tag	UNP O53649
A	-29	HIS	-	expression tag	UNP O53649
A	-28	HIS	-	expression tag	UNP O53649
A	-27	HIS	-	expression tag	UNP O53649
A	-26	HIS	-	expression tag	UNP O53649
A	-25	GLY	-	expression tag	UNP O53649
A	-24	MET	-	expression tag	UNP O53649
A	-23	ALA	-	expression tag	UNP O53649
A	-22	SER	-	expression tag	UNP O53649
A	-21	MET	-	expression tag	UNP O53649
A	-20	THR	-	expression tag	UNP O53649
A	-19	GLY	-	expression tag	UNP O53649
A	-18	GLY	-	expression tag	UNP O53649
A	-17	GLN	-	expression tag	UNP O53649
A	-16	GLN	-	expression tag	UNP O53649
A	-15	MET	-	expression tag	UNP O53649
A	-14	GLY	-	expression tag	UNP O53649
A	-13	ARG	-	expression tag	UNP O53649
A	-12	ASP	-	expression tag	UNP O53649
A	-11	LEU	-	expression tag	UNP O53649

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-10	TYR	-	expression tag	UNP O53649
A	-9	ASP	-	expression tag	UNP O53649
A	-8	ASP	-	expression tag	UNP O53649
A	-7	ASP	-	expression tag	UNP O53649
A	-6	ASP	-	expression tag	UNP O53649
A	-5	LYS	-	expression tag	UNP O53649
A	-4	ASP	-	expression tag	UNP O53649
A	-3	HIS	-	expression tag	UNP O53649
A	-2	PRO	-	expression tag	UNP O53649
A	-1	PHE	-	expression tag	UNP O53649
A	0	THR	-	expression tag	UNP O53649
B	-35	MET	-	expression tag	UNP O53649
B	-34	ARG	-	expression tag	UNP O53649
B	-33	GLY	-	expression tag	UNP O53649
B	-32	SER	-	expression tag	UNP O53649
B	-31	HIS	-	expression tag	UNP O53649
B	-30	HIS	-	expression tag	UNP O53649
B	-29	HIS	-	expression tag	UNP O53649
B	-28	HIS	-	expression tag	UNP O53649
B	-27	HIS	-	expression tag	UNP O53649
B	-26	HIS	-	expression tag	UNP O53649
B	-25	GLY	-	expression tag	UNP O53649
B	-24	MET	-	expression tag	UNP O53649
B	-23	ALA	-	expression tag	UNP O53649
B	-22	SER	-	expression tag	UNP O53649
B	-21	MET	-	expression tag	UNP O53649
B	-20	THR	-	expression tag	UNP O53649
B	-19	GLY	-	expression tag	UNP O53649
B	-18	GLY	-	expression tag	UNP O53649
B	-17	GLN	-	expression tag	UNP O53649
B	-16	GLN	-	expression tag	UNP O53649
B	-15	MET	-	expression tag	UNP O53649
B	-14	GLY	-	expression tag	UNP O53649
B	-13	ARG	-	expression tag	UNP O53649
B	-12	ASP	-	expression tag	UNP O53649
B	-11	LEU	-	expression tag	UNP O53649
B	-10	TYR	-	expression tag	UNP O53649
B	-9	ASP	-	expression tag	UNP O53649
B	-8	ASP	-	expression tag	UNP O53649
B	-7	ASP	-	expression tag	UNP O53649
B	-6	ASP	-	expression tag	UNP O53649
B	-5	LYS	-	expression tag	UNP O53649

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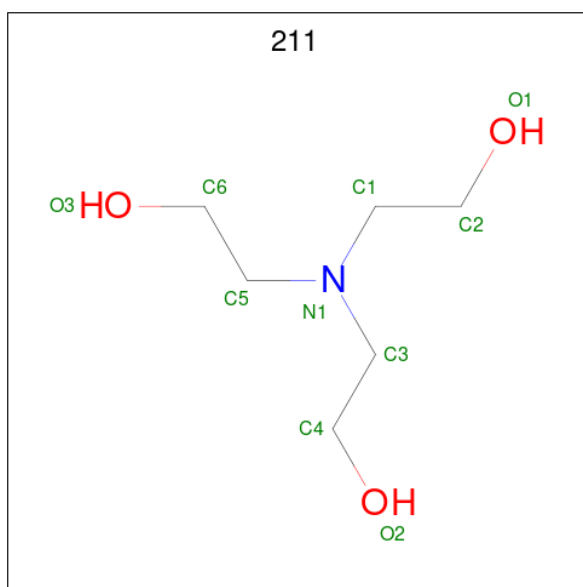
Chain	Residue	Modelled	Actual	Comment	Reference
B	-4	ASP	-	expression tag	UNP O53649
B	-3	HIS	-	expression tag	UNP O53649
B	-2	PRO	-	expression tag	UNP O53649
B	-1	PHE	-	expression tag	UNP O53649
B	0	THR	-	expression tag	UNP O53649

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

- # RDF

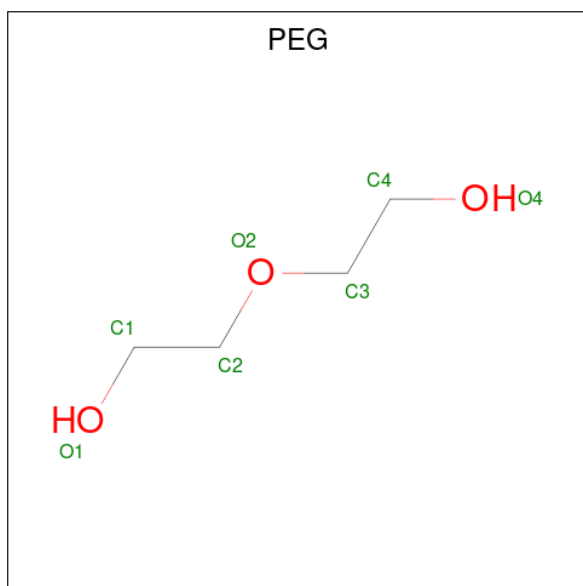
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total 37	C 23	N 3	O 10	P 1	0	0
3	B	1	Total 37	C 23	N 3	O 10	P 1	0	0

- 
- WORLD WIDE
PDB
PROTEIN DATA BANK



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

- Molecule 5 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



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

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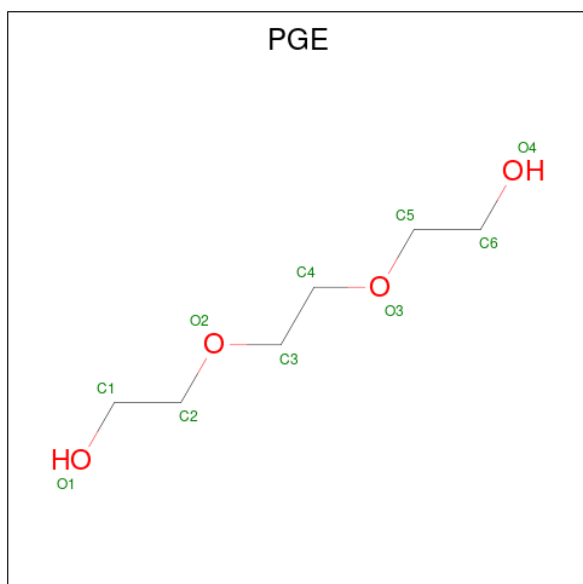
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			7	4	3		
5	A	1	Total	C	O	0	0
			7	4	3		
5	A	1	Total	C	O	0	0
			7	4	3		
5	A	1	Total	C	O	0	0
			7	4	3		
5	A	1	Total	C	O	0	0
			7	4	3		
5	A	1	Total	C	O	0	0
			7	4	3		
5	A	1	Total	C	O	0	0
			7	4	3		
5	A	1	Total	C	O	0	0
			7	4	3		
5	A	1	Total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

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	O	0	0
			7	4	3		
5	B	1	Total	C	O	0	0
			7	4	3		
5	B	1	Total	C	O	0	0
			7	4	3		
5	B	1	Total	C	O	0	0
			7	4	3		
5	B	1	Total	C	O	0	0
			7	4	3		

- Molecule 6 is TRIETHYLENE GLYCOL (CCD ID: PGE) (formula: $C_6H_{14}O_4$).



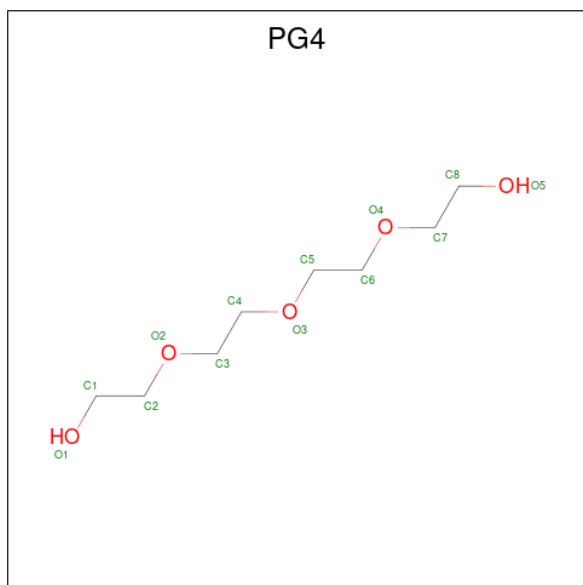
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			10	6	4		
6	A	1	Total	C	O	0	0
			10	6	4		
6	A	1	Total	C	O	0	0
			10	6	4		
6	A	1	Total	C	O	0	0
			10	6	4		
6	A	1	Total	C	O	0	0
			10	6	4		

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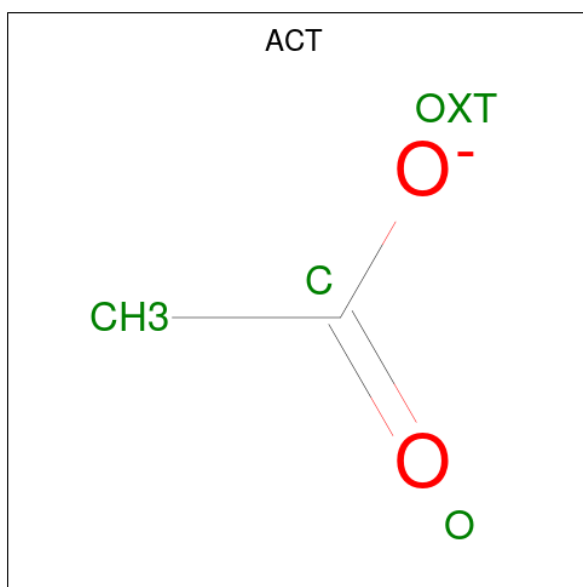
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	B	1	Total	C	O	0	0
			10	6	4		
6	B	1	Total	C	O	0	0
			10	6	4		
6	B	1	Total	C	O	0	0
			10	6	4		

- Molecule 7 is TETRAETHYLENE GLYCOL (CCD ID: PG4) (formula: $C_8H_{18}O_5$).



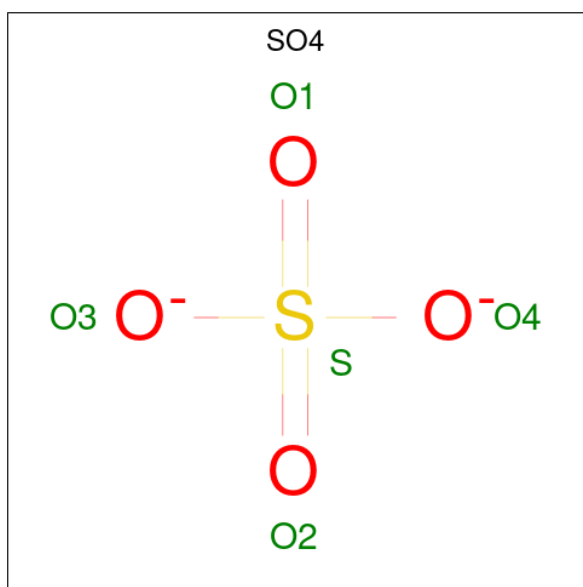
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			13	8	5		
7	B	1	Total	C	O	0	0
			13	8	5		
7	B	1	Total	C	O	0	0
			13	8	5		
7	B	1	Total	C	O	0	0
			13	8	5		

- Molecule 8 is ACETATE ION (CCD ID: ACT) (formula: $C_2H_3O_2$).



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

- Molecule 9 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

- Molecule 10 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

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

- Molecule 11 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	B	1	Total	Ca	0	0
			1	1		

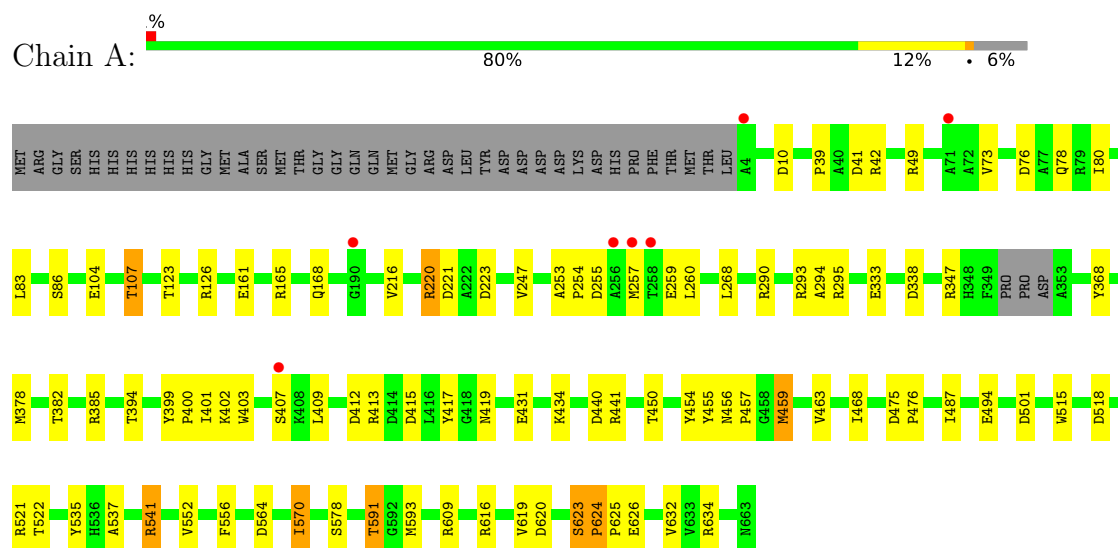
- Molecule 12 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
12	A	159	Total 159	O 159	0	0
12	B	173	Total 173	O 173	0	0

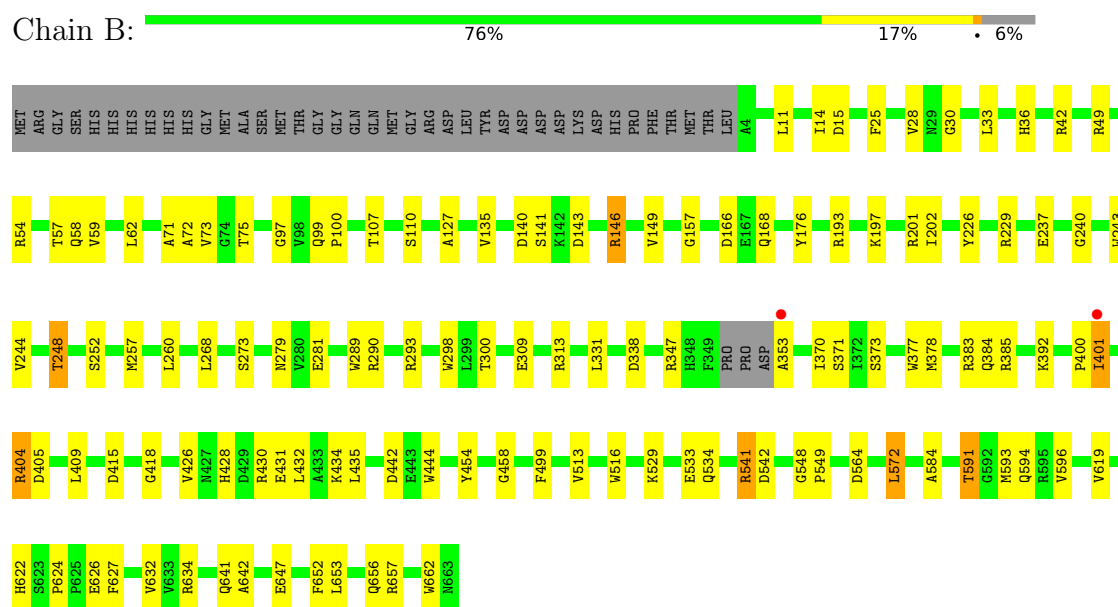
3 Residue-property plots [i](#)

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: ENDOPEPTIDASE, PEPTIDASE FAMILY M13



• Molecule 1: ENDOPEPTIDASE, PEPTIDASE FAMILY M13



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	128.29Å 198.77Å 63.27Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.60 30.00 – 2.60	Depositor EDS
% Data completeness (in resolution range)	92.8 (30.00-2.60) 92.7 (30.00-2.60)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.32 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.175 , 0.250 0.185 , 0.312	Depositor DCC
R_{free} test set	1458 reflections (2.00%)	wwPDB-VP
Wilson B-factor (Å ²)	27.0	Xtriage
Anisotropy	0.101	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 42.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.37$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	11220	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.45% 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, 211, ACT, CA, PEG, SO4, PG4, PGE, CL, RDF

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	0.88	1/5317 (0.0%)	1.04	8/7236 (0.1%)
1	B	0.89	0/5317	1.03	3/7236 (0.0%)
All	All	0.88	1/10634 (0.0%)	1.04	11/14472 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	468	ILE	CA-CB	5.85	1.61	1.54

The worst 5 of 11 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	399	TYR	N-CA-C	6.06	113.67	108.22
1	B	454	TYR	N-CA-C	5.88	117.19	108.60
1	B	240	GLY	N-CA-C	-5.85	107.32	113.58
1	A	624	PRO	CA-C-N	5.48	125.57	119.32
1	A	624	PRO	C-N-CA	5.48	125.57	119.32

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	5187	0	4968	60	0
1	B	5187	0	4968	81	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	37	0	32	8	0
3	B	37	0	32	5	0
4	A	10	0	15	1	0
5	A	105	0	150	14	0
5	B	105	0	150	10	0
6	A	50	0	70	1	0
6	B	30	0	42	2	0
7	A	13	0	18	0	0
7	B	39	0	54	12	0
8	A	20	0	15	0	0
8	B	32	0	24	2	0
9	A	10	0	0	0	0
9	B	20	0	0	0	0
10	A	1	0	0	0	0
10	B	2	0	0	0	0
11	B	1	0	0	0	0
12	A	159	0	0	8	0
12	B	173	0	0	9	0
All	All	11220	0	10538	156	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 7.

The worst 5 of 156 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)
3:B:1665:RDF:HD1	3:B:1665:RDF:CD11	0.97	1.15
3:A:1665:RDF:CD11	3:A:1665:RDF:HD1	0.97	1.13
1:B:541:ARG:HH11	1:B:541:ARG:HG2	1.16	1.05
1:B:591:THR:HG22	1:B:594:GLN:H	1.29	0.95
1:B:548:GLY:H	5:B:1670:PEG:H31	1.31	0.92

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	653/699 (93%)	624 (96%)	27 (4%)	2 (0%)	36	58
1	B	653/699 (93%)	622 (95%)	31 (5%)	0	100	100
All	All	1306/1398 (93%)	1246 (95%)	58 (4%)	2 (0%)	43	66

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	168	GLN
1	A	254	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	522/558 (94%)	500 (96%)	22 (4%)	26	52
1	B	522/558 (94%)	497 (95%)	25 (5%)	23	47
All	All	1044/1116 (94%)	997 (96%)	47 (4%)	24	50

5 of 47 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	237	GLU
1	B	383	ARG
1	B	248	THR
1	B	338	ASP

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Mol	Chain	Res	Type
1	B	385	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 10 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	456	ASN
1	B	546	HIS
1	B	576	GLN
1	A	546	HIS
1	B	58	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 70 ligands modelled in this entry, 6 are monoatomic - leaving 64 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	PEG	B	1667	-	6,6,6	0.65	0	5,5,5	1.45	0
7	PG4	B	1685	-	12,12,12	1.09	0	11,11,11	1.70	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	PGE	B	1682	-	9,9,9	0.61	0	8,8,8	1.66	1 (12%)
5	PEG	A	1676	-	6,6,6	0.48	0	5,5,5	2.23	3 (60%)
5	PEG	B	1680	-	6,6,6	0.55	0	5,5,5	1.57	2 (40%)
5	PEG	A	1671	-	6,6,6	0.55	0	5,5,5	1.57	1 (20%)
5	PEG	B	1672	-	6,6,6	0.53	0	5,5,5	1.77	2 (40%)
6	PGE	A	1684	-	9,9,9	0.64	0	8,8,8	1.60	1 (12%)
5	PEG	A	1678	-	6,6,6	0.47	0	5,5,5	1.95	2 (40%)
5	PEG	A	1677	-	6,6,6	0.61	0	5,5,5	1.45	0
6	PGE	A	1685	-	9,9,9	0.51	0	8,8,8	1.94	3 (37%)
5	PEG	B	1669	-	6,6,6	0.57	0	5,5,5	1.41	0
8	ACT	A	1690	-	3,3,3	0.85	0	3,3,3	0.81	0
5	PEG	A	1672	-	6,6,6	0.67	0	5,5,5	1.43	0
6	PGE	A	1683	-	9,9,9	0.56	0	8,8,8	1.91	3 (37%)
9	SO4	B	1695	-	4,4,4	0.39	0	6,6,6	0.51	0
5	PEG	B	1676	-	6,6,6	0.56	0	5,5,5	1.94	2 (40%)
5	PEG	B	1671	-	6,6,6	0.56	0	5,5,5	1.60	0
5	PEG	A	1670	-	6,6,6	0.49	0	5,5,5	1.70	1 (20%)
6	PGE	A	1682	-	9,9,9	0.53	0	8,8,8	1.89	4 (50%)
5	PEG	B	1677	-	6,6,6	0.48	0	5,5,5	1.79	0
5	PEG	B	1678	-	6,6,6	0.62	0	5,5,5	1.40	0
5	PEG	A	1669	-	6,6,6	0.56	0	5,5,5	1.68	0
7	PG4	A	1687	-	12,12,12	1.13	1 (8%)	11,11,11	1.82	3 (27%)
5	PEG	B	1666	-	6,6,6	0.52	0	5,5,5	1.67	2 (40%)
8	ACT	B	1690	-	3,3,3	0.82	0	3,3,3	0.82	0
5	PEG	B	1679	-	6,6,6	0.48	0	5,5,5	1.89	2 (40%)
5	PEG	A	1681	-	6,6,6	0.53	0	5,5,5	1.78	1 (20%)
8	ACT	B	1694	-	3,3,3	0.76	0	3,3,3	0.96	0
5	PEG	A	1675	-	6,6,6	0.52	0	5,5,5	1.64	1 (20%)
5	PEG	A	1679	-	6,6,6	0.60	0	5,5,5	1.58	1 (20%)
7	PG4	B	1684	-	12,12,12	1.33	3 (25%)	11,11,11	2.62	10 (90%)
8	ACT	A	1688	-	3,3,3	0.94	0	3,3,3	0.55	0
6	PGE	B	1683	-	9,9,9	0.56	0	8,8,8	1.61	2 (25%)
4	211	A	1666	-	9,9,9	0.83	0	9,9,9	1.32	0
9	SO4	B	1697	-	4,4,4	0.35	0	6,6,6	0.32	0
9	SO4	B	1698	-	4,4,4	0.46	0	6,6,6	0.47	0
5	PEG	B	1670	-	6,6,6	0.54	0	5,5,5	1.55	1 (20%)
5	PEG	B	1674	-	6,6,6	0.51	0	5,5,5	1.73	2 (40%)
8	ACT	A	1692	-	3,3,3	0.81	0	3,3,3	0.90	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	PEG	A	1668	-	6,6,6	0.63	0	5,5,5	1.42	0
6	PGE	A	1686	-	9,9,9	0.66	0	8,8,8	1.37	0
8	ACT	B	1687	-	3,3,3	0.71	0	3,3,3	1.02	0
8	ACT	B	1693	-	3,3,3	0.75	0	3,3,3	1.16	0
5	PEG	B	1668	-	6,6,6	0.60	0	5,5,5	1.66	1 (20%)
8	ACT	A	1689	-	3,3,3	0.87	0	3,3,3	0.96	0
5	PEG	A	1673	-	6,6,6	0.42	0	5,5,5	2.26	3 (60%)
5	PEG	B	1675	-	6,6,6	0.59	0	5,5,5	1.70	1 (20%)
6	PGE	B	1681	-	9,9,9	0.53	0	8,8,8	1.82	3 (37%)
5	PEG	A	1674	-	6,6,6	0.56	0	5,5,5	1.53	1 (20%)
5	PEG	B	1673	-	6,6,6	0.57	0	5,5,5	1.65	1 (20%)
3	RDF	A	1665	2	38,39,39	2.20	11 (28%)	53,57,57	2.54	15 (28%)
8	ACT	A	1691	-	3,3,3	0.79	0	3,3,3	0.99	0
5	PEG	A	1667	-	6,6,6	0.57	0	5,5,5	1.57	1 (20%)
8	ACT	B	1688	-	3,3,3	0.87	0	3,3,3	0.90	0
8	ACT	B	1689	-	3,3,3	0.97	0	3,3,3	0.80	0
9	SO4	A	1694	-	4,4,4	0.46	0	6,6,6	0.44	0
9	SO4	A	1693	-	4,4,4	0.47	0	6,6,6	0.26	0
5	PEG	A	1680	-	6,6,6	0.60	0	5,5,5	1.45	0
3	RDF	B	1665	2	38,39,39	2.25	11 (28%)	53,57,57	2.52	15 (28%)
8	ACT	B	1692	-	3,3,3	0.84	0	3,3,3	0.64	0
9	SO4	B	1696	-	4,4,4	0.44	0	6,6,6	0.48	0
8	ACT	B	1691	-	3,3,3	0.87	0	3,3,3	0.98	0
7	PG4	B	1686	-	12,12,12	1.11	1 (8%)	11,11,11	1.71	4 (36%)

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	PEG	B	1667	-	-	2/4/4/4	-
7	PG4	B	1685	-	-	8/10/10/10	-
6	PGE	B	1682	-	-	1/7/7/7	-
5	PEG	A	1676	-	-	3/4/4/4	-
5	PEG	B	1680	-	-	2/4/4/4	-
5	PEG	A	1671	-	-	4/4/4/4	-
5	PEG	B	1672	-	-	4/4/4/4	-
6	PGE	A	1684	-	-	5/7/7/7	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	PEG	A	1678	-	-	2/4/4/4	-
5	PEG	A	1677	-	-	2/4/4/4	-
6	PGE	A	1685	-	-	4/7/7/7	-
5	PEG	B	1669	-	-	3/4/4/4	-
5	PEG	A	1672	-	-	2/4/4/4	-
6	PGE	A	1683	-	-	4/7/7/7	-
5	PEG	B	1676	-	-	2/4/4/4	-
5	PEG	B	1671	-	-	3/4/4/4	-
5	PEG	A	1670	-	-	4/4/4/4	-
6	PGE	A	1682	-	-	2/7/7/7	-
5	PEG	B	1677	-	-	3/4/4/4	-
5	PEG	B	1678	-	-	3/4/4/4	-
5	PEG	A	1669	-	-	4/4/4/4	-
7	PG4	A	1687	-	-	6/10/10/10	-
5	PEG	B	1666	-	-	4/4/4/4	-
5	PEG	B	1679	-	-	1/4/4/4	-
5	PEG	A	1681	-	-	1/4/4/4	-
5	PEG	A	1675	-	-	2/4/4/4	-
5	PEG	A	1679	-	-	4/4/4/4	-
7	PG4	B	1684	-	-	8/10/10/10	-
6	PGE	B	1683	-	-	3/7/7/7	-
4	211	A	1666	-	-	4/9/9/9	-
5	PEG	B	1670	-	-	3/4/4/4	-
5	PEG	B	1674	-	-	2/4/4/4	-
5	PEG	A	1668	-	-	2/4/4/4	-
6	PGE	A	1686	-	-	3/7/7/7	-
5	PEG	B	1668	-	-	4/4/4/4	-
5	PEG	A	1673	-	-	4/4/4/4	-
5	PEG	B	1675	-	-	3/4/4/4	-
6	PGE	B	1681	-	-	4/7/7/7	-
5	PEG	A	1674	-	-	3/4/4/4	-
5	PEG	B	1673	-	-	4/4/4/4	-
3	RDF	A	1665	2	-	3/27/50/50	0/3/3/3
5	PEG	A	1667	-	-	2/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	PEG	A	1680	-	-	2/4/4/4	-
3	RDF	B	1665	2	-	5/27/50/50	0/3/3/3
7	PG4	B	1686	-	-	7/10/10/10	-

The worst 5 of 27 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1665	RDF	C-N1	6.10	1.47	1.34
3	A	1665	RDF	P-O1P	5.66	1.54	1.46
3	A	1665	RDF	C-N1	5.56	1.46	1.34
3	B	1665	RDF	P-O1P	5.50	1.54	1.46
3	B	1665	RDF	CD21-CG1	-4.88	1.36	1.44

The worst 5 of 95 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1665	RDF	O1-P-O1P	-10.21	97.48	115.32
3	A	1665	RDF	O1-P-O1P	-9.82	98.16	115.32
3	A	1665	RDF	O1P-P-N	-8.81	98.81	112.75
3	B	1665	RDF	O1P-P-N	-7.16	101.42	112.75
3	B	1665	RDF	CD21-CG1-CD11	5.41	111.69	106.16

There are no chirality outliers.

5 of 151 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1665	RDF	C7-CA1-CB1-CG1
5	A	1681	PEG	C1-C2-O2-C3
7	B	1686	PG4	C5-C6-O4-C7
5	A	1675	PEG	C4-C3-O2-C2
5	A	1676	PEG	C1-C2-O2-C3

There are no ring outliers.

25 monomers are involved in 53 short contacts:

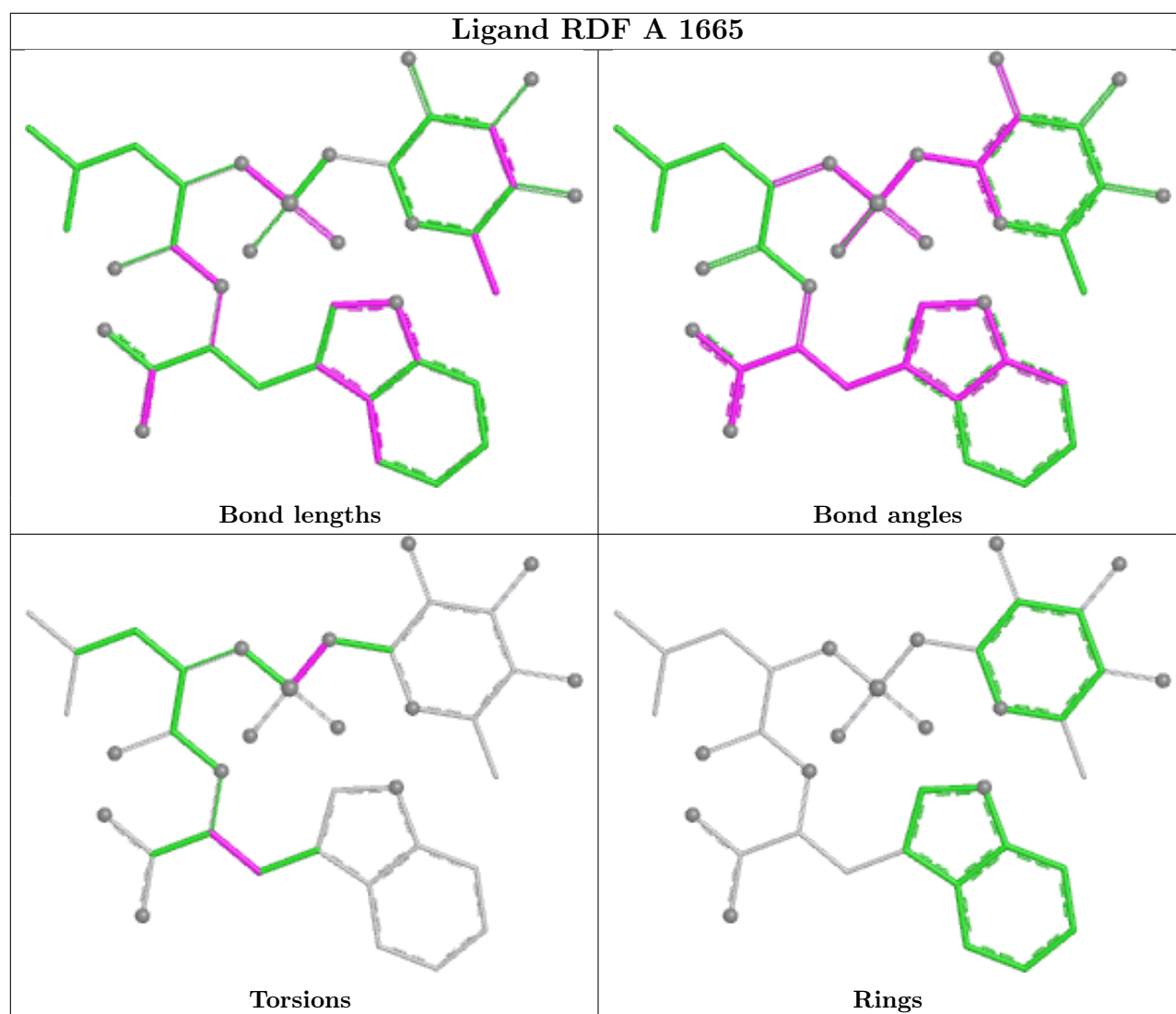
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	1667	PEG	1	0
7	B	1685	PG4	1	0
6	B	1682	PGE	1	0
5	A	1676	PEG	6	0

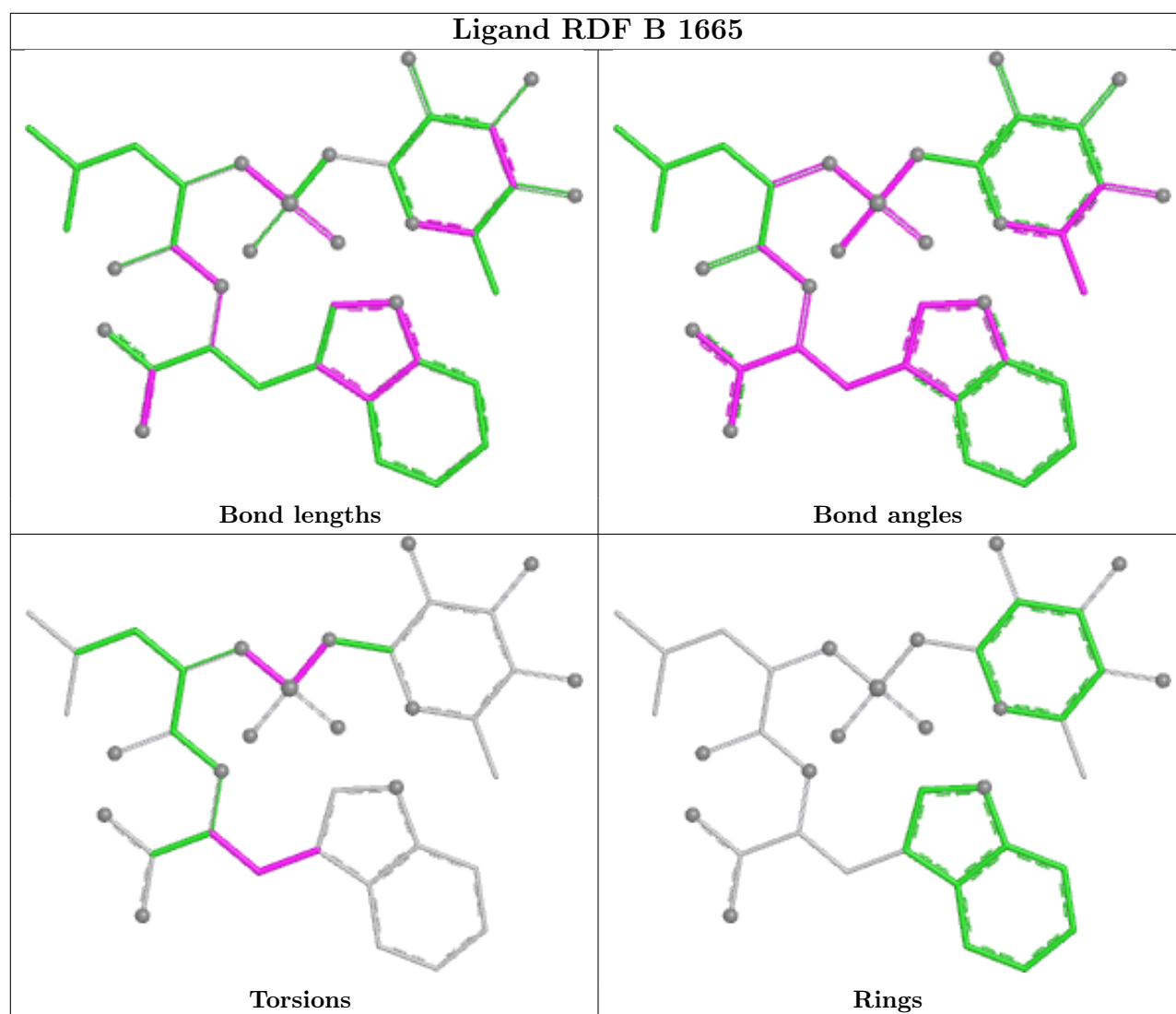
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	1680	PEG	1	0
6	A	1684	PGE	1	0
5	B	1676	PEG	2	0
5	B	1671	PEG	1	0
5	B	1678	PEG	1	0
5	A	1669	PEG	3	0
5	B	1679	PEG	1	0
5	A	1681	PEG	1	0
7	B	1684	PG4	8	0
6	B	1683	PGE	1	0
4	A	1666	211	1	0
5	B	1670	PEG	1	0
5	A	1668	PEG	2	0
5	B	1668	PEG	1	0
5	A	1673	PEG	1	0
5	B	1673	PEG	1	0
3	A	1665	RDF	8	0
5	A	1667	PEG	1	0
3	B	1665	RDF	5	0
8	B	1692	ACT	2	0
7	B	1686	PG4	3	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 [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	657/699 (93%)	-0.31	7 (1%)	78 74	11, 29, 53, 66	0
1	B	657/699 (93%)	-0.40	2 (0%)	90 88	10, 26, 45, 64	0
All	All	1314/1398 (93%)	-0.35	9 (0%)	84 82	10, 27, 51, 66	0

The worst 5 of 9 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	256	ALA	3.3
1	B	353	ALA	2.9
1	A	71	ALA	2.4
1	A	407	SER	2.3
1	A	190	GLY	2.3

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
5	PEG	B	1680	7/7	0.40	0.32	61,65,68,69	0
8	ACT	B	1689	4/4	0.56	0.23	48,48,48,48	0
8	ACT	B	1693	4/4	0.57	0.21	52,52,53,53	0
5	PEG	A	1675	7/7	0.60	0.28	61,63,66,67	0
5	PEG	A	1671	7/7	0.63	0.28	72,76,80,80	0
8	ACT	A	1691	4/4	0.63	0.27	63,63,64,64	0
8	ACT	B	1688	4/4	0.66	0.22	51,51,51,52	0
8	ACT	A	1688	4/4	0.66	0.17	34,35,36,36	0
8	ACT	B	1687	4/4	0.66	0.21	71,72,72,72	0
5	PEG	B	1679	7/7	0.67	0.23	55,59,63,63	0
8	ACT	A	1689	4/4	0.69	0.23	48,49,49,49	0
5	PEG	A	1677	7/7	0.69	0.27	68,71,72,72	0
8	ACT	B	1692	4/4	0.70	0.20	48,48,48,48	0
5	PEG	A	1674	7/7	0.71	0.18	50,53,55,55	0
5	PEG	A	1673	7/7	0.73	0.18	47,52,54,55	0
5	PEG	A	1678	7/7	0.73	0.23	59,65,67,67	0
6	PGE	B	1681	10/10	0.73	0.18	44,49,52,52	0
7	PG4	B	1685	13/13	0.73	0.24	65,78,87,88	0
8	ACT	B	1691	4/4	0.73	0.23	53,54,54,54	0
5	PEG	B	1672	7/7	0.73	0.24	48,54,55,56	0
5	PEG	B	1677	7/7	0.73	0.28	74,74,75,75	0
5	PEG	A	1681	7/7	0.74	0.24	69,70,70,70	0
5	PEG	B	1669	7/7	0.75	0.17	48,51,54,55	0
5	PEG	A	1670	7/7	0.75	0.21	48,54,57,57	0
6	PGE	B	1683	10/10	0.77	0.16	45,49,49,50	0
7	PG4	B	1684	13/13	0.77	0.20	36,43,47,48	0
5	PEG	A	1669	7/7	0.77	0.17	49,52,54,54	0
5	PEG	B	1674	7/7	0.77	0.18	58,59,61,61	0
6	PGE	A	1685	10/10	0.77	0.22	63,65,70,70	0
5	PEG	B	1670	7/7	0.77	0.15	45,46,48,48	0
5	PEG	B	1666	7/7	0.78	0.18	49,53,54,54	0
5	PEG	B	1678	7/7	0.78	0.18	50,54,56,57	0
8	ACT	B	1690	4/4	0.79	0.17	51,52,52,52	0
8	ACT	A	1692	4/4	0.79	0.20	51,52,52,52	0
5	PEG	A	1668	7/7	0.80	0.16	46,48,51,51	0
5	PEG	B	1667	7/7	0.80	0.21	56,58,61,62	0
5	PEG	B	1676	7/7	0.80	0.21	49,53,54,55	0
6	PGE	A	1683	10/10	0.80	0.17	41,51,57,57	0
5	PEG	B	1671	7/7	0.80	0.20	51,54,57,58	0
6	PGE	A	1686	10/10	0.81	0.20	52,53,57,57	0
6	PGE	A	1682	10/10	0.81	0.16	50,56,58,59	0
7	PG4	B	1686	13/13	0.82	0.15	43,53,58,59	0
5	PEG	B	1668	7/7	0.82	0.19	47,48,51,52	0

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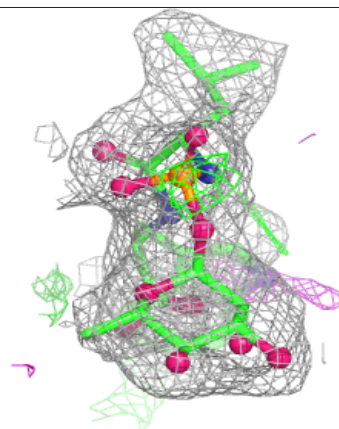
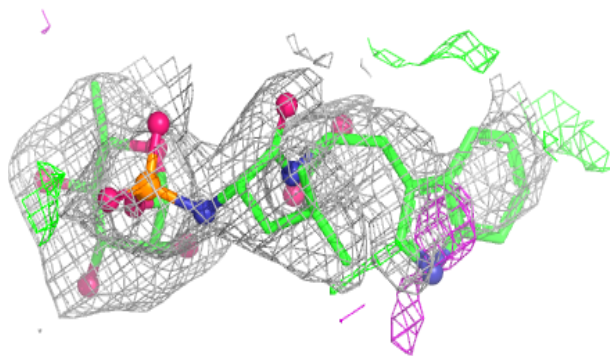
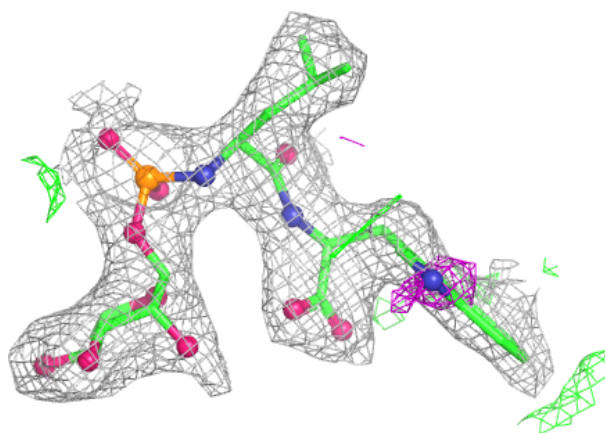
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	211	A	1666	10/10	0.82	0.17	49,55,57,57	0
6	PGE	B	1682	10/10	0.82	0.14	51,55,57,57	0
5	PEG	A	1667	7/7	0.83	0.17	55,56,57,58	0
7	PG4	A	1687	13/13	0.83	0.16	54,58,59,59	0
9	SO4	B	1695	5/5	0.83	0.20	91,91,92,92	0
5	PEG	B	1675	7/7	0.84	0.16	53,53,56,56	0
8	ACT	A	1690	4/4	0.85	0.17	53,53,54,54	0
8	ACT	B	1694	4/4	0.86	0.19	49,50,50,50	0
5	PEG	A	1680	7/7	0.86	0.16	59,59,60,61	0
5	PEG	B	1673	7/7	0.87	0.13	49,50,53,55	0
5	PEG	A	1679	7/7	0.88	0.12	41,42,46,48	0
5	PEG	A	1676	7/7	0.89	0.23	46,48,50,51	0
9	SO4	B	1697	5/5	0.89	0.11	68,69,70,70	0
5	PEG	A	1672	7/7	0.90	0.13	57,60,63,63	0
6	PGE	A	1684	10/10	0.91	0.12	32,40,42,42	0
10	CL	A	1695	1/1	0.91	0.12	59,59,59,59	0
10	CL	B	1701	1/1	0.91	0.10	57,57,57,57	0
10	CL	B	1699	1/1	0.92	0.09	56,56,56,56	0
9	SO4	B	1698	5/5	0.93	0.17	52,53,55,55	0
9	SO4	A	1694	5/5	0.93	0.14	51,52,54,54	0
3	RDF	B	1665	37/37	0.95	0.09	16,19,53,53	0
3	RDF	A	1665	37/37	0.95	0.09	15,19,52,52	0
9	SO4	B	1696	5/5	0.98	0.05	26,26,28,31	0
9	SO4	A	1693	5/5	0.98	0.05	24,24,25,26	0
2	ZN	B	1664	1/1	1.00	0.02	16,16,16,16	0
2	ZN	A	1664	1/1	1.00	0.01	18,18,18,18	0
11	CA	B	1700	1/1	1.00	0.03	19,19,19,19	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.

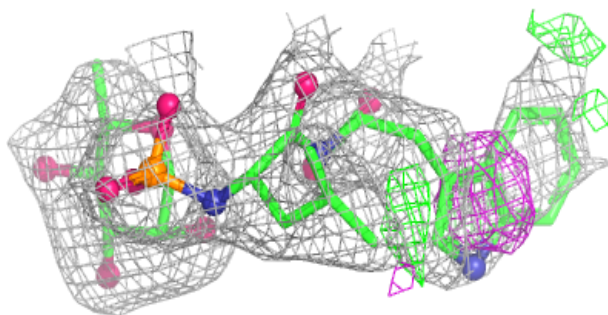
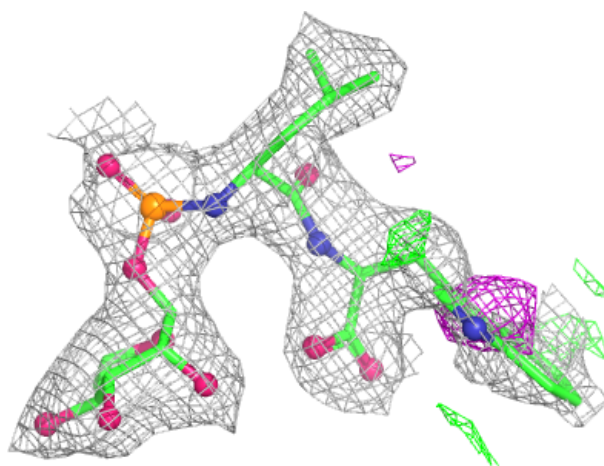
Electron density around RDF B 1665:

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



Electron density around RDF A 1665:

$2mF_o-DF_c$ (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.