



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 5, 2026 – 08:48 PM UTC

PDB ID : 3MDJ / pdb_00003mdj
Title : ER Aminopeptidase, ERAP1, Bound to the Zinc Aminopeptidase Inhibitor, Bestatin
Authors : Nguyen, T.T.; Stern, L.J.
Deposited on : 2010-03-30
Resolution : 2.95 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)
Xtrriage (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

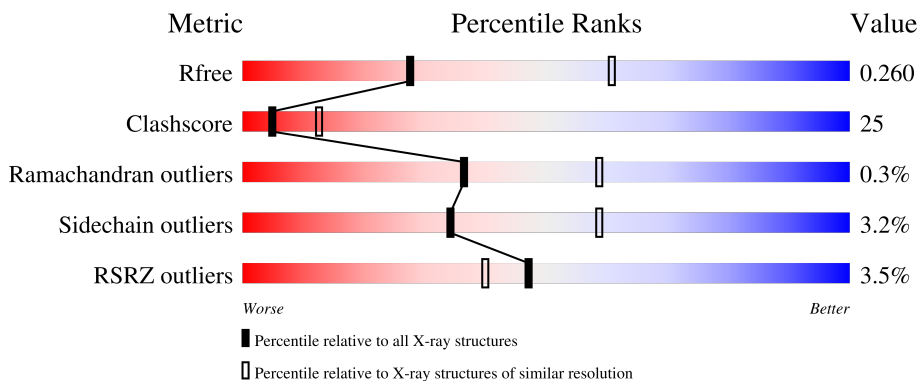
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.95 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	1130 (2.98-2.94)
Clashscore	190562	1157 (2.98-2.94)
Ramachandran outliers	187476	1101 (2.98-2.94)
Sidechain outliers	187428	1101 (2.98-2.94)
RSRZ outliers	180081	1130 (2.98-2.94)

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	921	
1	B	921	
1	C	921	
2	D	5	
2	F	5	

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Mol	Chain	Length	Quality of chain
3	E	2	

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
2	NAG	F	1	X	-	-	-
5	BES	B	1001	-	-	X	-
6	NAG	A	6001	X	-	-	-

2 Entry composition i

There are 6 unique types of molecules in this entry. The entry contains 20163 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 Endoplasmic reticulum aminopeptidase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	818	Total 6625	C 4271	N 1096	O 1227	S 31	55	0	0
1	B	821	Total 6643	C 4281	N 1099	O 1232	S 31	50	0	0
1	C	819	Total 6634	C 4276	N 1097	O 1230	S 31	68	0	0

There are 66 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	346	ASP	GLY	SEE REMARK 999	UNP Q9NZ08
A	514	ARG	GLY	SEE REMARK 999	UNP Q9NZ08
A	528	ARG	LYS	SEE REMARK 999	UNP Q9NZ08
A	730	GLU	GLN	SEE REMARK 999	UNP Q9NZ08
A	940	HIS	-	expression tag	UNP Q9NZ08
A	941	ASP	-	expression tag	UNP Q9NZ08
A	942	PRO	-	expression tag	UNP Q9NZ08
A	943	GLU	-	expression tag	UNP Q9NZ08
A	944	ALA	-	expression tag	UNP Q9NZ08
A	945	ASP	-	expression tag	UNP Q9NZ08
A	946	ALA	-	expression tag	UNP Q9NZ08
A	947	THR	-	expression tag	UNP Q9NZ08
A	948	GLY	-	expression tag	UNP Q9NZ08
A	949	LEU	-	expression tag	UNP Q9NZ08
A	950	GLU	-	expression tag	UNP Q9NZ08
A	951	ARG	-	expression tag	UNP Q9NZ08
A	952	MET	-	expression tag	UNP Q9NZ08
A	953	LEU	-	expression tag	UNP Q9NZ08
A	954	GLU	-	expression tag	UNP Q9NZ08
A	955	SER	-	expression tag	UNP Q9NZ08
A	956	ARG	-	expression tag	UNP Q9NZ08
A	957	GLY	-	expression tag	UNP Q9NZ08
B	346	ASP	GLY	SEE REMARK 999	UNP Q9NZ08

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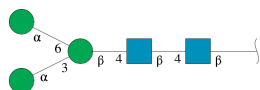
Chain	Residue	Modelled	Actual	Comment	Reference
B	514	ARG	GLY	SEE REMARK 999	UNP Q9NZ08
B	528	ARG	LYS	SEE REMARK 999	UNP Q9NZ08
B	730	GLU	GLN	SEE REMARK 999	UNP Q9NZ08
B	940	HIS	-	expression tag	UNP Q9NZ08
B	941	ASP	-	expression tag	UNP Q9NZ08
B	942	PRO	-	expression tag	UNP Q9NZ08
B	943	GLU	-	expression tag	UNP Q9NZ08
B	944	ALA	-	expression tag	UNP Q9NZ08
B	945	ASP	-	expression tag	UNP Q9NZ08
B	946	ALA	-	expression tag	UNP Q9NZ08
B	947	THR	-	expression tag	UNP Q9NZ08
B	948	GLY	-	expression tag	UNP Q9NZ08
B	949	LEU	-	expression tag	UNP Q9NZ08
B	950	GLU	-	expression tag	UNP Q9NZ08
B	951	ARG	-	expression tag	UNP Q9NZ08
B	952	MET	-	expression tag	UNP Q9NZ08
B	953	LEU	-	expression tag	UNP Q9NZ08
B	954	GLU	-	expression tag	UNP Q9NZ08
B	955	SER	-	expression tag	UNP Q9NZ08
B	956	ARG	-	expression tag	UNP Q9NZ08
B	957	GLY	-	expression tag	UNP Q9NZ08
C	346	ASP	GLY	SEE REMARK 999	UNP Q9NZ08
C	514	ARG	GLY	SEE REMARK 999	UNP Q9NZ08
C	528	ARG	LYS	SEE REMARK 999	UNP Q9NZ08
C	730	GLU	GLN	SEE REMARK 999	UNP Q9NZ08
C	940	HIS	-	expression tag	UNP Q9NZ08
C	941	ASP	-	expression tag	UNP Q9NZ08
C	942	PRO	-	expression tag	UNP Q9NZ08
C	943	GLU	-	expression tag	UNP Q9NZ08
C	944	ALA	-	expression tag	UNP Q9NZ08
C	945	ASP	-	expression tag	UNP Q9NZ08
C	946	ALA	-	expression tag	UNP Q9NZ08
C	947	THR	-	expression tag	UNP Q9NZ08
C	948	GLY	-	expression tag	UNP Q9NZ08
C	949	LEU	-	expression tag	UNP Q9NZ08
C	950	GLU	-	expression tag	UNP Q9NZ08
C	951	ARG	-	expression tag	UNP Q9NZ08
C	952	MET	-	expression tag	UNP Q9NZ08
C	953	LEU	-	expression tag	UNP Q9NZ08
C	954	GLU	-	expression tag	UNP Q9NZ08
C	955	SER	-	expression tag	UNP Q9NZ08
C	956	ARG	-	expression tag	UNP Q9NZ08

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Chain	Residue	Modelled	Actual	Comment	Reference
C	957	GLY	-	expression tag	UNP Q9NZ08

- Molecule 2 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	D	5	61	34	2	25	0	0	0
2	F	5	61	34	2	25	0	0	0

- Molecule 3 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.

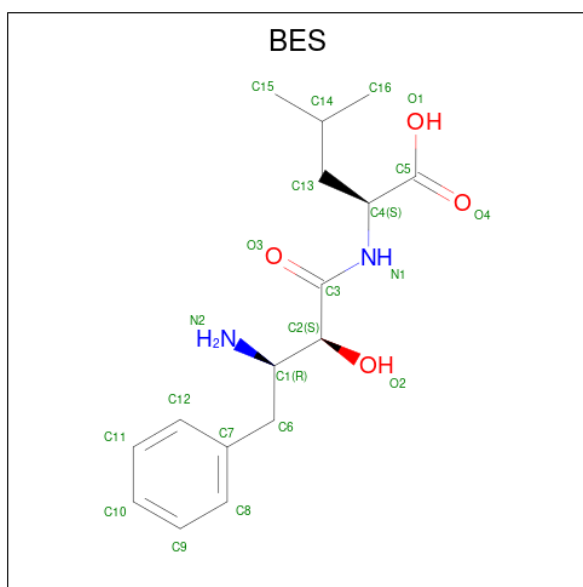


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
3	E	2	28	16	2	10	0	0	0

- Molecule 4 is ZINC ION (CCD ID: ZN) (formula: Zn).

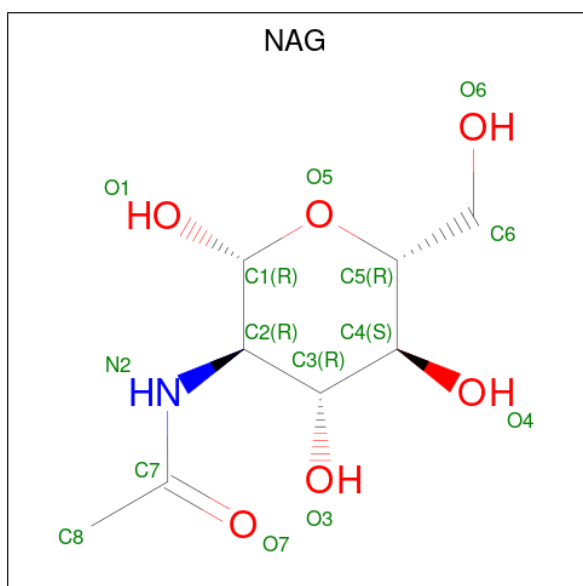
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Zn	0	0
			1	1		
4	B	1	Total	Zn	0	0
			1	1		
4	C	1	Total	Zn	0	0
			1	1		

- Molecule 5 is 2-(3-AMINO-2-HYDROXY-4-PHENYL-BUTYRYLAMINO)-4-METHYL-PENTANOIC ACID (CCD ID: BES) (formula: C₁₆H₂₄N₂O₄).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
5	A	1	22	16	2	4	0	0
5	B	1	22	16	2	4	0	0
5	C	1	22	16	2	4	0	0

- Molecule 6 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).

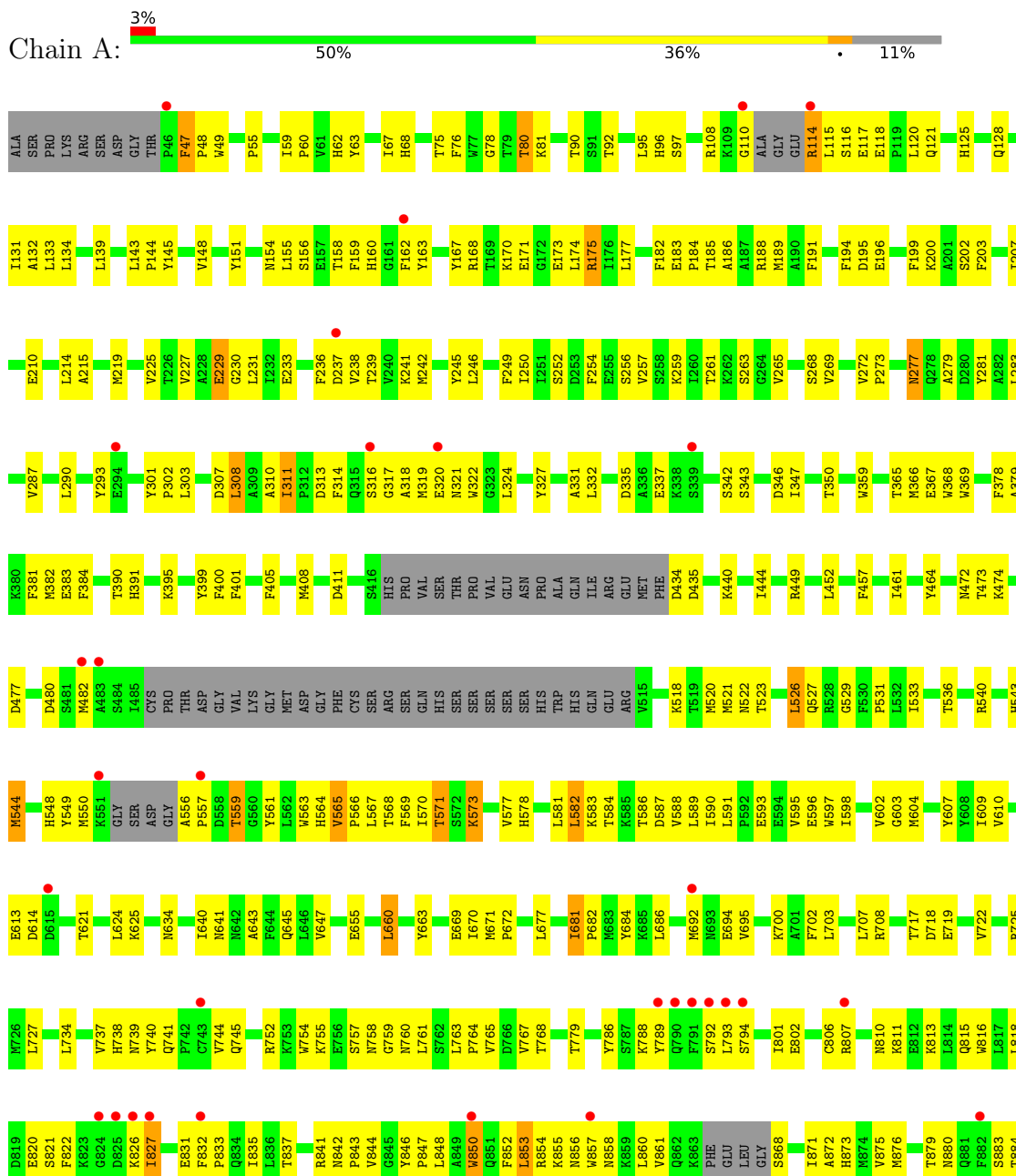


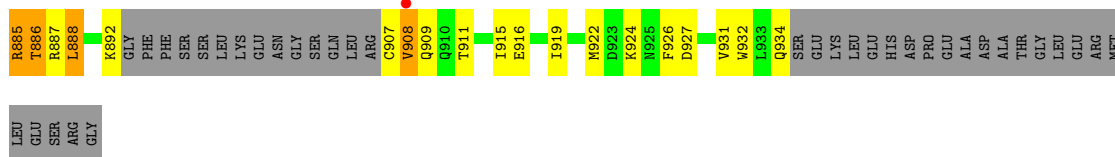
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total 14	C 8	N 1	O 5	0	0
6	A	1	Total 14	C 8	N 1	O 5	0	0
6	B	1	Total 14	C 8	N 1	O 5	0	0

3 Residue-property plots

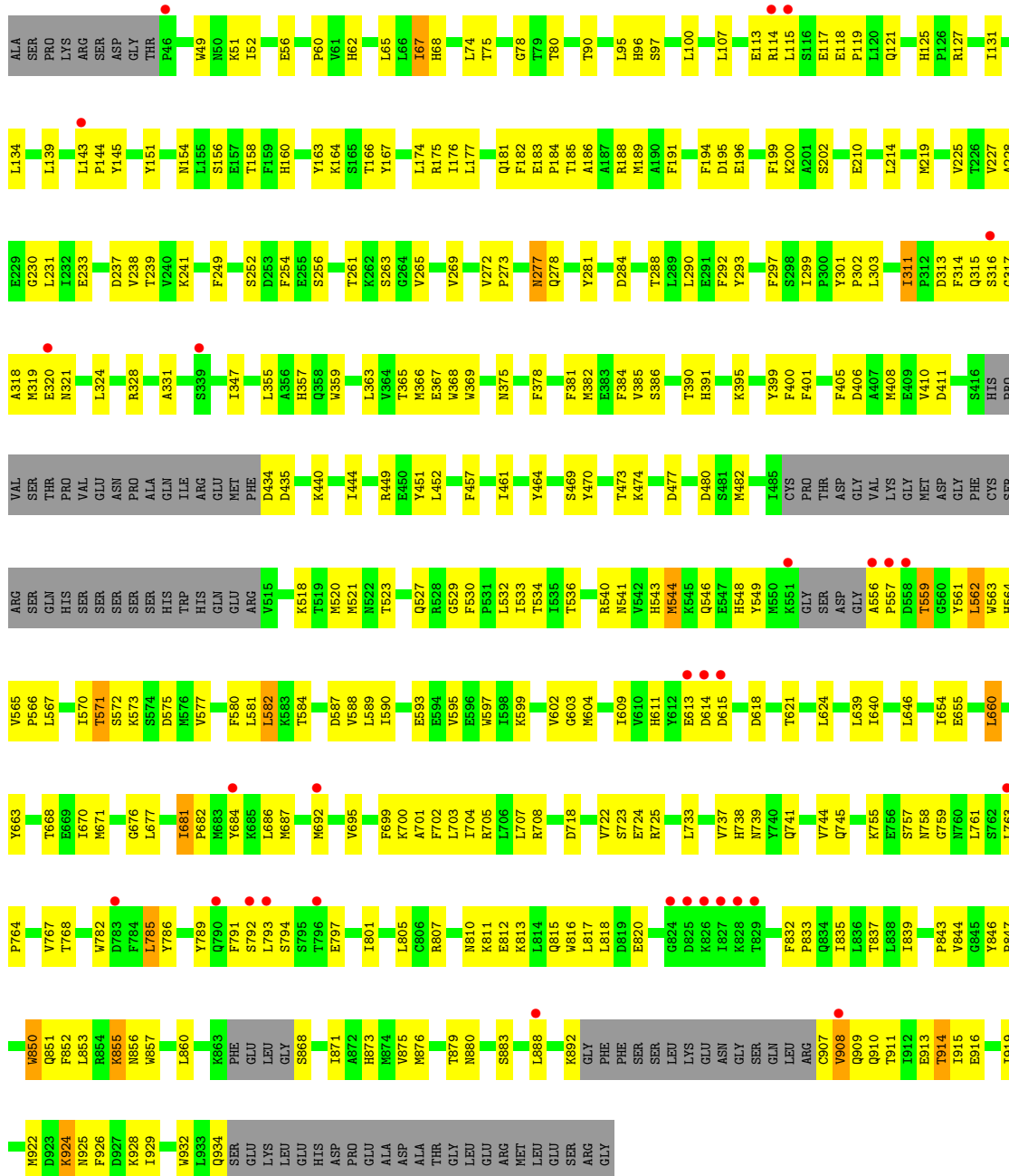
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: Endoplasmic reticulum aminopeptidase 1





• Molecule 1: Endoplasmic reticulum aminopeptidase 1



• Molecule 1: Endoplasmic reticulum aminopeptidase 1

- Molecule 2: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain F:  40% 20% 40%



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain E:  50% 50%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	71.03Å 234.63Å 95.86Å 90.00° 103.59° 90.00°	Depositor
Resolution (Å)	38.11 – 2.95 38.11 – 2.95	Depositor EDS
% Data completeness (in resolution range)	99.8 (38.11-2.95) 99.8 (38.11-2.95)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.53 (at 2.95Å)	Xtrriage
Refinement program	PHENIX 1.5_2	Depositor
R, R_{free}	0.199 , 0.264 0.195 , 0.260	Depositor DCC
R_{free} test set	3208 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	55.4	Xtrriage
Anisotropy	0.439	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 67.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	20163	wwPDB-VP
Average B, all atoms (Å ²)	69.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.77% 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, BMA, BES, NAG, MAN

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	0.33	0/6785	0.74	6/9195 (0.1%)
1	B	0.35	0/6804	0.75	2/9222 (0.0%)
1	C	0.33	0/6794	0.75	10/9207 (0.1%)
All	All	0.34	0/20383	0.75	18/27624 (0.1%)

There are no bond length outliers.

The worst 5 of 18 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	125	HIS	CA-C-N	5.81	125.95	119.32
1	C	125	HIS	C-N-CA	5.81	125.95	119.32
1	C	301	TYR	CA-C-N	5.79	125.65	119.28
1	C	301	TYR	C-N-CA	5.79	125.65	119.28
1	C	614	ASP	CB-CA-C	-5.78	109.89	116.54

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	6625	0	6567	342	0
1	B	6643	0	6582	334	0
1	C	6634	0	6574	336	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	61	0	52	4	0
2	F	61	0	52	1	0
3	E	28	0	25	1	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
5	A	22	0	22	2	0
5	B	22	0	22	9	0
5	C	22	0	23	6	0
6	A	28	0	26	1	0
6	B	14	0	13	0	0
All	All	20163	0	19958	1003	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 25.

The worst 5 of 1003 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:640:ILE:HD11	1:C:663:TYR:HE2	1.13	1.12
1:B:113:GLU:O	1:B:117:GLU:HG3	1.49	1.12
1:A:210:GLU:HG3	2:D:1:NAG:H83	1.37	1.07
1:A:875:VAL:HG11	1:A:908:VAL:CG2	1.86	1.04
1:C:640:ILE:HD11	1:C:663:TYR:CE2	1.95	1.02

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	804/921 (87%)	750 (93%)	50 (6%)	4 (0%)	24 49

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	B	809/921 (88%)	755 (93%)	52 (6%)	2 (0%)	43 66
1	C	805/921 (87%)	753 (94%)	51 (6%)	1 (0%)	48 72
All	All	2418/2763 (88%)	2258 (93%)	153 (6%)	7 (0%)	36 59

5 of 7 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	908	VAL
1	A	229	GLU
1	A	908	VAL
1	B	908	VAL
1	A	855	LYS

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	733/819 (90%)	706 (96%)	27 (4%)	30 55
1	B	734/819 (90%)	711 (97%)	23 (3%)	35 59
1	C	734/819 (90%)	714 (97%)	20 (3%)	39 64
All	All	2201/2457 (90%)	2131 (97%)	70 (3%)	34 59

5 of 70 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	395	LYS
1	C	478	LEU
1	C	593	GLU
1	A	888	LEU
1	A	886	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 62 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	278	GLN
1	C	645	GLN
1	B	810	ASN
1	C	641	ASN
1	C	810	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

12 monosaccharides are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	NAG	D	1	1,2	14,14,15	0.88	0	17,19,21	1.95	5 (29%)
2	NAG	D	2	2	14,14,15	0.63	0	17,19,21	2.18	3 (17%)
2	BMA	D	3	2	11,11,12	0.41	0	15,15,17	0.71	0
2	MAN	D	4	2	11,11,12	0.60	0	15,15,17	0.81	0
2	MAN	D	5	2	11,11,12	0.93	1 (9%)	15,15,17	1.61	3 (20%)
3	NAG	E	1	1,3	14,14,15	0.78	0	17,19,21	1.18	1 (5%)
3	NAG	E	2	3	14,14,15	0.65	0	17,19,21	2.23	3 (17%)
2	NAG	F	1	1,2	14,14,15	1.03	1 (7%)	17,19,21	1.86	6 (35%)
2	NAG	F	2	2	14,14,15	0.73	0	17,19,21	2.56	4 (23%)
2	BMA	F	3	2	11,11,12	0.39	0	15,15,17	0.70	0
2	MAN	F	4	2	11,11,12	0.46	0	15,15,17	0.73	0
2	MAN	F	5	2	11,11,12	0.79	0	15,15,17	1.16	2 (13%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	D	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	D	2	2	-	0/6/23/26	0/1/1/1
2	BMA	D	3	2	-	0/2/19/22	0/1/1/1
2	MAN	D	4	2	-	0/2/19/22	0/1/1/1
2	MAN	D	5	2	-	0/2/19/22	0/1/1/1
3	NAG	E	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	E	2	3	-	0/6/23/26	0/1/1/1
2	NAG	F	1	1,2	1/1/5/7	0/6/23/26	0/1/1/1
2	NAG	F	2	2	-	0/6/23/26	0/1/1/1
2	BMA	F	3	2	-	0/2/19/22	0/1/1/1
2	MAN	F	4	2	-	0/2/19/22	0/1/1/1
2	MAN	F	5	2	-	0/2/19/22	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	5	MAN	O5-C5	2.15	1.47	1.43
2	F	1	NAG	C4-C3	-2.07	1.47	1.52

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

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	2	NAG	O5-C5-C6	-7.68	92.71	107.66
2	D	2	NAG	O5-C5-C6	-6.81	94.40	107.66
3	E	2	NAG	O5-C5-C6	-6.74	94.55	107.66
2	F	1	NAG	O4-C4-C3	-4.09	100.75	110.38
2	F	2	NAG	C6-C5-C4	4.03	122.93	113.02

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	F	1	NAG	C1

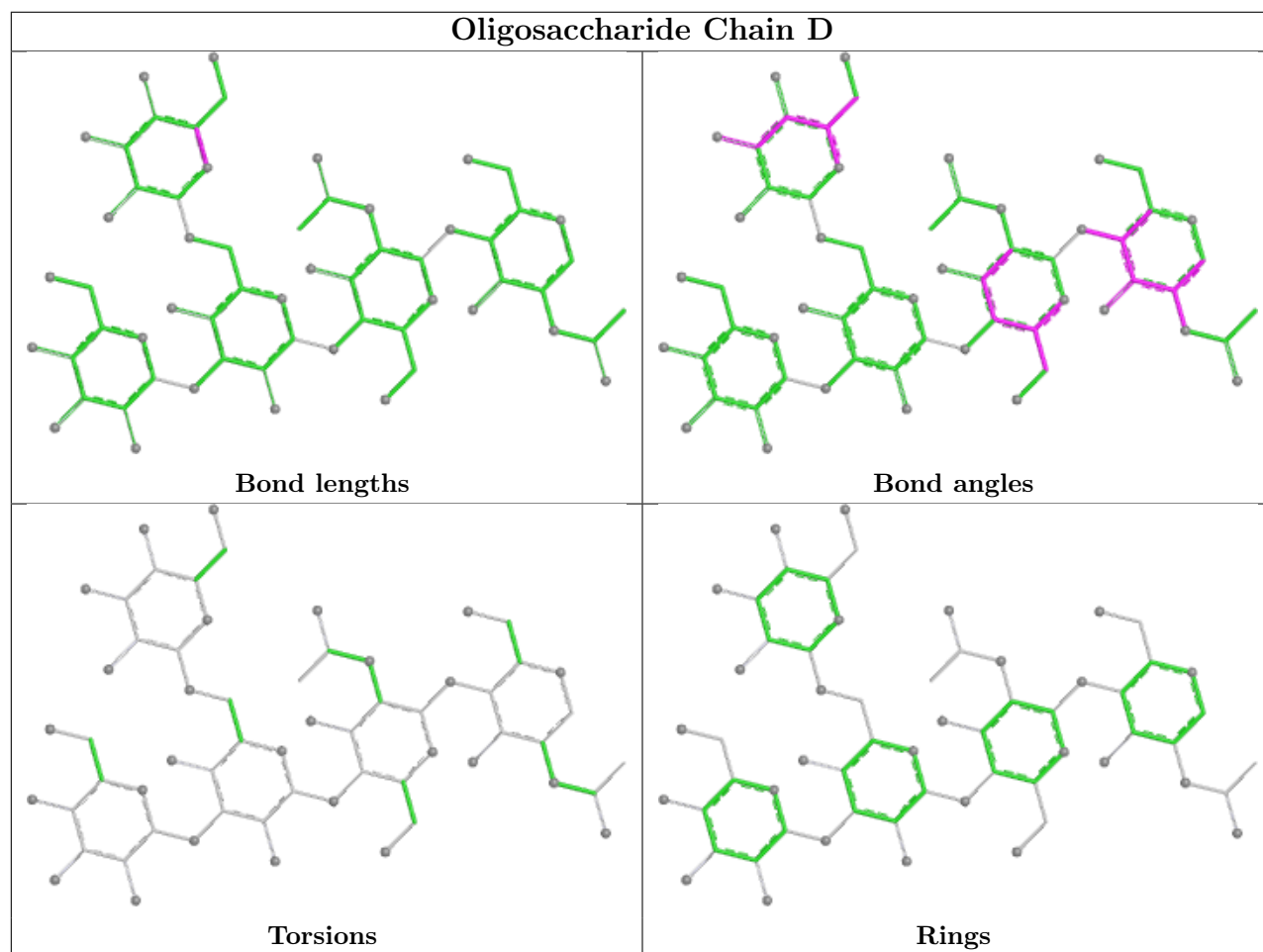
There are no torsion outliers.

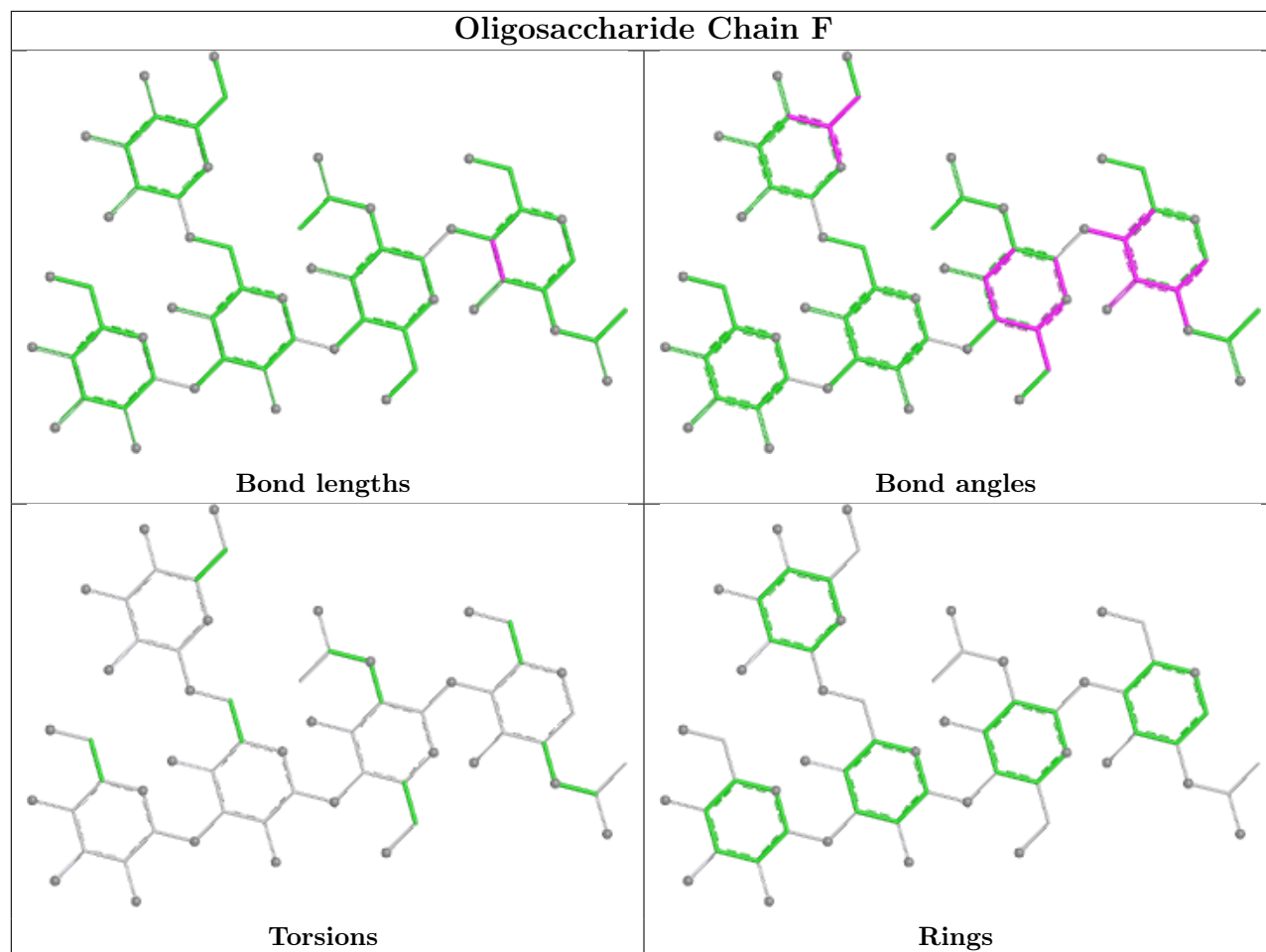
There are no ring outliers.

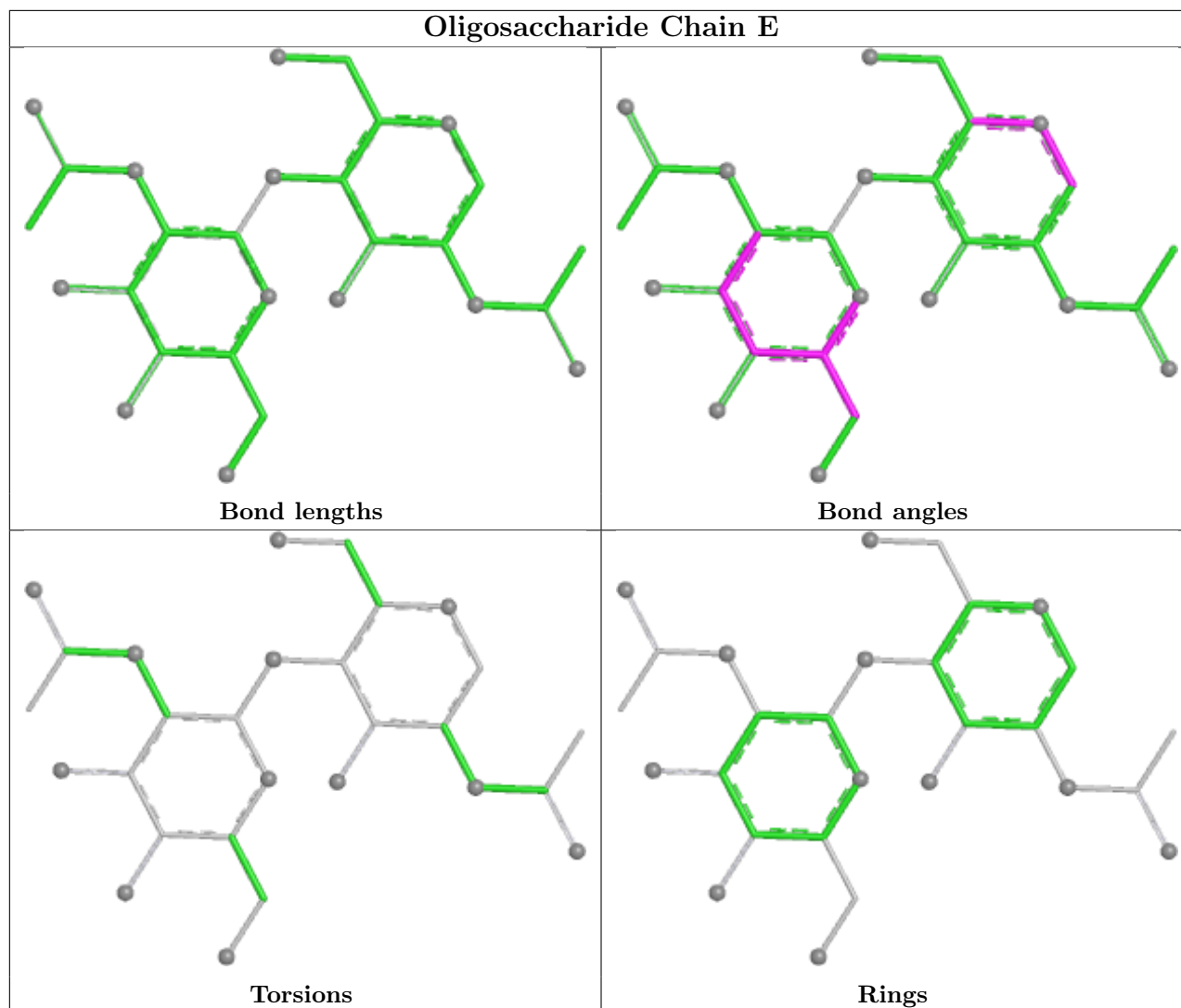
5 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	F	2	NAG	1	0
3	E	1	NAG	1	0
2	D	1	NAG	3	0
2	F	5	MAN	1	0
2	D	4	MAN	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 oligosaccharide.







5.6 Ligand geometry [i](#)

Of 9 ligands modelled in this entry, 3 are monoatomic - leaving 6 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	BES	C	1001	4	22,22,22	0.60	0	26,29,29	0.87	1 (3%)
5	BES	A	1001	4	22,22,22	0.62	0	26,29,29	0.84	1 (3%)
6	NAG	A	6001	1	14,14,15	1.29	2 (14%)	17,19,21	1.47	4 (23%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	NAG	A	6000	1	14,14,15	0.93	1 (7%)	17,19,21	1.39	2 (11%)
5	BES	B	1001	4	22,22,22	0.60	0	26,29,29	0.89	0
6	NAG	B	6000	1	14,14,15	0.67	0	17,19,21	0.99	1 (5%)

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	BES	C	1001	4	-	2/24/24/24	0/1/1/1
6	NAG	A	6001	1	1/1/5/7	2/6/23/26	0/1/1/1
5	BES	A	1001	4	-	2/24/24/24	0/1/1/1
6	NAG	A	6000	1	-	2/6/23/26	0/1/1/1
5	BES	B	1001	4	-	2/24/24/24	0/1/1/1
6	NAG	B	6000	1	-	2/6/23/26	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	6001	NAG	C1-C2	3.34	1.56	1.52
6	A	6001	NAG	O5-C1	2.62	1.48	1.43
6	A	6000	NAG	O5-C1	2.22	1.47	1.43

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

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	6000	NAG	C1-O5-C5	3.48	116.85	112.19
6	A	6001	NAG	O5-C1-C2	-3.12	106.46	111.29
6	A	6001	NAG	C1-O5-C5	2.77	115.90	112.19
6	A	6001	NAG	C6-C5-C4	-2.75	106.26	113.02
6	A	6000	NAG	O5-C1-C2	-2.28	107.77	111.29

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
6	A	6001	NAG	C1

5 of 12 torsion outliers are listed below:

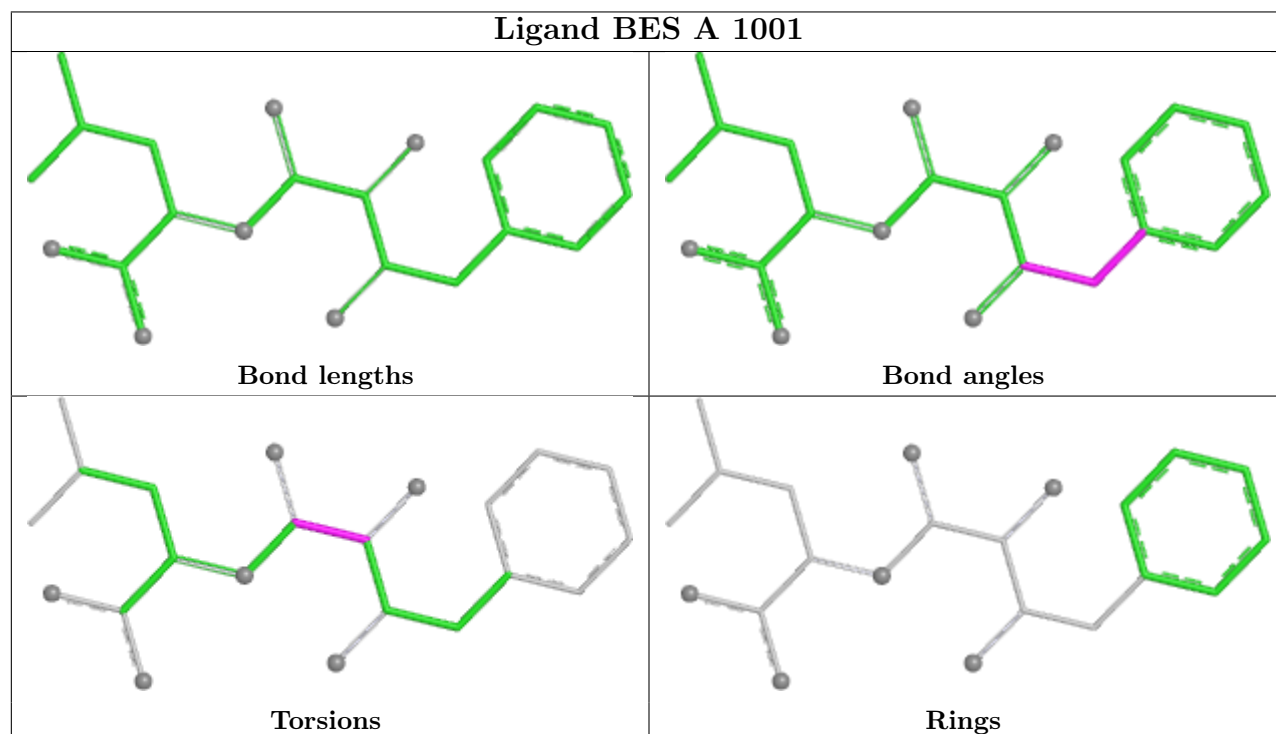
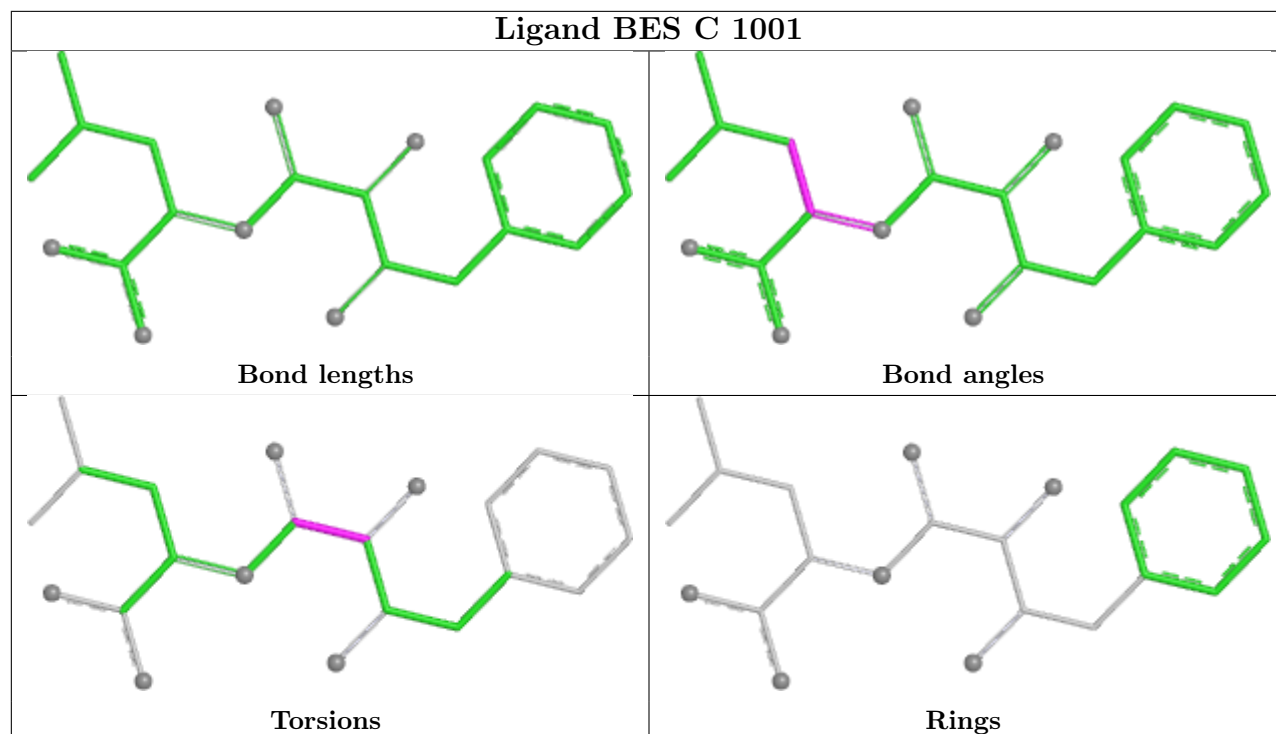
Mol	Chain	Res	Type	Atoms
6	A	6000	NAG	C8-C7-N2-C2
6	A	6000	NAG	O7-C7-N2-C2
6	A	6001	NAG	C8-C7-N2-C2
6	A	6001	NAG	O7-C7-N2-C2
6	B	6000	NAG	C8-C7-N2-C2

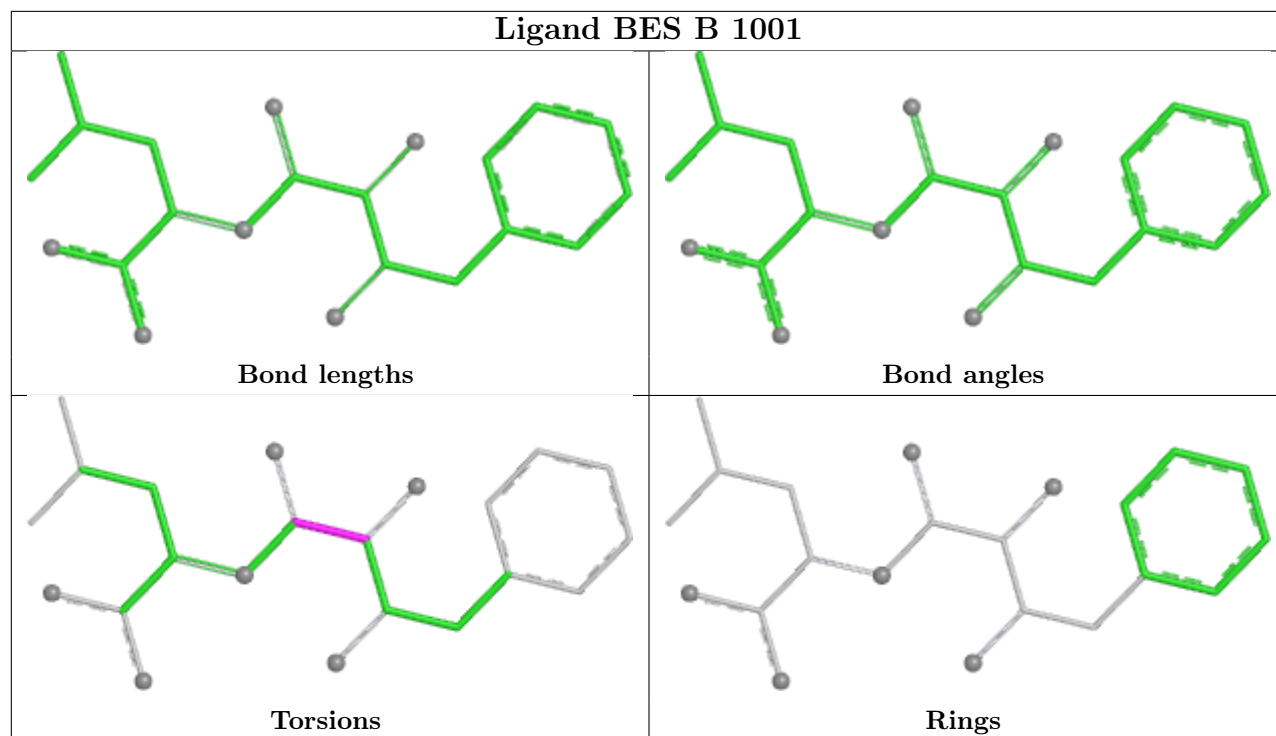
There are no ring outliers.

4 monomers are involved in 18 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	C	1001	BES	6	0
5	A	1001	BES	2	0
6	A	6001	NAG	1	0
5	B	1001	BES	9	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	818/921 (88%)	-0.05	32 (3%) 43 35	24, 56, 128, 238	11 (1%)
1	B	821/921 (89%)	-0.02	30 (3%) 45 37	24, 56, 127, 223	11 (1%)
1	C	818/921 (88%)	-0.06	25 (3%) 51 44	23, 57, 127, 222	12 (1%)
All	All	2457/2763 (88%)	-0.05	87 (3%) 47 40	23, 56, 127, 238	34 (1%)

The worst 5 of 87 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	114	ARG	4.6
1	B	824	GLY	4.6
1	B	825	ASP	4.2
1	B	613	GLU	4.0
1	C	320	GLU	3.9

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. 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.

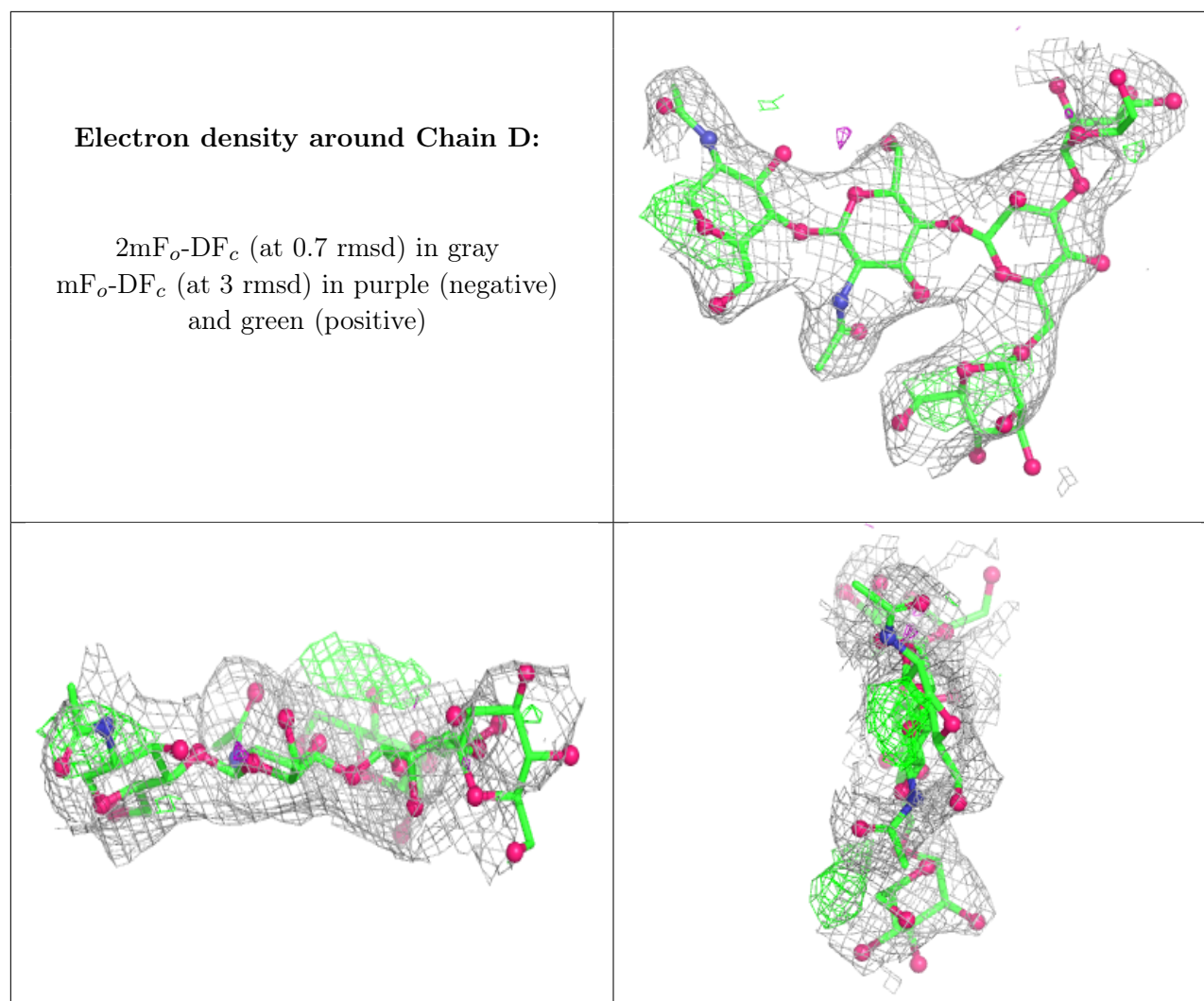
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	NAG	E	2	14/15	0.40	0.16	118,161,189,212	0
2	MAN	D	4	11/12	0.50	0.16	111,168,187,189	0
2	BMA	F	3	11/12	0.63	0.21	43,135,191,223	0

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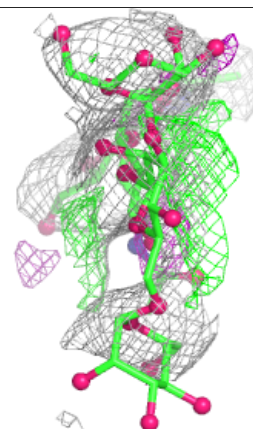
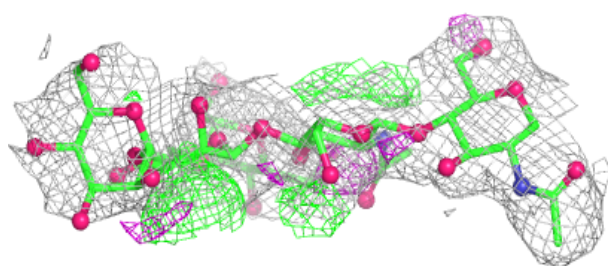
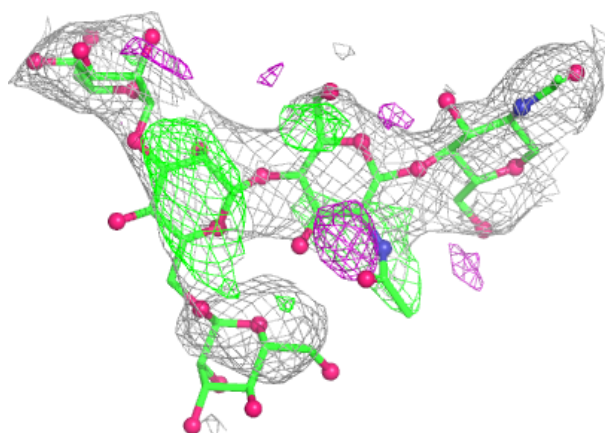
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	MAN	F	5	11/12	0.70	0.16	86,142,212,223	0
2	MAN	D	5	11/12	0.72	0.14	33,95,172,209	0
2	NAG	F	2	14/15	0.78	0.25	47,135,257,284	0
2	MAN	F	4	11/12	0.81	0.17	75,95,103,110	0
2	NAG	D	2	14/15	0.82	0.10	60,85,108,108	0
2	NAG	F	1	14/15	0.86	0.15	20,79,110,127	0
2	NAG	D	1	14/15	0.86	0.15	49,90,110,161	0
3	NAG	E	1	14/15	0.87	0.17	50,93,110,139	0
2	BMA	D	3	11/12	0.87	0.11	51,86,129,131	0

The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

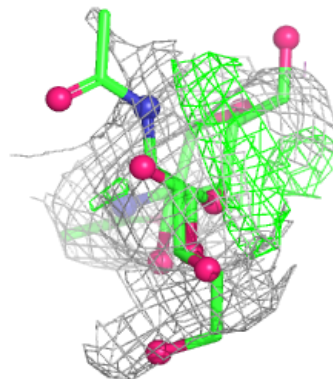
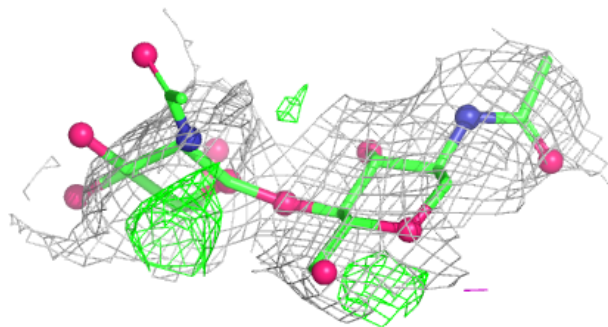
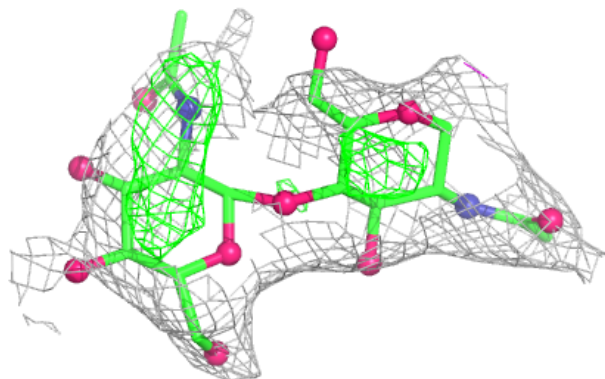


Electron density around Chain F:

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

**Electron density around Chain E:**

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

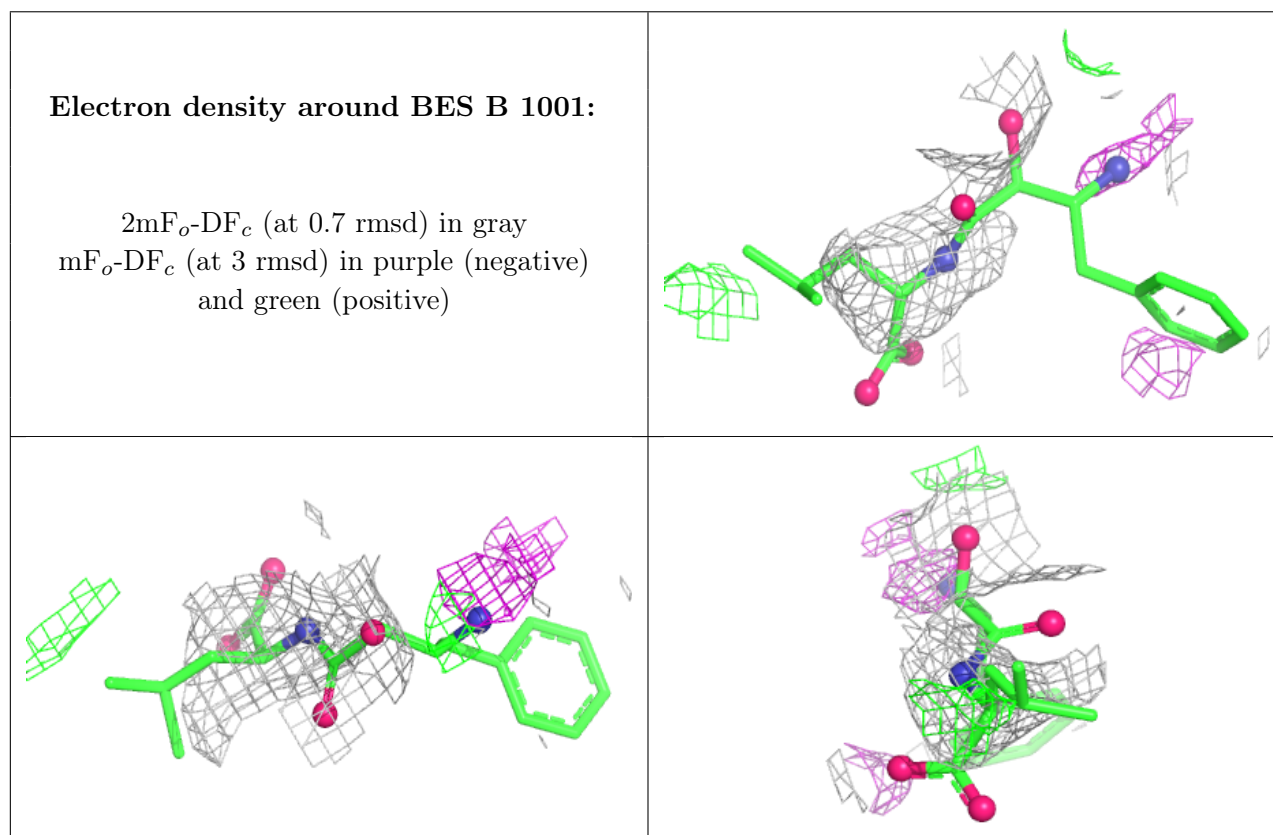


6.4 Ligands [i](#)

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.

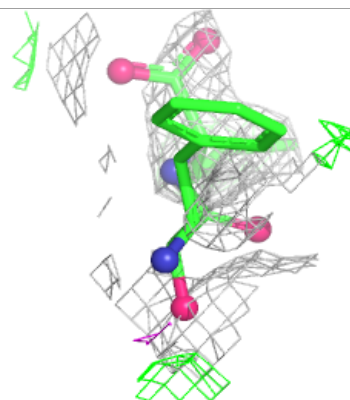
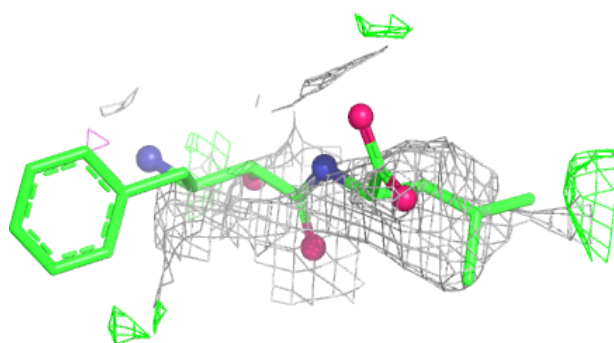
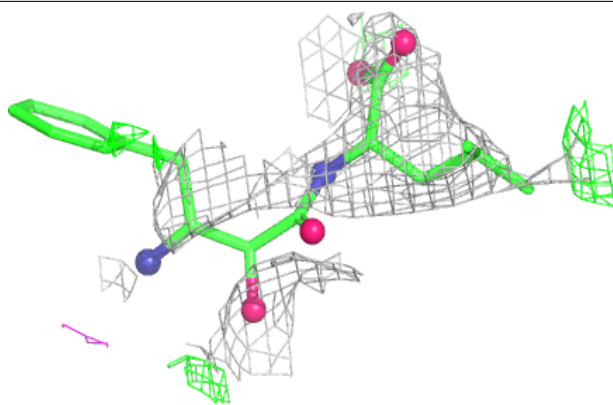
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
6	NAG	B	6000	14/15	0.55	0.14	59,128,146,153	0
6	NAG	A	6001	14/15	0.57	0.13	72,140,150,153	0
6	NAG	A	6000	14/15	0.61	0.17	76,119,141,145	0
5	BES	B	1001	22/22	0.80	0.26	121,167,185,218	0
5	BES	C	1001	22/22	0.81	0.24	79,178,213,218	0
5	BES	A	1001	22/22	0.81	0.28	86,184,208,217	0
4	ZN	A	1000	1/1	1.00	0.03	40,40,40,40	0
4	ZN	B	1000	1/1	1.00	0.01	38,38,38,38	0
4	ZN	C	1000	1/1	1.00	0.03	38,38,38,38	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.

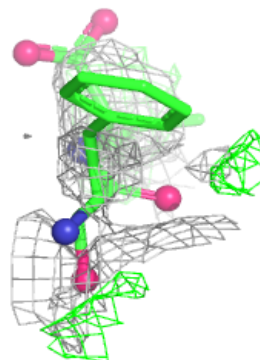
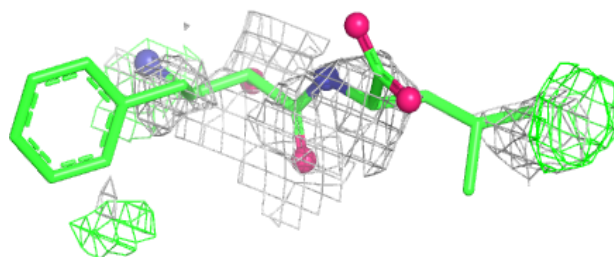
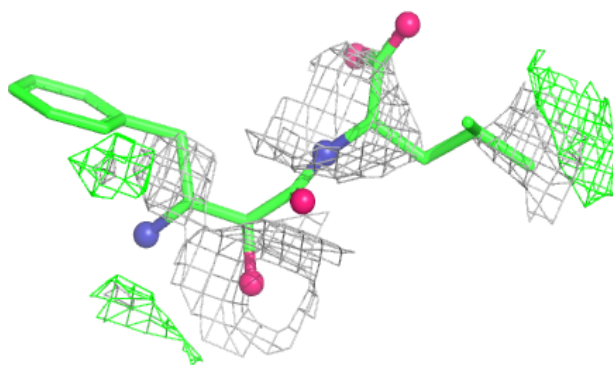


Electron density around BES C 1001:

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

**Electron density around BES 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)



6.5 Other polymers [i](#)

There are no such residues in this entry.