



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 10:49 AM UTC

PDB ID : 6R4L / pdb_00006r4l
Title : Crystal structure of S. cerevisia Niemann-Pick type C protein NCR1
Authors : Winkler, M.B.L.; Kidmose, R.T.; Pedersen, B.P.
Deposited on : 2019-03-22
Resolution : 3.50 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

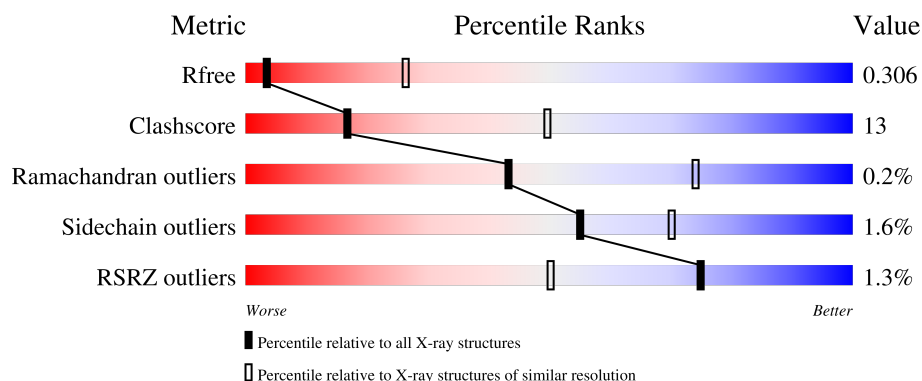
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

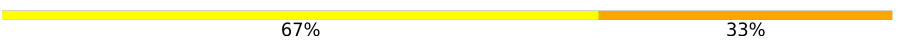
The reported resolution of this entry is 3.50 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1085 (3.54-3.46)
Clashscore	190562	1140 (3.54-3.46)
Ramachandran outliers	187476	1113 (3.54-3.46)
Sidechain outliers	187428	1114 (3.54-3.46)
RSRZ outliers	180081	1084 (3.54-3.46)

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	1198	 67% 24% 9%
2	B	3	 67% 33%
2	C	3	 67% 33%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 8894 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 NPC intracellular cholesterol transporter 1-related protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	1096	Total	C	N	O	S	0	0	0
			8759	5706	1390	1606	57			

There are 28 discrepancies between the modelled and reference sequences:

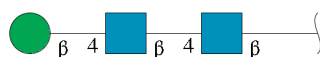
Chain	Residue	Modelled	Actual	Comment	Reference
A	1171	LEU	-	expression tag	UNP Q12200
A	1172	VAL	-	expression tag	UNP Q12200
A	1173	PRO	-	expression tag	UNP Q12200
A	1174	ARG	-	expression tag	UNP Q12200
A	1175	GLY	-	expression tag	UNP Q12200
A	1176	SER	-	expression tag	UNP Q12200
A	1177	GLY	-	expression tag	UNP Q12200
A	1178	GLY	-	expression tag	UNP Q12200
A	1179	GLY	-	expression tag	UNP Q12200
A	1180	GLY	-	expression tag	UNP Q12200
A	1181	SER	-	expression tag	UNP Q12200
A	1182	GLY	-	expression tag	UNP Q12200
A	1183	GLY	-	expression tag	UNP Q12200
A	1184	GLY	-	expression tag	UNP Q12200
A	1185	GLY	-	expression tag	UNP Q12200
A	1186	SER	-	expression tag	UNP Q12200
A	1187	GLY	-	expression tag	UNP Q12200
A	1188	GLY	-	expression tag	UNP Q12200
A	1189	HIS	-	expression tag	UNP Q12200
A	1190	HIS	-	expression tag	UNP Q12200
A	1191	HIS	-	expression tag	UNP Q12200
A	1192	HIS	-	expression tag	UNP Q12200
A	1193	HIS	-	expression tag	UNP Q12200
A	1194	HIS	-	expression tag	UNP Q12200
A	1195	HIS	-	expression tag	UNP Q12200
A	1196	HIS	-	expression tag	UNP Q12200
A	1197	HIS	-	expression tag	UNP Q12200

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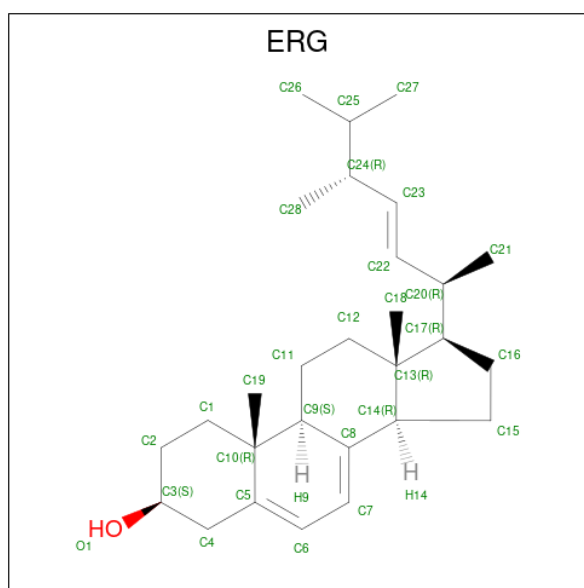
Chain	Residue	Modelled	Actual	Comment	Reference
A	1198	HIS	-	expression tag	UNP Q12200

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



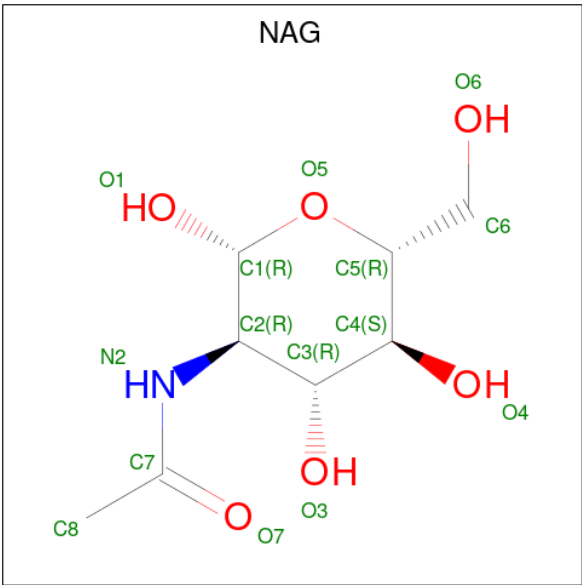
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	3	Total	C	N	O	0	0	0
			39	22	2	15			
2	C	3	Total	C	N	O	0	0	0
			39	22	2	15			

- Molecule 3 is ERGOSTEROL (CCD ID: ERG) (formula: $C_{28}H_{44}O$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			29	28	1		

- Molecule 4 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).

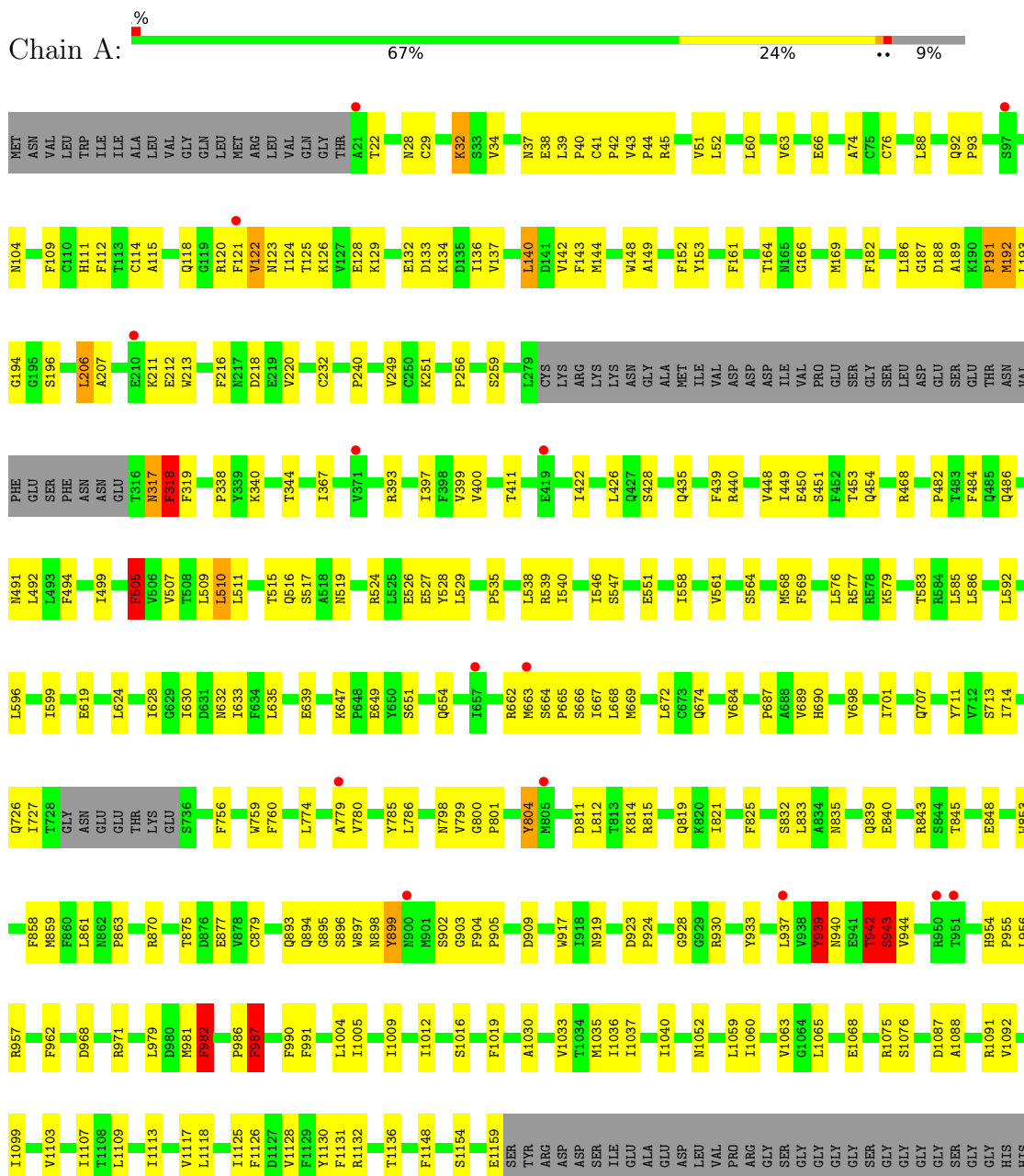


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			14	8	1	5		
4	A	1	Total	C	N	O	0	0
			14	8	1	5		

3 Residue-property plots [i](#)

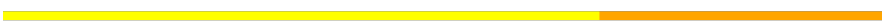
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: NPC intracellular cholesterol transporter 1-related protein 1



HIS
HIS
HIS
HIS
HIS
HIS
HIS
HIS

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

Chain B:  67% 33%

NAG1
NAG2
BMA3

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

Chain C:  67% 33%

NAG1
NAG2
BMA3

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	148.67Å 90.06Å 161.68Å 90.00° 110.31° 90.00°	Depositor
Resolution (Å)	49.12 – 3.50 49.12 – 3.50	Depositor EDS
% Data completeness (in resolution range)	93.6 (49.12-3.50) 93.5 (49.12-3.50)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.18 (at 3.48Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.266 , 0.301 0.267 , 0.306	Depositor DCC
R_{free} test set	1209 reflections (4.75%)	wwPDB-VP
Wilson B-factor (Å ²)	134.5	Xtriage
Anisotropy	0.368	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.25 , 97.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	8894	wwPDB-VP
Average B, all atoms (Å ²)	157.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.16% 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: BMA, NAG, ERG

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	0.41	4/8977 (0.0%)	0.65	7/12185 (0.1%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	7

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	317	ASN	CB-CG	5.49	1.65	1.52
1	A	468	ARG	CG-CD	-5.40	1.36	1.52
1	A	1159	GLU	CG-CD	5.09	1.64	1.52
1	A	1159	GLU	CD-OE2	5.02	1.34	1.25

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	943	SER	N-CA-CB	-7.72	97.44	110.49
1	A	192	MET	N-CA-CB	-7.12	98.47	110.49
1	A	32	LYS	CA-CB-CG	6.85	127.81	114.10
1	A	192	MET	CA-CB-CG	6.12	126.33	114.10
1	A	942	THR	N-CA-C	-5.68	102.11	110.52
1	A	982	PHE	CB-CA-C	5.18	121.32	111.17
1	A	468	ARG	NE-CZ-NH1	-5.04	116.47	121.50

There are no chirality outliers.

All (7) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	191	PRO	Peptide
1	A	318	PHE	Sidechain
1	A	505	PHE	Sidechain
1	A	899	TYR	Sidechain
1	A	939	TYR	Sidechain
1	A	942	THR	Peptide
1	A	987	PHE	Sidechain

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	8759	0	8670	223	0
2	B	39	0	34	1	0
2	C	39	0	34	3	0
3	A	29	0	44	5	0
4	A	28	0	26	4	0
All	All	8894	0	8808	226	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 13.

All (226) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:212:GLU:HG2	4:A:1208:NAG:H82	1.50	0.93
1:A:517:SER:HB3	2:C:1:NAG:H83	1.53	0.90
1:A:51:VAL:HG21	1:A:74:ALA:H	1.42	0.85
1:A:426:LEU:HD13	1:A:528:TYR:HB2	1.60	0.83
1:A:451:SER:HB3	1:A:507:VAL:HG23	1.65	0.79
1:A:632:ASN:HD21	1:A:667:ILE:HD12	1.49	0.78
1:A:450:GLU:OE2	3:A:1201:ERG:H191	1.84	0.76
1:A:393:ARG:HD2	1:A:511:LEU:HD11	1.67	0.75
1:A:39:LEU:HD13	1:A:193:LEU:HD22	1.67	0.74
1:A:122:VAL:HG11	1:A:212:GLU:CD	2.13	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:774:LEU:HB2	1:A:1052:ASN:HD22	1.53	0.74
1:A:684:VAL:HG21	1:A:689:VAL:HG13	1.72	0.72
1:A:897:TRP:CZ3	1:A:899:TYR:HD2	2.08	0.71
1:A:454:GLN:HB3	1:A:482:PRO:HB3	1.72	0.71
1:A:635:LEU:HD22	1:A:663:MET:HE1	1.72	0.71
1:A:111:HIS:HA	1:A:115:ALA:HB3	1.71	0.71
1:A:422:ILE:HD12	1:A:509:LEU:HD21	1.71	0.71
1:A:940:ASN:HB3	1:A:942:THR:O	1.90	0.70
1:A:1128:VAL:HG22	1:A:1132:ARG:HH12	1.56	0.70
1:A:212:GLU:HG2	4:A:1208:NAG:C8	2.21	0.70
1:A:164:THR:HG23	1:A:166:GLY:H	1.56	0.69
1:A:129:LYS:HE3	1:A:132:GLU:HB2	1.73	0.69
1:A:801:PRO:HB3	3:A:1201:ERG:H272	1.75	0.69
1:A:492:LEU:HD12	1:A:799:VAL:HG21	1.74	0.69
1:A:576:LEU:HD11	1:A:579:LYS:HD2	1.76	0.68
1:A:206:LEU:HD23	1:A:206:LEU:H	1.59	0.67
1:A:256:PRO:HB2	1:A:259:SER:HB2	1.75	0.67
1:A:51:VAL:HG12	1:A:52:LEU:H	1.59	0.67
1:A:125:THR:H	1:A:140:LEU:HD11	1.59	0.67
1:A:632:ASN:ND2	1:A:667:ILE:HD12	2.09	0.67
1:A:812:LEU:HG	1:A:943:SER:HB3	1.77	0.67
1:A:939:TYR:O	1:A:939:TYR:HD1	1.80	0.65
1:A:122:VAL:HG11	1:A:212:GLU:OE2	1.97	0.64
1:A:630:ILE:HD13	1:A:1109:LEU:HD21	1.80	0.64
1:A:968:ASP:HA	1:A:971:ARG:HG2	1.79	0.64
1:A:318:PHE:HD1	1:A:318:PHE:O	1.80	0.64
1:A:940:ASN:H	1:A:944:VAL:HA	1.62	0.63
1:A:393:ARG:HB2	1:A:511:LEU:HG	1.78	0.63
1:A:987:PHE:CD1	1:A:987:PHE:C	2.76	0.63
1:A:515:THR:HG22	1:A:516:GLN:H	1.62	0.63
1:A:917:TRP:HE1	1:A:928:GLY:HA3	1.64	0.63
1:A:92:GLN:HG3	1:A:93:PRO:HD3	1.81	0.62
1:A:756:PHE:CD2	1:A:1148:PHE:HB2	2.33	0.62
1:A:428:SER:HA	1:A:524:ARG:HG2	1.81	0.62
1:A:129:LYS:HE3	1:A:132:GLU:CB	2.29	0.62
1:A:812:LEU:HG	1:A:943:SER:CB	2.30	0.62
1:A:863:PRO:HG3	1:A:894:GLN:HA	1.82	0.62
1:A:435:GLN:O	1:A:440:ARG:NH1	2.31	0.61
1:A:120:ARG:HD3	1:A:216:PHE:HE1	1.65	0.61
1:A:774:LEU:HB2	1:A:1052:ASN:ND2	2.14	0.61
1:A:249:VAL:HG12	1:A:251:LYS:H	1.65	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:505:PHE:CD1	1:A:505:PHE:C	2.78	0.60
1:A:780:VAL:HG11	1:A:786:LEU:HB2	1.83	0.60
1:A:919:ASN:O	1:A:930:ARG:NH1	2.34	0.60
1:A:666:SER:HB2	1:A:1075:ARG:HG2	1.83	0.60
1:A:118:GLN:OE1	1:A:120:ARG:NH2	2.35	0.59
1:A:547:SER:HB3	1:A:786:LEU:HD11	1.84	0.58
1:A:982:PHE:CD1	1:A:982:PHE:C	2.81	0.58
1:A:711:TYR:HA	1:A:714:ILE:HG22	1.85	0.58
1:A:1009:ILE:HD12	1:A:1033:VAL:HG13	1.86	0.58
1:A:937:LEU:CD1	1:A:939:TYR:HB3	2.34	0.57
1:A:726:GLN:OE1	1:A:726:GLN:N	2.33	0.57
1:A:212:GLU:CG	4:A:1208:NAG:H82	2.29	0.57
1:A:651:SER:HB3	1:A:654:GLN:OE1	2.05	0.57
1:A:125:THR:N	1:A:140:LEU:HD11	2.20	0.56
1:A:515:THR:HB	2:C:1:NAG:H81	1.85	0.56
1:A:923:ASP:HB3	1:A:924:PRO:HD3	1.88	0.56
1:A:111:HIS:NE2	1:A:218:ASP:OD2	2.38	0.56
1:A:367:ILE:HD12	1:A:990:PHE:HZ	1.69	0.56
1:A:524:ARG:NH2	1:A:527:GLU:OE2	2.39	0.56
2:B:1:NAG:H82	2:B:1:NAG:H3	1.87	0.56
1:A:962:PHE:HB3	1:A:991:PHE:CE2	2.40	0.56
1:A:28:ASN:HA	1:A:40:PRO:HA	1.88	0.55
1:A:647:LYS:HE2	1:A:649:GLU:OE2	2.06	0.55
1:A:568:MET:SD	1:A:1113:ILE:HB	2.47	0.55
1:A:448:VAL:CG2	1:A:510:LEU:HD11	2.36	0.55
1:A:451:SER:OG	1:A:453:THR:OG1	2.24	0.55
1:A:484:PHE:HD2	1:A:486:GLN:HB2	1.71	0.55
1:A:39:LEU:HB2	1:A:136:ILE:HG21	1.89	0.55
1:A:624:LEU:O	1:A:628:ILE:HG22	2.07	0.55
1:A:870:ARG:HG3	1:A:879:CYS:HB2	1.89	0.55
1:A:987:PHE:C	1:A:987:PHE:HD1	2.15	0.55
1:A:1076:SER:HB3	1:A:1099:ILE:HD12	1.89	0.54
1:A:411:THR:HG22	1:A:538:LEU:HD21	1.90	0.53
1:A:484:PHE:CD2	1:A:486:GLN:HB2	2.43	0.53
1:A:397:ILE:HB	1:A:507:VAL:HG13	1.90	0.53
1:A:88:LEU:HD21	1:A:109:PHE:HD2	1.74	0.53
1:A:448:VAL:HG23	1:A:510:LEU:HD11	1.91	0.53
1:A:129:LYS:HB3	1:A:132:GLU:HB2	1.90	0.53
1:A:505:PHE:C	1:A:505:PHE:HD1	2.16	0.53
1:A:982:PHE:C	1:A:982:PHE:HD1	2.14	0.53
1:A:42:PRO:HG3	1:A:137:VAL:HB	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:188:ASP:HA	1:A:194:GLY:HA3	1.91	0.53
1:A:801:PRO:HD2	1:A:956:LEU:HD12	1.91	0.52
1:A:624:LEU:HD12	1:A:1065:LEU:HD21	1.92	0.52
1:A:122:VAL:HG11	1:A:212:GLU:OE1	2.10	0.51
1:A:979:LEU:HB2	1:A:981:MET:HE2	1.92	0.51
1:A:526:GLU:OE2	1:A:785:TYR:OH	2.27	0.51
1:A:187:GLY:HA2	1:A:196:SER:HA	1.93	0.51
1:A:1118:LEU:HB2	1:A:1131:PHE:CE1	2.46	0.51
1:A:367:ILE:HD12	1:A:990:PHE:CZ	2.44	0.51
1:A:875:THR:HG23	1:A:877:GLU:H	1.74	0.51
1:A:232:CYS:HB3	1:A:835:ASN:OD1	2.10	0.50
1:A:811:ASP:HB2	1:A:814:LYS:HE3	1.93	0.50
1:A:954:HIS:O	1:A:956:LEU:N	2.44	0.50
1:A:207:ALA:HB2	1:A:212:GLU:HG3	1.94	0.50
1:A:639:GLU:OE2	1:A:662:ARG:NH2	2.42	0.50
1:A:672:LEU:HD22	1:A:1019:PHE:HE2	1.77	0.50
1:A:804:TYR:N	1:A:804:TYR:CD1	2.80	0.50
1:A:1009:ILE:HA	1:A:1012:ILE:HD12	1.94	0.50
1:A:189:ALA:HB1	1:A:191:PRO:HD2	1.94	0.50
1:A:519:ASN:HB3	1:A:546:ILE:HD13	1.94	0.50
1:A:43:VAL:HG22	1:A:44:PRO:HD2	1.94	0.49
1:A:144:MET:HE3	1:A:149:ALA:HB2	1.94	0.49
1:A:535:PRO:HB2	1:A:538:LEU:HD13	1.93	0.49
1:A:1040:ILE:HG12	1:A:1136:THR:HG22	1.94	0.49
1:A:104:ASN:HD22	1:A:220:VAL:HG23	1.77	0.49
1:A:148:TRP:NE1	1:A:220:VAL:HB	2.28	0.49
1:A:152:PHE:HE2	1:A:182:PHE:HE2	1.59	0.49
1:A:630:ILE:HA	1:A:633:ILE:HD12	1.94	0.49
1:A:669:MET:HE3	1:A:1016:SER:HA	1.95	0.49
1:A:687:PRO:HA	1:A:690:HIS:HB3	1.95	0.49
1:A:32:LYS:CE	1:A:134:LYS:HE2	2.42	0.49
1:A:124:ILE:HA	1:A:140:LEU:HG	1.95	0.49
1:A:845:THR:OG1	1:A:968:ASP:OD2	2.23	0.48
1:A:1004:LEU:HD21	1:A:1060:ILE:HD11	1.95	0.48
1:A:112:PHE:HE1	1:A:121:PHE:CD1	2.31	0.48
1:A:1088:ALA:HA	1:A:1091:ARG:HE	1.78	0.48
1:A:904:PHE:CG	1:A:905:PRO:HD2	2.49	0.48
1:A:111:HIS:O	1:A:120:ARG:HD2	2.14	0.48
1:A:898:ASN:ND2	1:A:903:GLY:HA3	2.28	0.48
1:A:1005:ILE:HD13	1:A:1037:ILE:HG21	1.94	0.48
1:A:893:GLN:HG3	1:A:895:GLY:H	1.78	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:76:CYS:HB3	1:A:114:CYS:HA	1.96	0.47
1:A:596:LEU:O	1:A:599:ILE:HG13	2.13	0.47
1:A:897:TRP:CH2	1:A:899:TYR:HD2	2.32	0.47
1:A:169:MET:HE1	1:A:182:PHE:CD1	2.49	0.47
1:A:577:ARG:HG3	1:A:583:THR:O	2.14	0.47
1:A:815:ARG:NH2	1:A:902:SER:O	2.47	0.47
1:A:585:LEU:HD13	1:A:586:LEU:HD13	1.95	0.47
1:A:400:VAL:HG13	1:A:539:ARG:HB2	1.97	0.47
1:A:619:GLU:HG2	1:A:1126:PHE:HE1	1.79	0.47
1:A:815:ARG:O	1:A:819:GLN:HG3	2.15	0.47
1:A:1092:VAL:HG21	1:A:1154:SER:HB2	1.96	0.47
1:A:825:PHE:HE1	1:A:859:MET:HE2	1.80	0.47
1:A:897:TRP:CZ3	1:A:899:TYR:CD2	2.98	0.47
1:A:109:PHE:CE1	1:A:186:LEU:HD11	2.51	0.46
1:A:207:ALA:CB	1:A:212:GLU:HG3	2.46	0.46
1:A:439:PHE:HD2	1:A:448:VAL:HG21	1.79	0.46
1:A:664:SER:N	1:A:665:PRO:HD2	2.31	0.46
1:A:211:LYS:O	1:A:212:GLU:HB3	2.16	0.46
1:A:1117:VAL:HG13	1:A:1126:PHE:HE2	1.80	0.46
4:A:1208:NAG:O7	4:A:1208:NAG:H3	2.15	0.46
1:A:318:PHE:O	1:A:318:PHE:CD1	2.64	0.46
1:A:801:PRO:HG3	3:A:1201:ERG:H24	1.96	0.46
1:A:897:TRP:H	1:A:897:TRP:CD1	2.32	0.46
1:A:547:SER:O	1:A:551:GLU:HG2	2.15	0.46
1:A:60:LEU:HD11	1:A:66:GLU:HA	1.97	0.46
1:A:499:ILE:HD12	1:A:499:ILE:H	1.81	0.46
1:A:564:SER:OG	1:A:1113:ILE:HG13	2.15	0.45
1:A:672:LEU:HD22	1:A:1019:PHE:CE2	2.51	0.45
1:A:825:PHE:HE2	1:A:858:PHE:HD2	1.64	0.45
1:A:1033:VAL:O	1:A:1036:ILE:HG13	2.16	0.45
1:A:120:ARG:HD3	1:A:216:PHE:CE1	2.49	0.45
1:A:140:LEU:HA	1:A:140:LEU:HD13	1.63	0.45
1:A:494:PHE:O	1:A:798:ASN:ND2	2.49	0.45
1:A:800:GLY:O	3:A:1201:ERG:H161	2.16	0.45
1:A:861:LEU:HD13	1:A:904:PHE:CZ	2.52	0.45
1:A:632:ASN:OD1	1:A:707:GLN:NE2	2.49	0.45
1:A:843:ARG:HH22	1:A:971:ARG:HH12	1.64	0.45
1:A:904:PHE:CD2	1:A:905:PRO:HD2	2.51	0.45
1:A:1087:ASP:O	1:A:1091:ARG:HD3	2.17	0.45
1:A:126:LYS:H	1:A:140:LEU:HD21	1.82	0.45
1:A:893:GLN:HB3	1:A:896:SER:OG	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1126:PHE:HA	1:A:1130:TYR:HD2	1.82	0.45
1:A:121:PHE:CD1	1:A:142:VAL:HG22	2.52	0.44
1:A:760:PHE:CD1	1:A:1035:MET:HG3	2.53	0.44
1:A:51:VAL:HG12	1:A:52:LEU:N	2.28	0.44
1:A:318:PHE:CD1	1:A:318:PHE:C	2.96	0.44
1:A:491:ASN:ND2	1:A:848:GLU:OE2	2.51	0.44
1:A:667:ILE:HD13	1:A:667:ILE:N	2.32	0.44
1:A:449:ILE:HG13	1:A:509:LEU:HG	1.99	0.44
1:A:338:PRO:HB2	1:A:713:SER:HB3	1.99	0.44
1:A:832:SER:OG	1:A:835:ASN:ND2	2.50	0.44
1:A:905:PRO:HB3	1:A:909:ASP:HB3	2.00	0.44
1:A:577:ARG:O	1:A:583:THR:OG1	2.33	0.43
1:A:1103:VAL:O	1:A:1107:ILE:HG13	2.18	0.43
1:A:212:GLU:CD	1:A:212:GLU:O	2.61	0.43
1:A:22:THR:HB	1:A:45:ARG:HB3	1.99	0.43
1:A:122:VAL:HG12	1:A:123:ASN:H	1.83	0.43
1:A:569:PHE:CD1	1:A:592:LEU:HD13	2.53	0.43
1:A:1125:ILE:HD12	1:A:1126:PHE:N	2.33	0.43
1:A:240:PRO:HD2	1:A:839:GLN:HG3	1.99	0.43
1:A:698:VAL:HA	1:A:701:ILE:HG12	2.00	0.43
1:A:821:ILE:HA	1:A:833:LEU:HB3	1.99	0.43
1:A:206:LEU:H	1:A:206:LEU:CD2	2.29	0.43
1:A:840:GLU:OE2	1:A:843:ARG:NH2	2.51	0.43
1:A:29:CYS:HB2	1:A:41:CYS:HB3	1.40	0.43
1:A:128:GLU:HG3	1:A:136:ILE:HG13	2.00	0.43
1:A:206:LEU:HG	1:A:207:ALA:H	1.84	0.43
1:A:779:ALA:HB2	1:A:1125:ILE:HG23	2.01	0.42
1:A:853:TRP:HB3	1:A:933:TYR:CD1	2.54	0.42
1:A:37:ASN:OD1	1:A:38:GLU:N	2.51	0.42
1:A:153:TYR:HE1	1:A:169:MET:HE2	1.84	0.42
1:A:986:PRO:HB2	3:A:1201:ERG:H271	2.01	0.42
1:A:206:LEU:CD1	1:A:213:TRP:HD1	2.32	0.42
1:A:317:ASN:HB3	1:A:319:PHE:CD2	2.54	0.42
1:A:668:LEU:HG	1:A:672:LEU:HD13	2.01	0.42
1:A:142:VAL:HG12	1:A:143:PHE:O	2.20	0.42
1:A:632:ASN:HD22	1:A:632:ASN:HA	1.51	0.42
1:A:940:ASN:CB	1:A:942:THR:O	2.65	0.41
1:A:161:PHE:N	1:A:166:GLY:O	2.54	0.41
1:A:399:VAL:HG22	1:A:540:ILE:HG22	2.02	0.41
1:A:756:PHE:HA	1:A:759:TRP:HB3	2.02	0.41
1:A:1059:LEU:O	1:A:1063:VAL:HG23	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:1:NAG:H4	2:C:2:NAG:H2	1.89	0.41
1:A:674:GLN:NE2	1:A:1068:GLU:HB2	2.35	0.41
1:A:937:LEU:HD12	1:A:939:TYR:HB3	2.01	0.41
1:A:798:ASN:C	1:A:955:PRO:HG3	2.45	0.41
1:A:558:ILE:HD12	1:A:561:VAL:HG21	2.03	0.41
1:A:1128:VAL:HA	1:A:1132:ARG:NH1	2.36	0.41
1:A:340:LYS:O	1:A:344:THR:HG23	2.21	0.41
1:A:845:THR:HG21	1:A:956:LEU:HD21	2.03	0.41
1:A:1030:ALA:HA	1:A:1033:VAL:HG12	2.03	0.41
1:A:109:PHE:CZ	1:A:186:LEU:HD11	2.56	0.40
1:A:955:PRO:HB2	1:A:957:ARG:CD	2.52	0.40

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	1090/1198 (91%)	1004 (92%)	84 (8%)	2 (0%)	43 74

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	943	SER
1	A	192	MET

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	981/1064 (92%)	965 (98%)	16 (2%)	55 70

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	34	VAL
1	A	63	VAL
1	A	122	VAL
1	A	133	ASP
1	A	140	LEU
1	A	206	LEU
1	A	318	PHE
1	A	505	PHE
1	A	510	LEU
1	A	529	LEU
1	A	727	ILE
1	A	804	TYR
1	A	939	TYR
1	A	943	SER
1	A	982	PHE
1	A	987	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	199	GLN
1	A	332	GLN
1	A	555	ASN
1	A	632	ASN
1	A	707	GLN
1	A	798	ASN
1	A	835	ASN
1	A	893	GLN

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 ⓘ

6 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	B	1	1,2	14,14,15	0.77	1 (7%)	17,19,21	1.04	1 (5%)
2	NAG	B	2	2	14,14,15	0.62	1 (7%)	17,19,21	0.40	0
2	BMA	B	3	2	11,11,12	0.59	0	15,15,17	1.06	1 (6%)
2	NAG	C	1	1,2	14,14,15	1.20	1 (7%)	17,19,21	1.10	1 (5%)
2	NAG	C	2	2	14,14,15	0.36	0	17,19,21	0.71	0
2	BMA	C	3	2	11,11,12	0.93	1 (9%)	15,15,17	0.93	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	B	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	B	2	2	-	0/6/23/26	0/1/1/1
2	BMA	B	3	2	-	2/2/19/22	0/1/1/1
2	NAG	C	1	1,2	-	1/6/23/26	0/1/1/1
2	NAG	C	2	2	-	2/6/23/26	0/1/1/1
2	BMA	C	3	2	-	0/2/19/22	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	1	NAG	O5-C1	-3.81	1.37	1.43
2	B	1	NAG	C1-C2	2.59	1.55	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	3	BMA	O5-C1	-2.47	1.39	1.43
2	B	2	NAG	C1-C2	2.06	1.55	1.52

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1	NAG	C3-C4-C5	3.66	116.87	110.23
2	B	1	NAG	C2-N2-C7	3.12	127.08	122.90
2	B	3	BMA	C1-O5-C5	2.61	115.68	112.19

There are no chirality outliers.

All (7) torsion outliers are listed below:

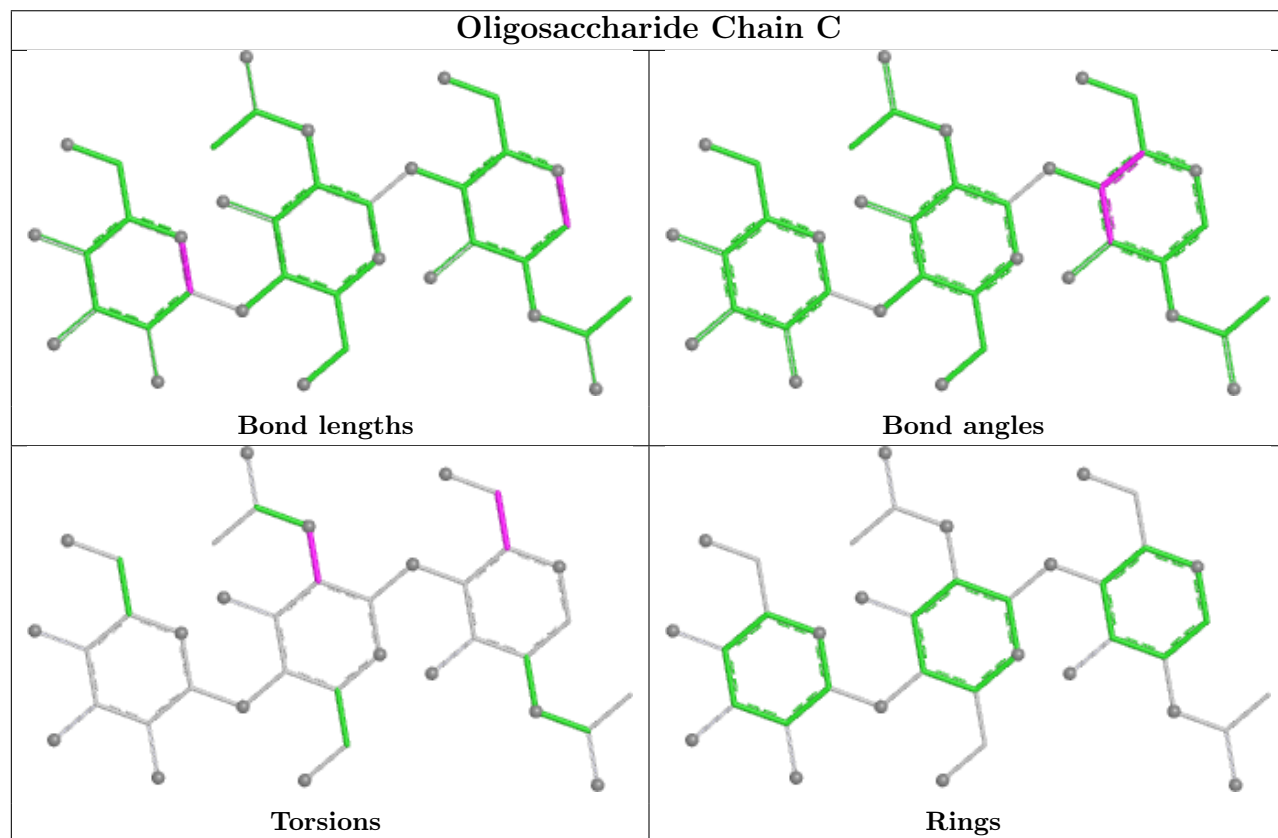
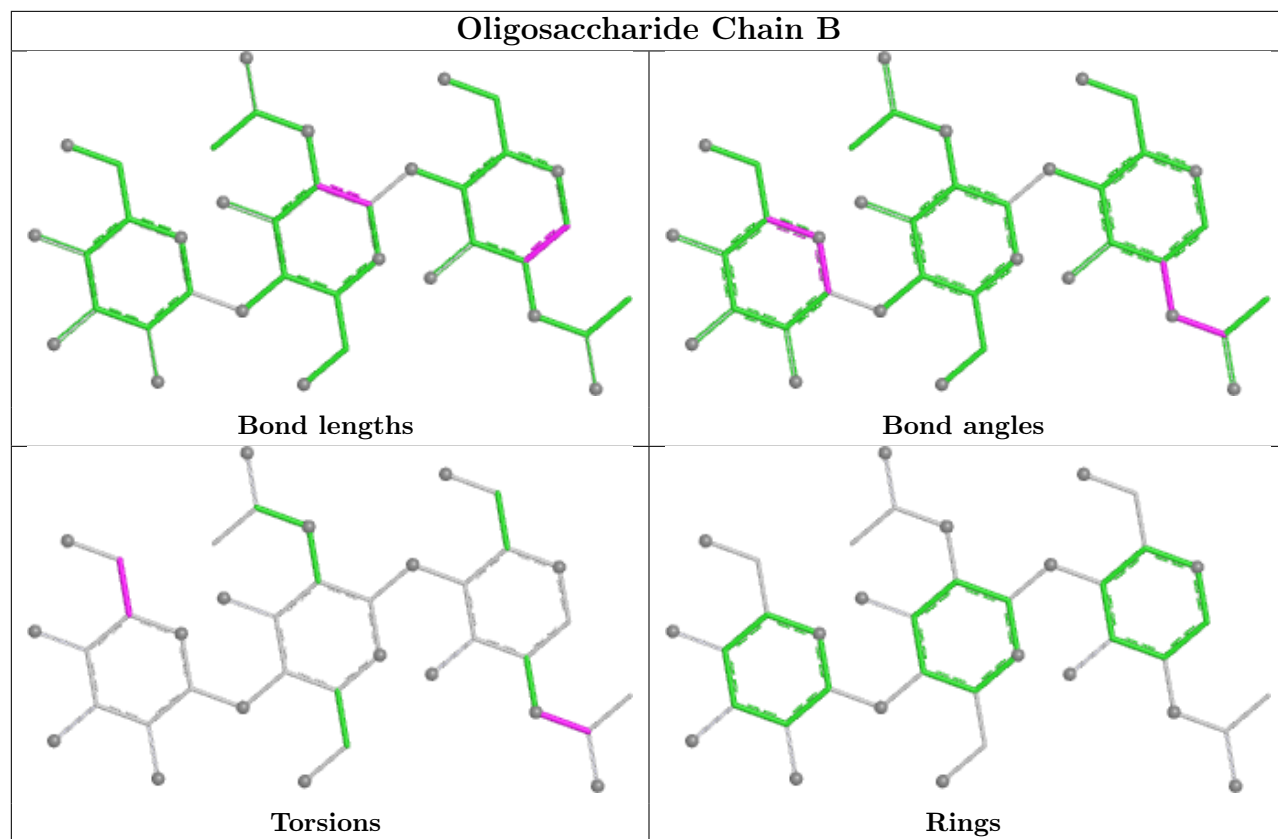
Mol	Chain	Res	Type	Atoms
2	B	1	NAG	C8-C7-N2-C2
2	B	1	NAG	O7-C7-N2-C2
2	C	1	NAG	O5-C5-C6-O6
2	B	3	BMA	C4-C5-C6-O6
2	B	3	BMA	O5-C5-C6-O6
2	C	2	NAG	C3-C2-N2-C7
2	C	2	NAG	C1-C2-N2-C7

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1	NAG	1	0
2	C	2	NAG	1	0
2	C	1	NAG	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 oligosaccharide.



5.6 Ligand geometry

3 ligands 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$
3	ERG	A	1201	-	32,32,32	0.19	0	45,50,50	0.80	2 (4%)
4	NAG	A	1208	1	14,14,15	1.28	2 (14%)	17,19,21	0.97	1 (5%)
4	NAG	A	1209	1	14,14,15	1.94	2 (14%)	17,19,21	1.21	2 (11%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	ERG	A	1201	-	-	5/13/71/71	0/4/4/4
4	NAG	A	1208	1	-	2/6/23/26	0/1/1/1
4	NAG	A	1209	1	-	2/6/23/26	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1209	NAG	O5-C1	6.17	1.54	1.43
4	A	1209	NAG	C1-C2	3.42	1.57	1.52
4	A	1208	NAG	C8-C7	-3.26	1.43	1.50
4	A	1208	NAG	O5-C1	-2.70	1.39	1.43

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1208	NAG	C1-C2-N2	-3.44	105.01	110.43
4	A	1209	NAG	C1-O5-C5	2.81	115.95	112.19
4	A	1209	NAG	O4-C4-C3	-2.41	104.70	110.38
3	A	1201	ERG	C9-C10-C5	2.27	113.28	109.66
3	A	1201	ERG	C14-C13-C17	-2.25	97.37	99.72

There are no chirality outliers.

All (9) torsion outliers are listed below:

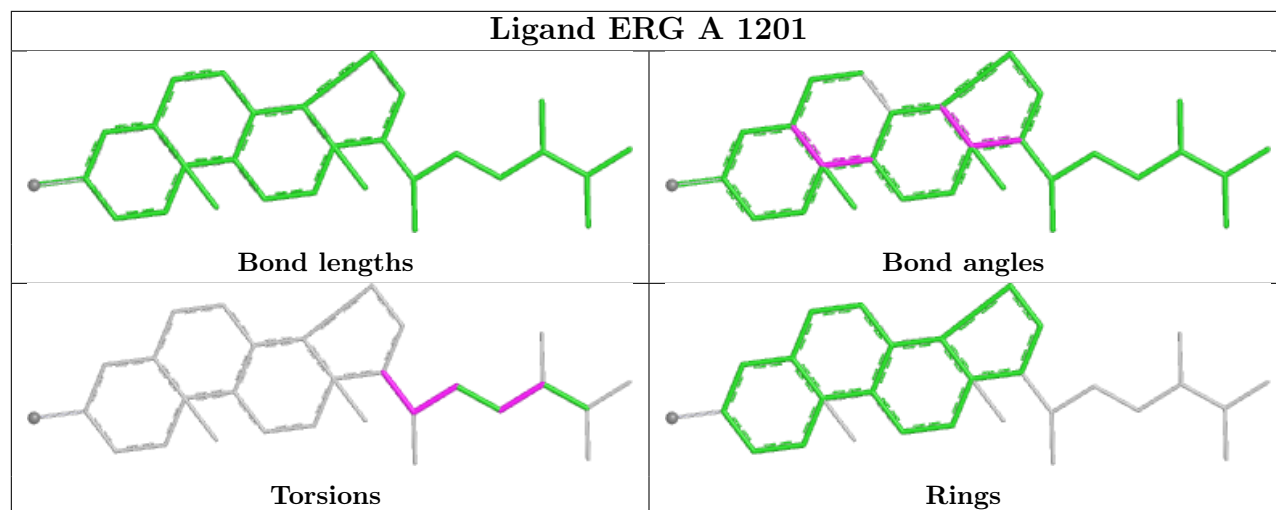
Mol	Chain	Res	Type	Atoms
4	A	1209	NAG	C8-C7-N2-C2
4	A	1209	NAG	O7-C7-N2-C2
3	A	1201	ERG	C17-C20-C22-C23
3	A	1201	ERG	C22-C23-C24-C25
3	A	1201	ERG	C16-C17-C20-C22
3	A	1201	ERG	C21-C20-C22-C23
4	A	1208	NAG	C3-C2-N2-C7
4	A	1208	NAG	C1-C2-N2-C7
3	A	1201	ERG	C22-C23-C24-C28

There are no ring outliers.

2 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1201	ERG	5	0
4	A	1208	NAG	4	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	1096/1198 (91%)	0.01	14 (1%)	75 50	74, 156, 215, 362	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	805	MET	3.7
1	A	779	ALA	3.3
1	A	937	LEU	3.0
1	A	950	ARG	3.0
1	A	663	MET	2.8
1	A	210	GLU	2.6
1	A	900	ASN	2.6
1	A	21	ALA	2.6
1	A	121	PHE	2.6
1	A	951	THR	2.4
1	A	371	VAL	2.2
1	A	97	SER	2.2
1	A	657	ILE	2.0
1	A	419	GLU	2.0

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

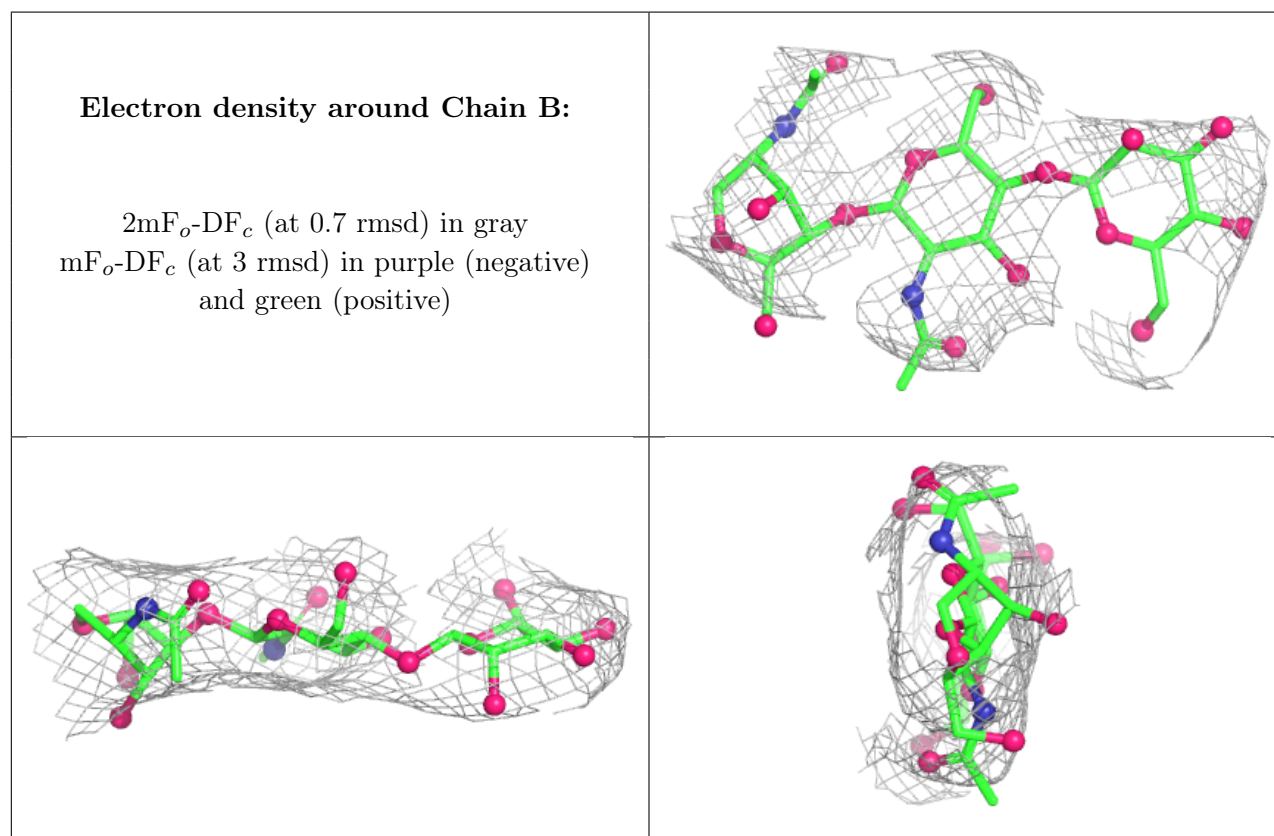
There are no non-standard protein/DNA/RNA residues in this entry.

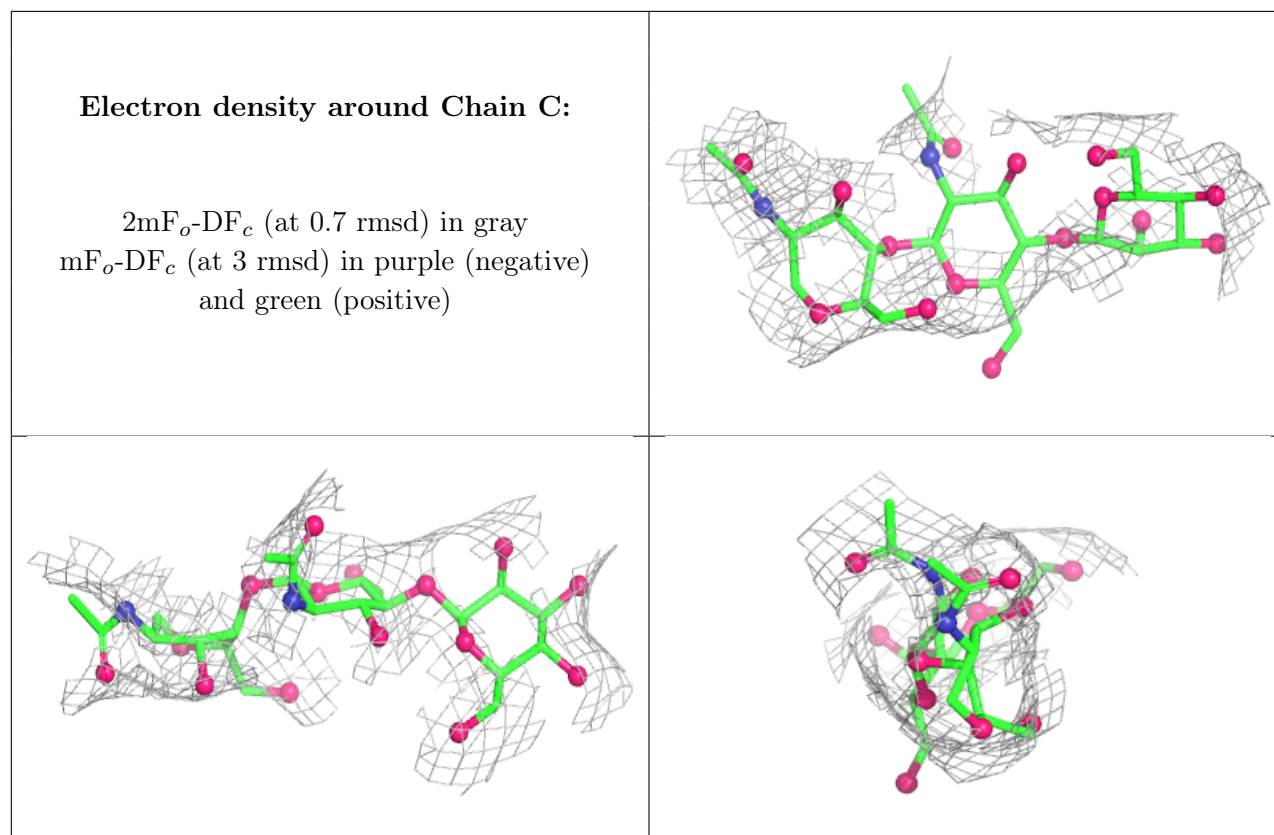
6.3 Carbohydrates [i](#)

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($\text{\AA}^2$)	Q<0.9
2	NAG	B	2	14/15	0.45	0.10	191,225,239,249	0
2	BMA	B	3	11/12	0.46	0.09	184,224,236,239	0
2	BMA	C	3	11/12	0.64	0.09	181,233,243,254	0
2	NAG	C	2	14/15	0.66	0.08	207,254,264,266	0
2	NAG	B	1	14/15	0.67	0.10	194,212,232,243	0
2	NAG	C	1	14/15	0.73	0.08	240,256,264,268	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.



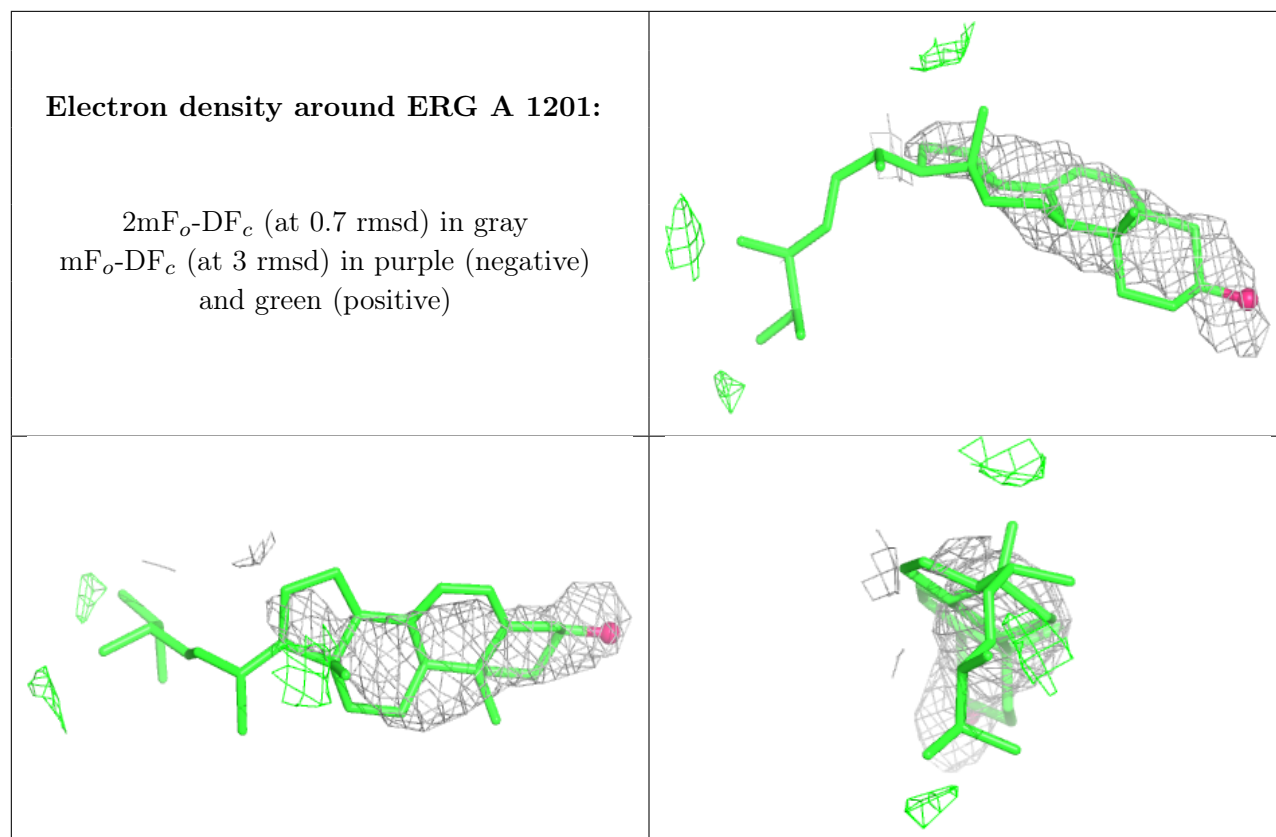


6.4 Ligands [i](#)

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	NAG	A	1208	14/15	0.60	0.09	210,237,265,275	0
4	NAG	A	1209	14/15	0.65	0.12	186,207,244,247	0
3	ERG	A	1201	29/29	0.91	0.25	142,156,172,175	29

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.



6.5 Other polymers [i](#)

There are no such residues in this entry.