



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 05:52 PM UTC

PDB ID : 3NWD / pdb_00003nwd
Title : Glycoprotein B from Herpes simplex virus type 1, Y179S mutant, low-pH
Authors : Stampfer, S.D.; Lou, H.; Cohen, G.H.; Eisenberg, R.J.; Heldwein, E.E.
Deposited on : 2010-07-09
Resolution : 2.88 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
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

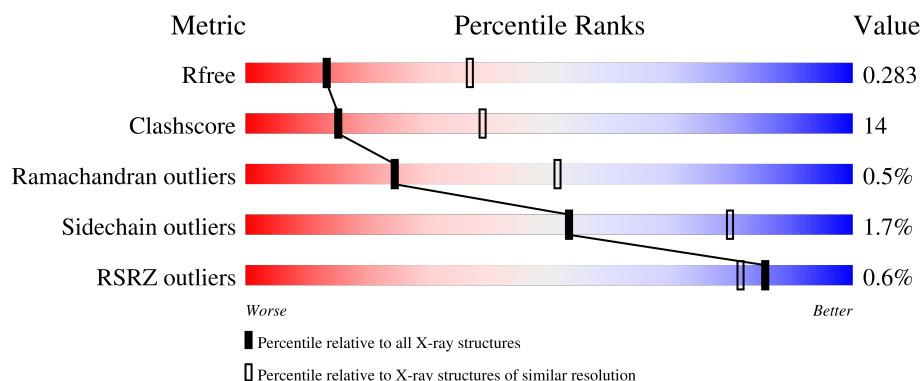
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.88 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	3557 (2.90-2.86)
Clashscore	190562	3801 (2.90-2.86)
Ramachandran outliers	187476	3699 (2.90-2.86)
Sidechain outliers	187428	3702 (2.90-2.86)
RSRZ outliers	180081	3558 (2.90-2.86)

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	703	
1	B	703	
1	C	703	
1	D	703	
2	E	2	

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Mol	Chain	Length	Quality of chain
2	F	2	 100%

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	E	1	X	-	-	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 19481 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 Envelope glycoprotein B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	597	Total	C	N	O	S	0	1	0
			4767	3000	835	910	22			
1	A	606	Total	C	N	O	S	0	1	0
			4843	3056	843	922	22			
1	C	601	Total	C	N	O	S	0	0	0
			4766	3003	831	910	22			
1	D	606	Total	C	N	O	S	0	1	0
			4829	3044	846	918	21			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	28	ASP	-	expression tag	UNP P06437
B	29	PRO	-	expression tag	UNP P06437
B	58	ALA	PRO	SEE REMARK 999	UNP P06437
B	179	SER	TYR	engineered mutation	UNP P06437
B	313	SER	THR	SEE REMARK 999	UNP P06437
B	443	LEU	GLN	SEE REMARK 999	UNP P06437
A	28	ASP	-	expression tag	UNP P06437
A	29	PRO	-	expression tag	UNP P06437
A	58	ALA	PRO	SEE REMARK 999	UNP P06437
A	179	SER	TYR	engineered mutation	UNP P06437
A	313	SER	THR	SEE REMARK 999	UNP P06437
A	443	LEU	GLN	SEE REMARK 999	UNP P06437
C	28	ASP	-	expression tag	UNP P06437
C	29	PRO	-	expression tag	UNP P06437
C	58	ALA	PRO	SEE REMARK 999	UNP P06437
C	179	SER	TYR	engineered mutation	UNP P06437
C	313	SER	THR	SEE REMARK 999	UNP P06437
C	443	LEU	GLN	SEE REMARK 999	UNP P06437
D	28	ASP	-	expression tag	UNP P06437
D	29	PRO	-	expression tag	UNP P06437
D	58	ALA	PRO	SEE REMARK 999	UNP P06437

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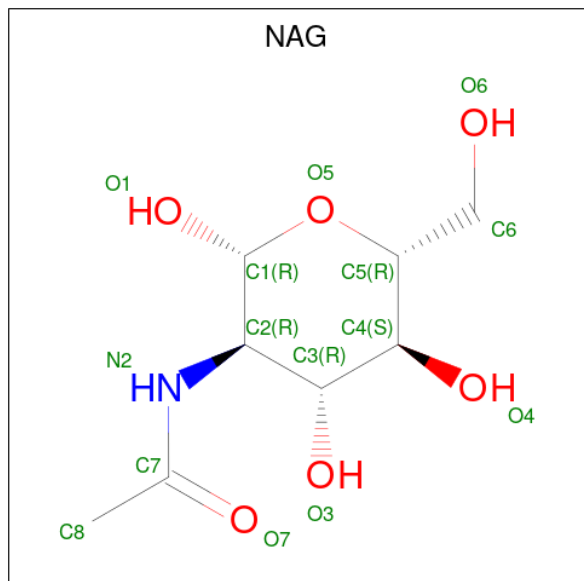
Chain	Residue	Modelled	Actual	Comment	Reference
D	179	SER	TYR	engineered mutation	UNP P06437
D	313	SER	THR	SEE REMARK 999	UNP P06437
D	443	LEU	GLN	SEE REMARK 999	UNP P06437

- Molecule 2 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
2	E	2	Total	C	N	O	0	0	0
			28	16	2	10			
2	F	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



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

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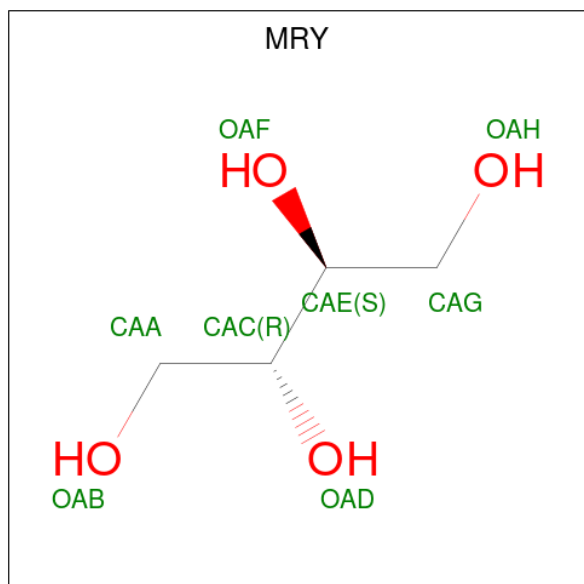
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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	C	1	Total	C	N	O	0	0
			14	8	1	5		
3	C	1	Total	C	N	O	0	0
			14	8	1	5		
3	C	1	Total	C	N	O	0	0
	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

- Molecule 4 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Cl	0	0
			1	1		
4	D	1	Total	Cl	0	0
			1	1		

- Molecule 5 is MESO-ERYTHRITOL (CCD ID: MRY) (formula: C₄H₁₀O₄).



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

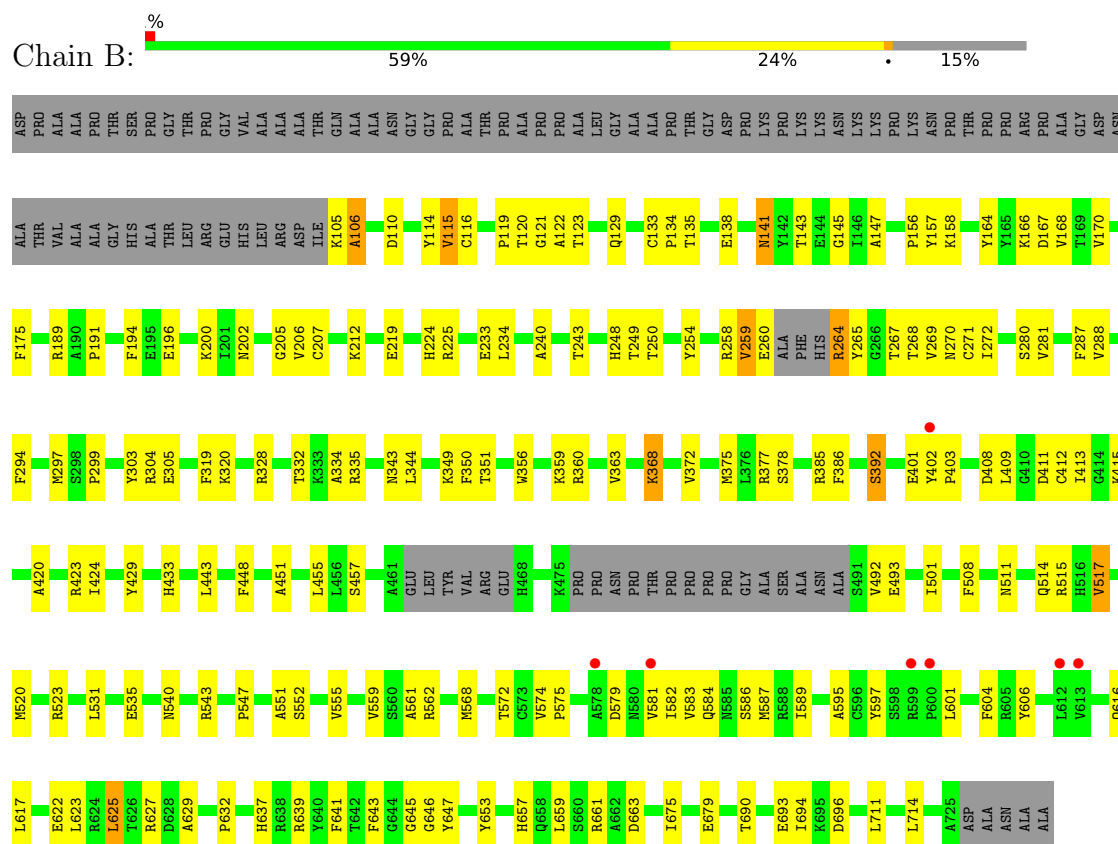
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	16	Total	O	0	0
			16	16		
6	A	25	Total	O	0	0
			25	25		
6	C	18	Total	O	0	0
			18	18		
6	D	25	Total	O	0	0
			25	25		

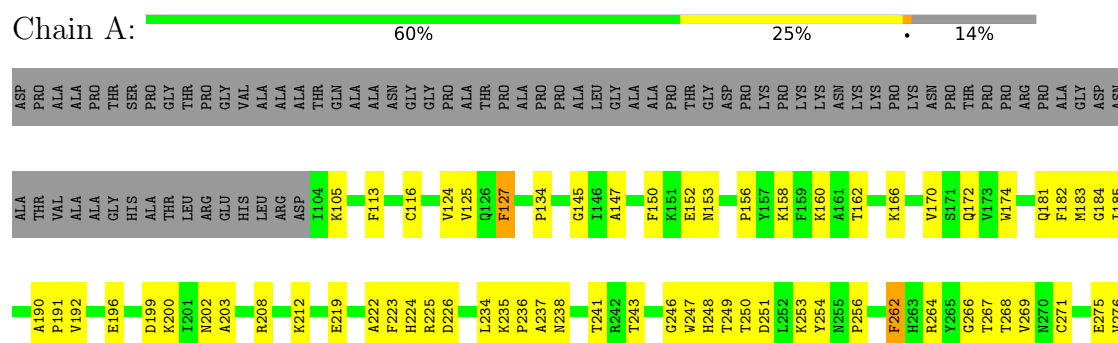
3 Residue-property plots [i](#)

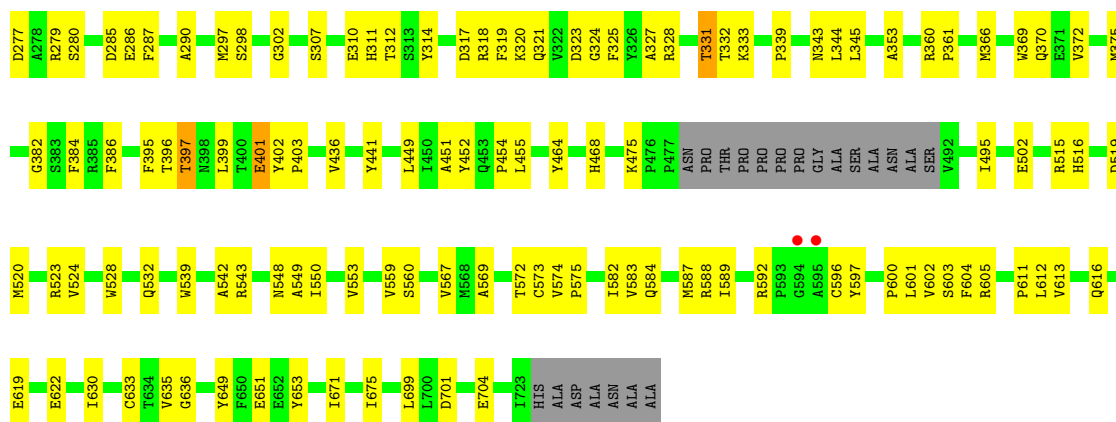
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density.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



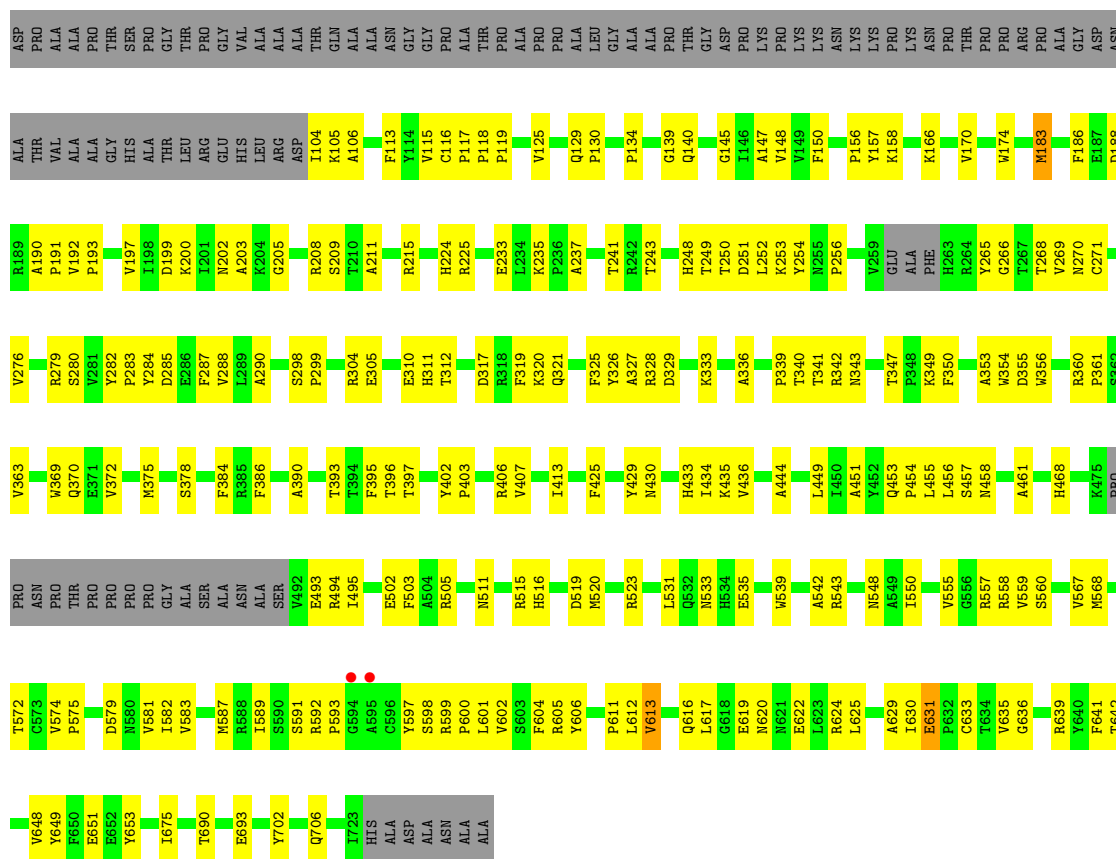
• Molecule 1: Envelope glycoprotein B





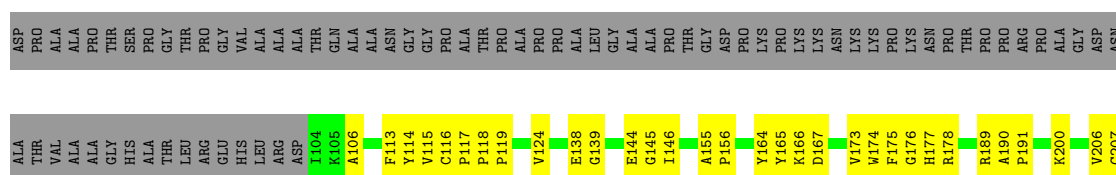
• Molecule 1: Envelope glycoprotein B

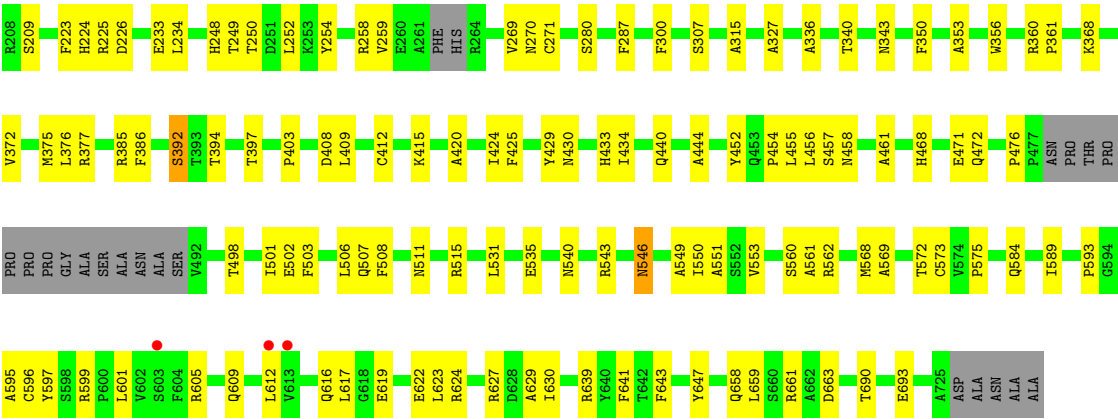
Chain C: 55% 30% 15%



• Molecule 1: Envelope glycoprotein B

Chain D: 64% 22% 14%





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



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



4 Data and refinement statistics

Property	Value	Source
Space group	P 3	Depositor
Cell constants a, b, c, α , β , γ	117.76Å 117.76Å 318.56Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.56 – 2.88 48.56 – 2.88	Depositor EDS
% Data completeness (in resolution range)	88.0 (48.56-2.88) 88.0 (48.56-2.88)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.32 (at 2.86Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6.1_357)	Depositor
R, R_{free}	0.238 , 0.283 0.234 , 0.283	Depositor DCC
R_{free} test set	5310 reflections (4.77%)	wwPDB-VP
Wilson B-factor (Å ²)	64.3	Xtriage
Anisotropy	0.014	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 66.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.020 for -h,-k,l 0.468 for h,-h-k,-l 0.022 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	19481	wwPDB-VP
Average B, all atoms (Å ²)	86.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 73.43 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.8395e-06. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical 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: MRY, NAG, CL

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	0.26	0/4966	0.70	2/6757 (0.0%)
1	B	0.25	0/4879	0.68	0/6631
1	C	0.25	0/4880	0.68	0/6639
1	D	0.25	0/4948	0.69	0/6731
All	All	0.25	0/19673	0.69	2/26758 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	475	LYS	CA-C-N	5.09	123.38	119.66
1	A	475	LYS	C-N-CA	5.09	123.38	119.66

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	4843	0	4604	121	0
1	B	4767	0	4540	130	0
1	C	4766	0	4514	148	0
1	D	4829	0	4602	115	0
2	E	28	0	25	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	F	28	0	25	4	0
3	A	42	0	39	0	0
3	B	28	0	26	0	0
3	C	42	0	39	0	0
3	D	14	0	13	0	0
4	A	1	0	0	0	0
4	D	1	0	0	0	0
5	C	8	0	10	1	0
6	A	25	0	0	0	0
6	B	16	0	0	0	0
6	C	18	0	0	1	0
6	D	25	0	0	0	0
All	All	19481	0	18437	515	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 14.

All (515) 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:183:MET:H	1:A:184:GLY:HA3	1.20	1.05
1:A:183:MET:N	1:A:184:GLY:HA3	1.88	0.88
1:B:259:VAL:HG22	1:B:260:GLU:H	1.45	0.82
1:C:256:PRO:HD3	1:C:266:GLY:HA2	1.64	0.79
1:D:403:PRO:HG3	1:D:476:PRO:HB3	1.65	0.79
1:C:393:THR:HG22	1:C:505:ARG:HG2	1.66	0.77
1:A:150:PHE:HB2	1:A:449:LEU:HB3	1.68	0.75
1:C:599:ARG:HH12	1:C:617:LEU:HG	1.53	0.73
1:A:436:VAL:HB	1:A:454:PRO:HB2	1.70	0.71
1:B:166:LYS:HG2	1:B:271:CYS:HA	1.72	0.71
1:C:174:TRP:HB2	1:C:183:MET:SD	2.30	0.70
1:C:397:THR:HG22	1:C:444:ALA:HA	1.73	0.69
1:C:116:CYS:HB3	1:C:560:SER:HB2	1.74	0.69
1:D:375:MET:SD	1:D:386:PHE:HB3	2.33	0.68
1:A:298:SER:HB3	1:A:310:GLU:HB3	1.76	0.67
1:B:385:ARG:HH22	1:B:515:ARG:NH2	1.93	0.67
1:C:156:PRO:HG2	1:C:279:ARG:NH2	2.10	0.67
1:D:456:LEU:HD21	1:D:461:ALA:HA	1.75	0.67
1:D:224:HIS:HB2	1:D:269:VAL:HB	1.77	0.66
1:A:156:PRO:HG2	1:A:279:ARG:NH2	2.11	0.66
1:A:174:TRP:HB2	1:A:183:MET:SD	2.36	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:546:ASN:HB3	1:D:549:ALA:HB3	1.78	0.66
1:C:202:ASN:O	1:C:328:ARG:HB3	1.95	0.65
1:D:584:GLN:HG3	1:D:601:LEU:HB2	1.79	0.65
1:C:158:LYS:HD3	1:C:279:ARG:NH1	2.12	0.65
1:D:166:LYS:HG2	1:D:271:CYS:HA	1.77	0.65
1:C:606:TYR:CD2	1:C:613:VAL:HG21	2.31	0.64
1:C:436:VAL:HB	1:C:454:PRO:HB2	1.79	0.64
1:D:166:LYS:HE3	1:D:207:CYS:SG	2.37	0.64
1:C:156:PRO:HG2	1:C:279:ARG:HH21	1.62	0.64
1:D:119:PRO:HG2	1:D:562:ARG:HG2	1.80	0.64
1:D:206:VAL:HG12	1:D:233:GLU:HG2	1.79	0.64
1:D:280:SER:HB2	1:D:287:PHE:HB3	1.80	0.63
1:A:105:LYS:HG2	1:A:582:ILE:HG22	1.81	0.63
1:C:166:LYS:HG2	1:C:271:CYS:HA	1.80	0.63
1:A:202:ASN:O	1:A:328:ARG:HB3	1.98	0.63
1:B:156:PRO:HB3	1:B:281:VAL:HG12	1.81	0.62
1:D:550:ILE:O	1:D:553:VAL:HG12	1.99	0.62
1:C:145:GLY:HA2	1:C:455:LEU:HG	1.81	0.62
1:C:583:VAL:HA	1:C:602:VAL:HG12	1.80	0.62
1:A:116:CYS:HB3	1:A:560:SER:HB2	1.82	0.61
1:B:224:HIS:HB2	1:B:269:VAL:HB	1.81	0.61
1:A:235:LYS:HG2	1:A:248:HIS:O	2.00	0.61
1:A:464:TYR:O	1:A:468:HIS:HB2	2.00	0.61
1:D:223:PHE:HB2	1:D:226:ASP:HA	1.83	0.60
1:B:189:ARG:HB2	1:B:349:LYS:HE2	1.83	0.60
1:A:325:PHE:HE2	1:A:327:ALA:HB2	1.66	0.60
1:D:119:PRO:HG3	1:D:561:ALA:HA	1.82	0.60
1:C:386:PHE:HB2	1:C:395:PHE:HB2	1.83	0.60
1:C:601:LEU:HD23	1:C:616:GLN:HB3	1.83	0.60
1:B:105:LYS:N	1:B:582:ILE:HG13	2.16	0.60
1:A:237:ALA:HA	1:A:248:HIS:CD2	2.37	0.60
1:D:377:ARG:HD3	1:D:386:PHE:CZ	2.38	0.59
1:D:616:GLN:HG3	1:D:624:ARG:HB2	1.83	0.59
1:A:235:LYS:HE3	1:A:248:HIS:CE1	2.37	0.59
1:C:649:TYR:CE2	1:C:651:GLU:HB2	2.37	0.59
1:D:360:ARG:HG2	1:D:409:LEU:HD23	1.84	0.58
1:B:343:ASN:HD21	1:B:356:TRP:HD1	1.51	0.58
1:C:235:LYS:HE3	1:C:248:HIS:CE1	2.38	0.58
1:D:252:LEU:H	1:D:252:LEU:HD23	1.68	0.58
1:B:280:SER:HB2	1:B:287:PHE:HB3	1.84	0.58
1:A:324:GLY:HA2	1:A:339:PRO:HB2	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:115:VAL:O	1:D:117:PRO:HD3	2.02	0.58
1:B:175:PHE:CD2	1:B:259:VAL:HG12	2.38	0.58
1:A:589:ILE:HG22	1:A:592:ARG:H	1.69	0.58
1:C:166:LYS:HD2	1:C:211:ALA:HB2	1.86	0.58
1:B:531:LEU:O	1:B:535:GLU:HG2	2.04	0.58
1:A:253:LYS:HA	1:A:268:THR:HG21	1.86	0.58
1:D:377:ARG:HB2	1:D:386:PHE:CE2	2.39	0.58
1:B:690:THR:OG1	1:B:693:GLU:HG3	2.04	0.57
1:A:314:TYR:HB3	1:A:318:ARG:HD2	1.84	0.57
1:C:511:ASN:O	1:C:515:ARG:HG2	2.04	0.57
1:A:600:PRO:O	1:A:616:GLN:HG3	2.04	0.57
1:B:170:VAL:HG22	1:B:267:THR:HG22	1.86	0.57
1:B:119:PRO:HG2	1:B:562:ARG:HG2	1.87	0.57
1:A:603:SER:HB3	1:A:612:LEU:HG	1.87	0.57
1:D:429:TYR:CE2	1:D:455:LEU:HD13	2.39	0.57
1:B:240:ALA:HB1	1:B:243:THR:OG1	2.05	0.57
1:D:248:HIS:HA	1:D:271:CYS:O	2.05	0.56
1:A:401:GLU:HG2	1:A:441:TYR:O	2.06	0.56
1:A:582:ILE:HD11	1:A:603:SER:HB2	1.87	0.56
1:A:588:ARG:HG2	1:A:596:CYS:SG	2.46	0.56
1:C:118:PRO:HD3	1:C:625:LEU:CD2	2.35	0.56
1:A:331:THR:HG22	1:A:332:THR:HG23	1.87	0.56
1:B:141:ASN:N	1:B:141:ASN:HD22	2.02	0.56
1:C:113:PHE:O	1:C:575:PRO:HA	2.04	0.56
1:D:343:ASN:HD21	1:D:356:TRP:HD1	1.54	0.56
1:A:325:PHE:CE2	1:A:327:ALA:HB2	2.41	0.56
1:D:139:GLY:C	2:F:1:NAG:H82	2.31	0.56
1:A:147:ALA:HA	1:A:451:ALA:O	2.06	0.56
1:B:616:GLN:CD	1:B:627:ARG:HA	2.31	0.56
1:B:343:ASN:CG	1:B:356:TRP:HB2	2.31	0.55
1:C:250:THR:HG22	1:C:252:LEU:H	1.72	0.55
1:A:539:TRP:O	1:A:543:ARG:HG3	2.06	0.55
1:B:205:GLY:O	1:B:233:GLU:HG3	2.06	0.55
1:B:540:ASN:O	1:B:543:ARG:HG2	2.07	0.55
1:B:625:LEU:H	1:B:625:LEU:HD22	1.72	0.55
1:B:595:ALA:HA	1:B:632:PRO:HA	1.89	0.55
1:C:241:THR:O	1:C:243:THR:HG23	2.07	0.55
1:A:597:TYR:HA	1:A:630:ILE:HA	1.89	0.55
1:A:649:TYR:CE2	1:A:651:GLU:HB2	2.41	0.55
1:A:366:MET:HE2	1:A:495:ILE:HB	1.88	0.54
1:C:516:HIS:O	1:C:520:MET:HG2	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:372:VAL:HG11	1:C:375:MET:HE2	1.89	0.54
1:C:543:ARG:HB2	1:C:568:MET:HE1	1.88	0.54
1:B:167:ASP:OD2	1:B:189:ARG:HG2	2.07	0.54
1:B:234:LEU:HD23	1:B:249:THR:HG23	1.89	0.54
1:A:223:PHE:HB2	1:A:226:ASP:HA	1.89	0.54
1:D:498:THR:HG21	1:D:503:PHE:HE2	1.73	0.54
1:B:372:VAL:HG11	1:B:375:MET:HE2	1.88	0.54
1:C:200:LYS:HE3	1:C:208:ARG:CZ	2.37	0.54
1:A:360:ARG:HB2	1:A:361:PRO:HD3	1.89	0.54
1:C:633:CYS:HA	1:C:653:TYR:OH	2.08	0.54
1:B:191:PRO:HA	1:B:350:PHE:HA	1.88	0.54
1:B:196:GLU:O	1:B:200:LYS:HB2	2.07	0.54
1:C:150:PHE:HB2	1:C:449:LEU:HB3	1.90	0.54
1:B:711:LEU:HD12	1:B:714:LEU:HD12	1.89	0.53
1:C:597:TYR:CE1	1:C:601:LEU:HD11	2.43	0.53
1:D:343:ASN:HD21	1:D:356:TRP:CD1	2.26	0.53
1:A:285:ASP:HA	1:A:298:SER:HB2	1.89	0.53
1:C:105:LYS:HE2	1:C:579:ASP:O	2.08	0.53
1:B:520:MET:HE2	1:B:520:MET:HA	1.91	0.53
1:B:551:ALA:O	1:B:555:VAL:HG23	2.09	0.53
1:D:225:ARG:HD2	1:D:254:TYR:CD1	2.43	0.53
1:B:559:VAL:HG12	1:B:572:THR:HA	1.91	0.53
1:B:647:TYR:HE2	1:B:661:ARG:HE	1.55	0.53
1:D:502:GLU:O	1:D:506:LEU:HB2	2.08	0.53
1:D:531:LEU:O	1:D:535:GLU:HG2	2.08	0.53
1:B:551:ALA:HB2	1:B:568:MET:SD	2.49	0.53
1:D:209:SER:HB2	1:D:224:HIS:HB3	1.91	0.53
1:B:250:THR:HB	1:B:270:ASN:HB2	1.90	0.52
1:B:639:ARG:HB3	1:B:641:PHE:CE1	2.44	0.52
1:D:124:VAL:HG12	1:D:569:ALA:HA	1.90	0.52
1:D:206:VAL:HA	1:D:233:GLU:HA	1.90	0.52
1:A:515:ARG:HD2	1:A:515:ARG:N	2.23	0.52
1:B:375:MET:SD	1:B:386:PHE:HB3	2.49	0.52
1:A:145:GLY:HA2	1:A:455:LEU:HG	1.92	0.52
1:A:212:LYS:HE2	1:A:219:GLU:OE2	2.10	0.52
1:A:280:SER:HB2	1:A:287:PHE:HB3	1.91	0.52
1:C:317:ASP:O	1:C:320:LYS:HE2	2.09	0.52
1:A:225:ARG:HD2	1:A:254:TYR:CD1	2.44	0.52
1:A:516:HIS:O	1:A:520:MET:HG2	2.10	0.52
1:B:206:VAL:HA	1:B:233:GLU:HA	1.91	0.52
1:C:225:ARG:HD2	1:C:254:TYR:CD1	2.45	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:343:ASN:O	1:A:353:ALA:HA	2.09	0.52
1:C:360:ARG:HB2	1:C:361:PRO:HD3	1.92	0.52
1:C:620:ASN:HA	1:C:642:THR:HG21	1.92	0.52
1:D:601:LEU:HB3	1:D:627:ARG:CZ	2.40	0.52
1:D:601:LEU:HD13	1:D:627:ARG:NH1	2.25	0.52
1:D:639:ARG:HB3	1:D:641:PHE:CE1	2.44	0.52
1:B:429:TYR:CD2	1:B:455:LEU:HD13	2.45	0.51
1:D:178:ARG:O	1:D:178:ARG:HG2	2.11	0.51
1:D:605:ARG:HB3	1:D:612:LEU:HD23	1.92	0.51
1:B:429:TYR:CE2	1:B:455:LEU:HD13	2.45	0.51
1:C:433:HIS:HD2	1:C:457:SER:HA	1.75	0.51
1:D:165:TYR:CE1	1:D:189:ARG:HD2	2.45	0.51
1:A:158:LYS:HD3	1:A:279:ARG:NH1	2.26	0.51
1:D:138:GLU:CB	2:F:1:NAG:H83	2.41	0.51
1:A:328:ARG:HH21	1:A:333:LYS:HG2	1.76	0.50
1:A:597:TYR:CE1	1:A:601:LEU:HD11	2.46	0.50
1:B:420:ALA:O	1:B:423:ARG:HB3	2.10	0.50
1:A:319:PHE:CZ	1:A:321:GLN:HB2	2.47	0.50
1:D:540:ASN:O	1:D:543:ARG:HG2	2.12	0.50
1:C:325:PHE:CD2	1:C:342:ARG:HB2	2.47	0.50
2:E:1:NAG:H3	2:E:1:NAG:O7	2.12	0.50
1:B:584:GLN:HG2	1:B:586:SER:H	1.76	0.50
1:A:605:ARG:HB3	1:A:611:PRO:O	2.12	0.50
1:D:616:GLN:CD	1:D:627:ARG:HA	2.36	0.50
1:A:319:PHE:O	1:A:320:LYS:HD3	2.11	0.50
1:C:117:PRO:HA	1:C:625:LEU:HD21	1.92	0.50
1:A:633:CYS:HA	1:A:653:TYR:OH	2.12	0.50
1:C:125:VAL:O	1:C:567:VAL:HG13	2.11	0.50
1:B:420:ALA:O	1:B:424:ILE:HG13	2.12	0.50
1:A:241:THR:O	1:A:243:THR:HG23	2.12	0.50
1:A:343:ASN:C	1:A:344:LEU:HD12	2.37	0.50
1:D:468:HIS:O	1:D:472:GLN:HG3	2.12	0.50
1:D:597:TYR:CZ	1:D:601:LEU:HD11	2.47	0.50
1:B:412:CYS:HA	1:B:415:LYS:HD2	1.94	0.49
1:B:552:SER:HA	1:B:559:VAL:HG22	1.93	0.49
1:D:343:ASN:ND2	1:D:356:TRP:HB2	2.28	0.49
1:B:377:ARG:HA	1:B:385:ARG:O	2.12	0.49
1:A:262:PHE:CD1	1:A:262:PHE:N	2.78	0.49
1:B:287:PHE:CD2	1:B:299:PRO:HG3	2.46	0.49
1:B:616:GLN:HB3	1:B:627:ARG:HG2	1.93	0.49
1:C:639:ARG:HB3	1:C:641:PHE:CE1	2.48	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:177:HIS:O	1:D:178:ARG:HB3	2.12	0.49
1:B:212:LYS:HE2	1:B:219:GLU:CD	2.38	0.49
1:C:702:TYR:O	1:C:706:GLN:HG2	2.13	0.49
1:A:183:MET:N	1:A:184:GLY:CA	2.71	0.49
1:C:567:VAL:HG12	1:C:568:MET:N	2.27	0.49
1:C:604:PHE:CZ	1:C:613:VAL:HB	2.47	0.49
1:D:425:PHE:CE2	1:D:430:ASN:HA	2.47	0.49
1:B:501:ILE:HG13	1:B:501:ILE:O	2.12	0.49
1:B:114:TYR:HA	1:B:574:VAL:O	2.13	0.49
1:B:212:LYS:HE2	1:B:219:GLU:OE2	2.13	0.49
1:B:617:LEU:HD12	1:B:622:GLU:O	2.11	0.49
1:A:372:VAL:HG11	1:A:375:MET:HE2	1.95	0.49
1:C:166:LYS:HZ1	1:C:192:VAL:HG22	1.77	0.49
1:D:501:ILE:HG13	1:D:501:ILE:O	2.13	0.49
1:B:511:ASN:O	1:B:515:ARG:HG2	2.13	0.48
1:A:285:ASP:HB2	1:A:311:HIS:HB3	1.93	0.48
1:C:287:PHE:CD2	1:C:299:PRO:HG3	2.48	0.48
1:B:265[B]:TYR:CD1	1:B:265[B]:TYR:C	2.91	0.48
1:A:311:HIS:CG	1:A:312:THR:N	2.81	0.48
1:C:304:ARG:HB3	1:C:305:GLU:OE1	2.12	0.48
1:C:494:ARG:HG2	1:C:495:ILE:N	2.27	0.48
1:B:116:CYS:HB2	1:B:622:GLU:CD	2.39	0.48
1:B:543:ARG:O	1:B:547:PRO:HG3	2.13	0.48
1:C:597:TYR:CE2	1:C:601:LEU:HD21	2.49	0.48
1:B:106:ALA:HA	1:B:643:PHE:CE2	2.49	0.48
1:A:386:PHE:HB2	1:A:395:PHE:HB2	1.95	0.48
1:D:119:PRO:HG3	1:D:561:ALA:CA	2.43	0.48
1:C:531:LEU:O	1:C:535:GLU:HG2	2.13	0.48
1:C:253:LYS:HA	1:C:268:THR:HG21	1.95	0.48
1:D:412:CYS:HA	1:D:415:LYS:NZ	2.29	0.48
1:D:551:ALA:HB2	1:D:568:MET:SD	2.53	0.48
1:B:305:GLU:O	1:B:305:GLU:HG2	2.14	0.48
1:A:286:GLU:HA	1:A:297:MET:O	2.13	0.48
1:D:377:ARG:HH21	1:D:454:PRO:HG3	1.78	0.48
1:D:507:GLN:HG3	1:D:511:ASN:OD1	2.14	0.48
1:D:647:TYR:CE2	1:D:661:ARG:HG2	2.49	0.48
1:B:166:LYS:HE3	1:B:207:CYS:SG	2.54	0.48
1:D:300:PHE:O	1:D:307:SER:HB2	2.14	0.48
1:D:377:ARG:HA	1:D:385:ARG:O	2.13	0.48
1:A:172:GLN:HB3	1:A:184:GLY:HA2	1.95	0.48
1:A:234:LEU:HD23	1:A:249:THR:HG23	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:105:LYS:HG3	1:C:582:ILE:HG22	1.95	0.48
1:D:599:ARG:NH2	1:D:619:GLU:HG2	2.29	0.48
1:B:304:ARG:HG2	1:A:320:LYS:NZ	2.29	0.47
1:A:199:ASP:O	1:A:203:ALA:HB3	2.14	0.47
1:C:205:GLY:HA3	1:C:328:ARG:NH1	2.29	0.47
1:C:559:VAL:HG12	1:C:572:THR:HA	1.97	0.47
1:A:619:GLU:O	1:A:622:GLU:HB2	2.15	0.47
1:C:311:HIS:CG	1:C:312:THR:N	2.83	0.47
1:C:435:LYS:HD2	1:C:453:GLN:CD	2.40	0.47
1:A:559:VAL:HA	1:A:573:CYS:H	1.80	0.47
1:D:434:ILE:HG23	1:D:458:ASN:HD21	1.80	0.47
1:D:596:CYS:O	1:D:630:ILE:HG23	2.13	0.47
1:B:360:ARG:HE	1:B:409:LEU:HD23	1.79	0.47
1:B:433:HIS:HA	1:B:457:SER:HA	1.97	0.47
1:C:158:LYS:HD3	1:C:279:ARG:HH12	1.77	0.47
1:C:616:GLN:HG3	1:C:624:ARG:HD2	1.96	0.47
1:D:106:ALA:HA	1:D:658:GLN:OE1	2.13	0.47
1:D:360:ARG:N	1:D:361:PRO:HD2	2.30	0.47
1:B:115:VAL:HB	1:B:623:LEU:HB2	1.96	0.47
1:A:172:GLN:HG2	1:A:183:MET:CE	2.45	0.47
1:A:519:ASP:O	1:A:523:ARG:HG3	2.14	0.47
1:B:259:VAL:HG13	1:B:260:GLU:N	2.30	0.47
1:B:264:ARG:HG3	1:B:265[A]:TYR:H	1.79	0.47
1:C:629:ALA:C	1:C:630:ILE:HD12	2.39	0.47
1:B:105:LYS:HE3	1:B:579:ASP:O	2.15	0.47
1:B:119:PRO:HG3	1:B:561:ALA:CA	2.44	0.47
1:A:583:VAL:HA	1:A:602:VAL:HG12	1.97	0.47
1:B:115:VAL:HG23	1:B:623:LEU:O	2.15	0.47
1:B:343:ASN:HD21	1:B:356:TRP:CD1	2.31	0.47
1:C:116:CYS:HB3	1:C:560:SER:CB	2.43	0.47
1:C:598:SER:HB3	1:C:631:GLU:HG2	1.97	0.47
1:A:113:PHE:O	1:A:575:PRO:HA	2.15	0.46
1:B:114:TYR:O	1:B:622:GLU:HA	2.15	0.46
1:B:194:PHE:HE1	1:B:344:LEU:HD13	1.79	0.46
1:C:599:ARG:NH1	1:C:617:LEU:HG	2.25	0.46
1:A:222:ALA:HB1	1:A:267:THR:HG21	1.96	0.46
1:C:543:ARG:HB3	1:C:550:ILE:HG21	1.97	0.46
1:D:138:GLU:HB2	2:F:1:NAG:H83	1.98	0.46
1:D:690:THR:OG1	1:D:693:GLU:HG3	2.14	0.46
1:D:174:TRP:CZ3	1:D:176:GLY:HA3	2.51	0.46
1:D:572:THR:HG22	1:D:573:CYS:H	1.80	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:120:THR:O	1:B:121:GLY:C	2.59	0.46
1:C:166:LYS:NZ	1:C:192:VAL:HG22	2.31	0.46
1:D:167:ASP:OD2	1:D:189:ARG:HG2	2.14	0.46
1:B:589:ILE:HG12	1:B:597:TYR:CE1	2.51	0.46
1:C:199:ASP:O	1:C:203:ALA:HB3	2.14	0.46
1:C:548:ASN:HB3	1:C:622:GLU:OE2	2.15	0.46
1:D:113:PHE:O	1:D:575:PRO:HA	2.15	0.46
1:B:601:LEU:HD13	1:B:627:ARG:NH1	2.30	0.46
1:A:548:ASN:OD1	1:A:560:SER:HA	2.15	0.46
1:A:559:VAL:HG12	1:A:572:THR:HA	1.97	0.46
1:B:250:THR:HB	1:B:270:ASN:CB	2.45	0.46
1:B:581:VAL:HG22	1:B:604:PHE:HB3	1.97	0.46
1:A:124:VAL:HG12	1:A:569:ALA:HA	1.98	0.46
1:C:190:ALA:HA	1:C:191:PRO:HD3	1.77	0.46
1:D:601:LEU:HA	1:D:616:GLN:HA	1.97	0.46
1:B:202:ASN:O	1:B:328:ARG:HB3	2.16	0.46
1:B:248:HIS:HA	1:B:271:CYS:O	2.16	0.46
1:B:270:ASN:O	1:B:272:ILE:HG13	2.16	0.46
1:C:157:TYR:CD1	1:C:363:VAL:HG12	2.51	0.46
1:C:329:ASP:O	1:C:333:LYS:HA	2.16	0.46
1:D:457:SER:O	1:D:458:ASN:HB2	2.16	0.46
1:B:646:GLY:HA3	1:B:659:LEU:O	2.17	0.45
1:A:520:MET:O	1:A:524:VAL:HG23	2.16	0.45
1:C:140:GLN:HE21	1:C:378:SER:HB2	1.80	0.45
1:C:285:ASP:HA	1:C:298:SER:HB2	1.97	0.45
1:D:252:LEU:HD23	1:D:252:LEU:N	2.31	0.45
1:D:440:GLN:NE2	1:D:471:GLU:HB3	2.32	0.45
1:C:147:ALA:HA	1:C:451:ALA:O	2.16	0.45
1:C:224:HIS:HB2	1:C:269:VAL:HB	1.97	0.45
1:C:319:PHE:CZ	1:C:321:GLN:HB2	2.51	0.45
1:D:498:THR:HG21	1:D:503:PHE:CE2	2.50	0.45
1:B:657:HIS:HD2	1:B:659:LEU:HG	1.81	0.45
1:C:502:GLU:HG3	1:C:503:PHE:N	2.31	0.45
1:D:584:GLN:CG	1:D:601:LEU:HB2	2.45	0.45
1:B:164:TYR:HD1	1:B:351:THR:O	2.00	0.45
1:B:332:THR:C	1:B:334:ALA:H	2.24	0.45
1:C:425:PHE:CZ	1:C:430:ASN:HA	2.51	0.45
1:A:701:ASP:OD1	1:A:704:GLU:HG3	2.16	0.45
1:A:202:ASN:OD1	1:A:327:ALA:HA	2.16	0.45
1:A:236:PRO:HA	1:A:246:GLY:O	2.17	0.45
1:D:164:TYR:HE1	1:D:353:ALA:HB3	1.81	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:394:THR:N	1:D:508:PHE:HB2	2.32	0.45
1:D:408:ASP:OD1	1:D:409:LEU:HG	2.16	0.45
1:A:162:THR:HA	1:A:275:GLU:HA	1.99	0.45
1:A:582:ILE:C	1:A:582:ILE:HD12	2.42	0.45
1:C:369:TRP:CD2	1:C:370:GLN:HG2	2.51	0.45
1:C:567:VAL:CG1	1:C:568:MET:N	2.80	0.45
1:B:194:PHE:CD1	1:B:320:LYS:HD2	2.51	0.45
1:B:378:SER:O	1:B:385:ARG:N	2.50	0.45
1:A:314:TYR:CE1	1:A:345:LEU:HD21	2.52	0.45
1:C:327:ALA:H	1:C:336:ALA:HB3	1.81	0.45
1:C:407:VAL:HA	1:C:493:GLU:O	2.17	0.45
1:A:328:ARG:HE	1:A:333:LYS:HA	1.82	0.45
1:C:325:PHE:O	1:C:339:PRO:HA	2.16	0.45
1:C:340:THR:HG21	1:C:355:ASP:OD2	2.17	0.45
1:D:207:CYS:HB3	1:D:249:THR:HG21	1.98	0.45
1:B:259:VAL:HG22	1:B:260:GLU:N	2.23	0.44
1:B:297:MET:HE1	1:B:319:PHE:HB2	1.99	0.44
1:B:443:LEU:HD12	1:B:448:PHE:O	2.17	0.44
1:C:342:ARG:HD2	1:C:354:TRP:O	2.17	0.44
1:B:202:ASN:O	1:B:328:ARG:HD3	2.18	0.44
1:A:160:LYS:HE3	1:A:277:ASP:OD1	2.17	0.44
1:A:200:LYS:HE3	1:A:208:ARG:CZ	2.48	0.44
1:C:115:VAL:O	1:C:117:PRO:HD3	2.17	0.44
1:C:605:ARG:HG2	1:C:611:PRO:O	2.16	0.44
1:B:328:ARG:HG3	1:B:335:ARG:HB3	1.99	0.44
1:B:129:GLN:H	1:B:129:GLN:CD	2.25	0.44
1:B:587:MET:HB2	1:B:653:TYR:HB3	1.99	0.44
1:D:191:PRO:HG3	1:D:350:PHE:N	2.32	0.44
1:B:147:ALA:HA	1:B:451:ALA:O	2.17	0.44
1:B:175:PHE:CE1	1:B:258:ARG:HA	2.53	0.44
1:A:125:VAL:O	1:A:567:VAL:HG13	2.18	0.44
1:A:584:GLN:HG2	1:A:601:LEU:C	2.43	0.44
1:A:635:VAL:HG12	1:A:636:GLY:N	2.32	0.44
1:D:659:LEU:HB2	1:D:663:ASP:OD1	2.17	0.44
1:C:186:PHE:CZ	1:C:188:ASP:HB2	2.52	0.44
1:C:406:ARG:HG2	1:C:406:ARG:O	2.17	0.44
1:B:343:ASN:OD1	1:B:356:TRP:HB2	2.18	0.44
1:B:679:GLU:H	1:B:679:GLU:CD	2.25	0.44
1:A:166:LYS:NZ	1:A:192:VAL:HG22	2.33	0.44
1:C:116:CYS:CB	1:C:560:SER:HB2	2.45	0.44
1:B:583:VAL:HG22	1:B:584:GLN:N	2.33	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:172:GLN:HG2	1:A:183:MET:HE2	1.99	0.44
1:C:600:PRO:O	1:C:616:GLN:HB2	2.18	0.44
1:D:250:THR:HB	1:D:270:ASN:HB2	2.00	0.44
1:D:375:MET:O	1:D:452:TYR:HE1	2.01	0.44
1:C:279:ARG:O	1:C:288:VAL:HG22	2.18	0.43
1:D:433:HIS:HA	1:D:456:LEU:O	2.17	0.43
1:B:145:GLY:HA2	1:B:455:LEU:HD12	2.00	0.43
1:A:332:THR:O	1:A:333:LYS:HB2	2.18	0.43
1:C:235:LYS:HG2	1:C:248:HIS:O	2.18	0.43
1:C:620:ASN:C	1:C:622:GLU:H	2.27	0.43
1:D:280:SER:HB2	1:D:287:PHE:CB	2.47	0.43
1:B:175:PHE:HD1	1:B:258:ARG:HH11	1.66	0.43
1:A:181:GLN:HG2	1:A:182:PHE:O	2.17	0.43
1:C:209:SER:HB2	1:C:224:HIS:HB3	2.00	0.43
1:B:168:VAL:HA	1:B:268:THR:O	2.18	0.43
1:A:250:THR:HG22	1:A:251:ASP:N	2.34	0.43
1:C:434:ILE:HD12	1:C:434:ILE:O	2.18	0.43
1:C:591:SER:C	1:C:593:PRO:HD3	2.44	0.43
1:A:382:GLY:HA2	1:A:399:LEU:HD11	2.00	0.43
1:A:384:PHE:CE2	1:A:399:LEU:HA	2.54	0.43
1:C:170:VAL:HA	1:C:266:GLY:O	2.18	0.43
1:C:559:VAL:HG12	1:C:572:THR:HG23	2.00	0.43
1:C:635:VAL:HG22	1:C:636:GLY:H	1.83	0.43
1:B:156:PRO:O	1:B:158:LYS:HG3	2.18	0.43
1:A:402:TYR:HA	1:A:403:PRO:HD3	1.72	0.43
1:C:384:PHE:O	1:C:396:THR:HA	2.19	0.43
2:F:1:NAG:O3	2:F:2:NAG:O5	2.34	0.43
1:C:193:PRO:O	1:C:197:VAL:HG23	2.18	0.43
1:B:106:ALA:HA	1:B:643:PHE:HE2	1.83	0.43
1:A:574:VAL:HA	1:A:575:PRO:HD3	1.86	0.43
1:C:327:ALA:H	1:C:336:ALA:CB	2.32	0.43
1:D:327:ALA:HB3	1:D:336:ALA:HB2	2.00	0.43
1:B:110:ASP:OD1	1:B:645:GLY:HA3	2.18	0.43
1:B:175:PHE:CD1	1:B:258:ARG:HD2	2.54	0.43
1:B:616:GLN:NE2	1:B:629:ALA:HB3	2.33	0.43
1:C:113:PHE:CD1	1:C:581:VAL:HG21	2.54	0.43
1:A:276:VAL:HB	1:A:290:ALA:HB3	2.01	0.42
1:C:104:ILE:O	1:C:582:ILE:HB	2.19	0.42
1:C:105:LYS:HB3	1:C:106:ALA:H	1.51	0.42
1:C:406:ARG:HG2	1:C:493:GLU:HB3	2.00	0.42
1:D:106:ALA:HB2	1:D:643:PHE:CZ	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:392:SER:HA	1:D:508:PHE:CZ	2.54	0.42
1:D:616:GLN:HE22	1:D:629:ALA:H	1.65	0.42
1:A:236:PRO:HA	1:A:247:TRP:CD1	2.54	0.42
1:A:256:PRO:HG3	1:A:266:GLY:N	2.35	0.42
1:C:319:PHE:O	1:C:320:LYS:HD3	2.20	0.42
1:C:340:THR:HG21	1:C:355:ASP:CG	2.43	0.42
1:C:458:ASN:HA	1:C:461:ALA:HB2	2.02	0.42
1:B:225:ARG:HD2	1:B:254:TYR:CD1	2.53	0.42
1:B:408:ASP:OD1	1:B:409:LEU:HG	2.20	0.42
1:B:574:VAL:HA	1:B:575:PRO:HD3	1.88	0.42
1:A:402:TYR:OH	1:A:495:ILE:HD11	2.19	0.42
1:C:619:GLU:O	1:C:622:GLU:HB3	2.19	0.42
1:D:234:LEU:HD23	1:D:249:THR:HG23	2.00	0.42
1:A:224:HIS:C	1:A:226:ASP:H	2.28	0.42
1:A:369:TRP:CD2	1:A:370:GLN:HG2	2.55	0.42
1:C:304:ARG:NH1	5:C:2001:MRY:HAA2	2.35	0.42
1:D:113:PHE:CD2	1:D:623:LEU:HD21	2.55	0.42
1:D:144:GLU:HA	1:D:376:LEU:HD23	2.01	0.42
1:D:343:ASN:O	1:D:353:ALA:HA	2.20	0.42
1:A:147:ALA:HB2	1:A:452:TYR:HD1	1.85	0.42
1:A:166:LYS:HG2	1:A:271:CYS:HA	2.00	0.42
1:A:298:SER:CB	1:A:310:GLU:HB3	2.48	0.42
1:C:148:VAL:HB	1:C:451:ALA:HB3	2.01	0.42
1:D:155:ALA:HA	1:D:156:PRO:HD3	1.82	0.42
1:A:396:THR:C	1:A:397:THR:HG22	2.45	0.42
1:C:276:VAL:HB	1:C:290:ALA:HB3	2.00	0.42
1:C:574:VAL:HA	1:C:575:PRO:HD3	1.90	0.42
1:C:589:ILE:HG22	1:C:592:ARG:H	1.84	0.42
1:D:616:GLN:NE2	1:D:629:ALA:HB3	2.35	0.42
1:D:617:LEU:HD12	1:D:622:GLU:O	2.20	0.42
1:A:127:PHE:HD2	1:A:127:PHE:HA	1.75	0.42
1:A:196:GLU:O	1:A:200:LYS:HB2	2.19	0.42
1:A:587:MET:O	1:A:596:CYS:HB3	2.19	0.42
1:C:237:ALA:HA	1:C:248:HIS:CD2	2.53	0.42
1:D:175:PHE:CD1	1:D:258:ARG:HD2	2.54	0.42
1:D:593:PRO:C	1:D:595:ALA:H	2.28	0.42
1:A:190:ALA:HA	1:A:191:PRO:HD3	1.74	0.42
1:C:157:TYR:HB2	1:C:284:TYR:CE2	2.55	0.42
1:C:604:PHE:O	1:C:612:LEU:HA	2.19	0.42
1:D:397:THR:HG22	1:D:444:ALA:HA	2.01	0.42
1:D:511:ASN:O	1:D:515:ARG:HG2	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:392:SER:HA	1:B:508:PHE:CE2	2.55	0.42
1:C:130:PRO:HB3	1:C:533:ASN:HB3	2.02	0.42
1:D:589:ILE:HG12	1:D:597:TYR:CE1	2.54	0.42
1:B:402:TYR:HA	1:B:403:PRO:HD3	1.89	0.41
1:B:492:VAL:HG12	1:B:493:GLU:N	2.35	0.41
1:C:119:PRO:HG3	1:C:560:SER:O	2.20	0.41
1:C:280:SER:HB2	1:C:287:PHE:HB3	2.02	0.41
1:C:606:TYR:HD2	1:C:613:VAL:HG21	1.81	0.41
1:B:637:HIS:CE1	1:B:639:ARG:HD2	2.54	0.41
1:A:542:ALA:C	1:A:550:ILE:HG12	2.45	0.41
1:A:671:ILE:HD12	1:A:671:ILE:N	2.35	0.41
1:C:233:GLU:O	1:C:249:THR:HG22	2.18	0.41
1:C:425:PHE:O	1:C:430:ASN:HB3	2.20	0.41
1:C:433:HIS:HA	1:C:456:LEU:O	2.20	0.41
1:C:557:ARG:HG3	1:C:559:VAL:HG13	2.02	0.41
1:B:288:VAL:HG12	1:B:294:PHE:HA	2.02	0.41
1:B:359:LYS:O	1:B:363:VAL:HG22	2.19	0.41
1:A:528:TRP:O	1:A:532:GLN:HG2	2.20	0.41
1:C:248:HIS:HE1	6:C:733:HOH:O	2.02	0.41
1:C:402:TYR:HA	1:C:403:PRO:HD3	1.79	0.41
1:C:558:ARG:NH2	1:C:622:GLU:HB2	2.36	0.41
1:A:248:HIS:HA	1:A:271:CYS:O	2.21	0.41
1:C:129:GLN:HB3	1:C:130:PRO:HD2	2.02	0.41
1:C:191:PRO:HA	1:C:350:PHE:HA	2.02	0.41
1:C:298:SER:HB3	1:C:310:GLU:HB3	2.02	0.41
1:D:572:THR:HG22	1:D:573:CYS:N	2.35	0.41
1:B:514:GLN:O	1:B:517:VAL:HG22	2.21	0.41
1:B:659:LEU:HB3	1:B:663:ASP:OD1	2.20	0.41
1:D:146:ILE:HD11	1:D:429:TYR:CD1	2.55	0.41
1:B:675:ILE:HD12	1:B:675:ILE:C	2.45	0.41
1:A:604:PHE:O	1:A:613:VAL:N	2.49	0.41
1:C:587:MET:HE3	1:C:653:TYR:CD2	2.56	0.41
1:D:145:GLY:HA3	1:D:452:TYR:CZ	2.55	0.41
1:B:411:ASP:HB3	1:B:415:LYS:HE3	2.02	0.41
1:B:520:MET:HE1	1:B:523:ARG:CZ	2.51	0.41
1:A:384:PHE:O	1:A:396:THR:HA	2.20	0.41
1:D:190:ALA:HA	1:D:191:PRO:HD3	1.87	0.41
1:D:420:ALA:O	1:D:424:ILE:HG13	2.21	0.41
1:D:605:ARG:HH11	1:D:609:GLN:HA	1.85	0.41
2:E:1:NAG:O7	2:E:1:NAG:C3	2.69	0.41
1:B:133:CYS:HA	1:B:134:PRO:HD3	1.91	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:282:TYR:CG	1:C:283:PRO:HA	2.56	0.41
1:C:519:ASP:O	1:C:523:ARG:HG3	2.21	0.41
1:D:647:TYR:HE2	1:D:661:ARG:HG2	1.86	0.41
1:B:122:ALA:HA	1:B:123:THR:HA	1.66	0.41
1:B:264:ARG:HG3	1:B:265[B]:TYR:H	1.84	0.41
1:B:368:LYS:NZ	1:B:368:LYS:HB3	2.35	0.41
1:B:401:GLU:HB3	1:B:402:TYR:H	1.75	0.41
1:A:152:GLU:HG2	1:A:153:ASN:N	2.36	0.41
1:C:326:TYR:HD1	1:C:336:ALA:O	2.04	0.41
1:C:539:TRP:HA	1:C:542:ALA:HB3	2.03	0.41
1:C:690:THR:OG1	1:C:693:GLU:HG3	2.21	0.41
1:D:372:VAL:HG11	1:D:375:MET:HE2	2.03	0.41
1:D:639:ARG:HB3	1:D:641:PHE:CZ	2.55	0.41
1:B:555:VAL:HG12	1:B:555:VAL:O	2.21	0.41
1:A:302:GLY:O	1:A:307:SER:HB2	2.21	0.41
1:A:441:TYR:HB3	1:A:449:LEU:HD11	2.02	0.41
1:C:321:GLN:NE2	1:C:341:THR:HG22	2.36	0.40
1:C:641:PHE:HB2	1:C:648:VAL:HG13	2.03	0.40
1:D:116:CYS:HB3	1:D:560:SER:HB3	2.03	0.40
1:D:175:PHE:CE1	1:D:258:ARG:HA	2.55	0.40
1:B:392:SER:HA	1:B:508:PHE:CZ	2.56	0.40
1:A:237:ALA:H	1:A:246:GLY:C	2.28	0.40
1:C:347:THR:N	1:C:350:PHE:O	2.54	0.40
1:D:114:TYR:HB3	1:D:573:CYS:HB3	2.04	0.40
1:D:117:PRO:O	1:D:118:PRO:C	2.64	0.40
1:D:174:TRP:CH2	1:D:176:GLY:HA3	2.56	0.40
1:D:659:LEU:HD23	1:D:659:LEU:N	2.36	0.40
1:A:200:LYS:HE3	1:A:208:ARG:NH1	2.36	0.40
1:A:549:ALA:O	1:A:553:VAL:HG12	2.22	0.40
1:C:340:THR:HG22	1:C:341:THR:N	2.36	0.40
1:D:200:LYS:HB3	1:D:206:VAL:HG23	2.04	0.40
1:B:157:TYR:CD1	1:B:363:VAL:HG12	2.57	0.40
1:B:224:HIS:CD2	1:B:225:ARG:HG2	2.56	0.40
1:B:303:TYR:O	1:A:317:ASP:HB3	2.21	0.40
1:B:694:ILE:C	1:B:696:ASP:N	2.79	0.40
1:A:224:HIS:HB2	1:A:269:VAL:HB	2.02	0.40
1:C:215:ARG:NH2	1:C:349:LYS:HG2	2.37	0.40
1:C:250:THR:HG22	1:C:251:ASP:N	2.36	0.40
1:C:343:ASN:O	1:C:353:ALA:HA	2.21	0.40
1:C:601:LEU:HA	1:C:616:GLN:HA	2.03	0.40
1:C:304:ARG:HD3	1:C:356:TRP:CZ3	2.57	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:597:TYR:HA	1:C:630:ILE:HA	2.04	0.40
1:D:175:PHE:HB2	1:D:259:VAL:O	2.21	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	603/703 (86%)	543 (90%)	56 (9%)	4 (1%)	18	44
1	B	590/703 (84%)	523 (89%)	64 (11%)	3 (0%)	24	51
1	C	595/703 (85%)	532 (89%)	58 (10%)	5 (1%)	16	41
1	D	601/703 (86%)	548 (91%)	52 (9%)	1 (0%)	43	70
All	All	2389/2812 (85%)	2146 (90%)	230 (10%)	13 (0%)	24	51

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	259	VAL
1	B	413	ILE
1	A	238	ASN
1	C	429	TYR
1	B	106	ALA
1	D	315	ALA
1	A	134	PRO
1	A	264	ARG
1	A	699	LEU
1	C	390	ALA
1	C	413	ILE
1	C	134	PRO
1	C	139	GLY

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	517/593 (87%)	507 (98%)	10 (2%)	50	77
1	B	509/593 (86%)	498 (98%)	11 (2%)	45	75
1	C	506/593 (85%)	498 (98%)	8 (2%)	55	81
1	D	514/593 (87%)	509 (99%)	5 (1%)	68	87
All	All	2046/2372 (86%)	2012 (98%)	34 (2%)	53	80

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	115	VAL
1	B	135	THR
1	B	138	GLU
1	B	141	ASN
1	B	143	THR
1	B	264	ARG
1	B	368	LYS
1	B	392	SER
1	B	517	VAL
1	B	606	TYR
1	B	625	LEU
1	A	127	PHE
1	A	170	VAL
1	A	185	ILE
1	A	262	PHE
1	A	323	ASP
1	A	331	THR
1	A	397	THR
1	A	401	GLU
1	A	502	GLU
1	A	675	ILE
1	C	183	MET
1	C	265	TYR
1	C	270	ASN

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Mol	Chain	Res	Type
1	C	468	HIS
1	C	555	VAL
1	C	613	VAL
1	C	631	GLU
1	C	675	ILE
1	D	173	VAL
1	D	340	THR
1	D	368	LYS
1	D	392	SER
1	D	546	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	321	GLN
1	B	453	GLN
1	B	516	HIS
1	B	533	ASN
1	B	580	ASN
1	B	584	GLN
1	A	228	HIS
1	A	440	GLN
1	A	620	ASN
1	C	140	GLN
1	C	270	ASN
1	C	468	HIS
1	D	228	HIS
1	D	472	GLN
1	D	532	GLN
1	D	637	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates

4 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	E	1	1,2	14,14,15	0.62	0	17,19,21	1.08	1 (5%)
2	NAG	E	2	2	14,14,15	0.46	0	17,19,21	0.68	0
2	NAG	F	1	1,2	14,14,15	0.54	0	17,19,21	1.51	4 (23%)
2	NAG	F	2	2	14,14,15	0.50	0	17,19,21	0.83	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
2	NAG	E	1	1,2	1/1/5/7	3/6/23/26	0/1/1/1
2	NAG	E	2	2	-	0/6/23/26	0/1/1/1
2	NAG	F	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	F	2	2	-	1/6/23/26	0/1/1/1

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	1	NAG	C2-N2-C7	-3.66	118.00	122.90
2	F	1	NAG	C1-O5-C5	2.57	115.63	112.19
2	E	1	NAG	C2-N2-C7	2.57	126.34	122.90
2	F	1	NAG	O7-C7-C8	-2.55	117.51	122.05
2	F	1	NAG	O7-C7-N2	2.36	126.15	121.98
2	F	2	NAG	C2-N2-C7	-2.12	120.06	122.90

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	E	1	NAG	C1

All (6) torsion outliers are listed below:

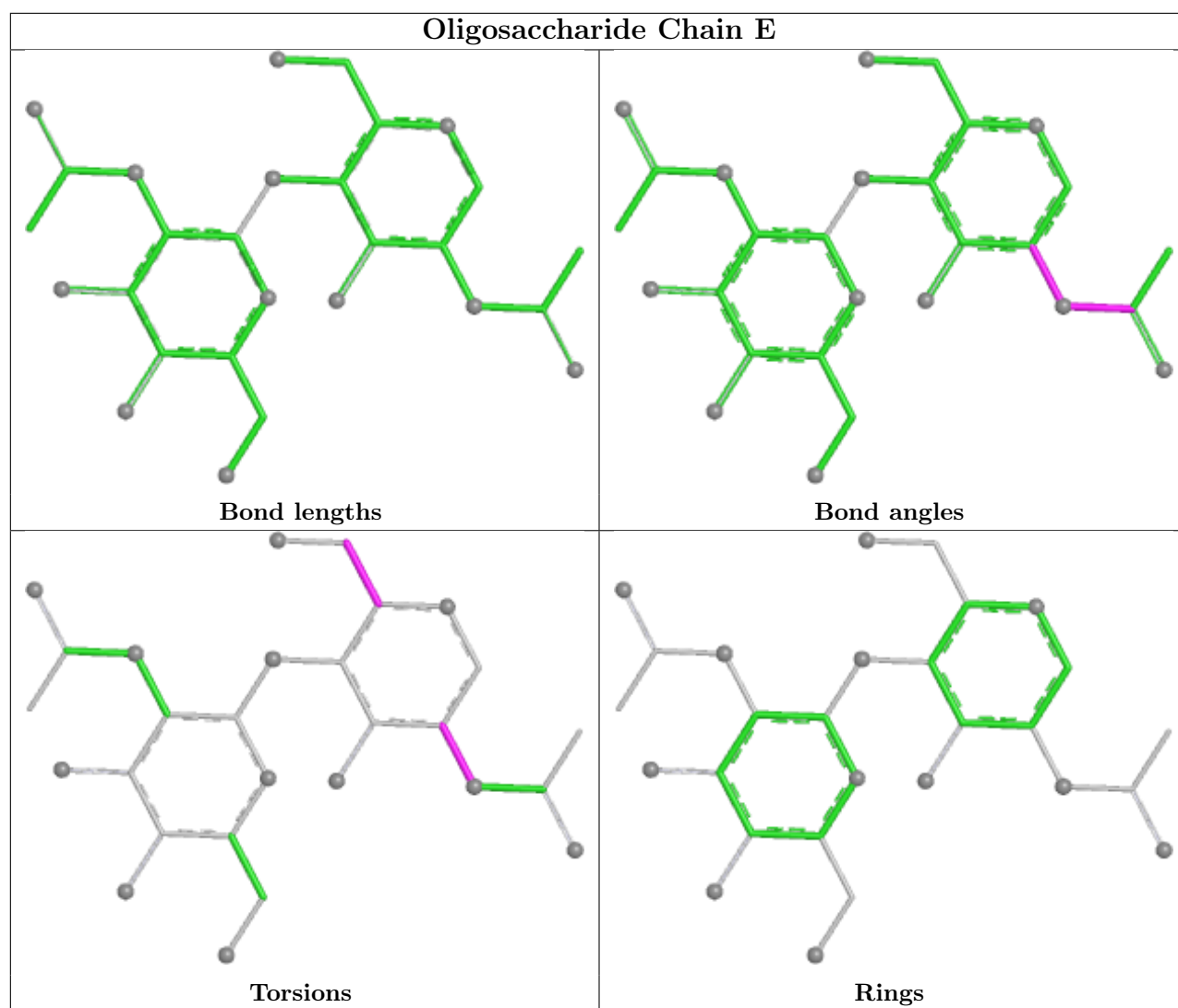
Mol	Chain	Res	Type	Atoms
2	E	1	NAG	C3-C2-N2-C7
2	F	1	NAG	C8-C7-N2-C2
2	F	1	NAG	O7-C7-N2-C2
2	E	1	NAG	C4-C5-C6-O6
2	F	2	NAG	O5-C5-C6-O6
2	E	1	NAG	O5-C5-C6-O6

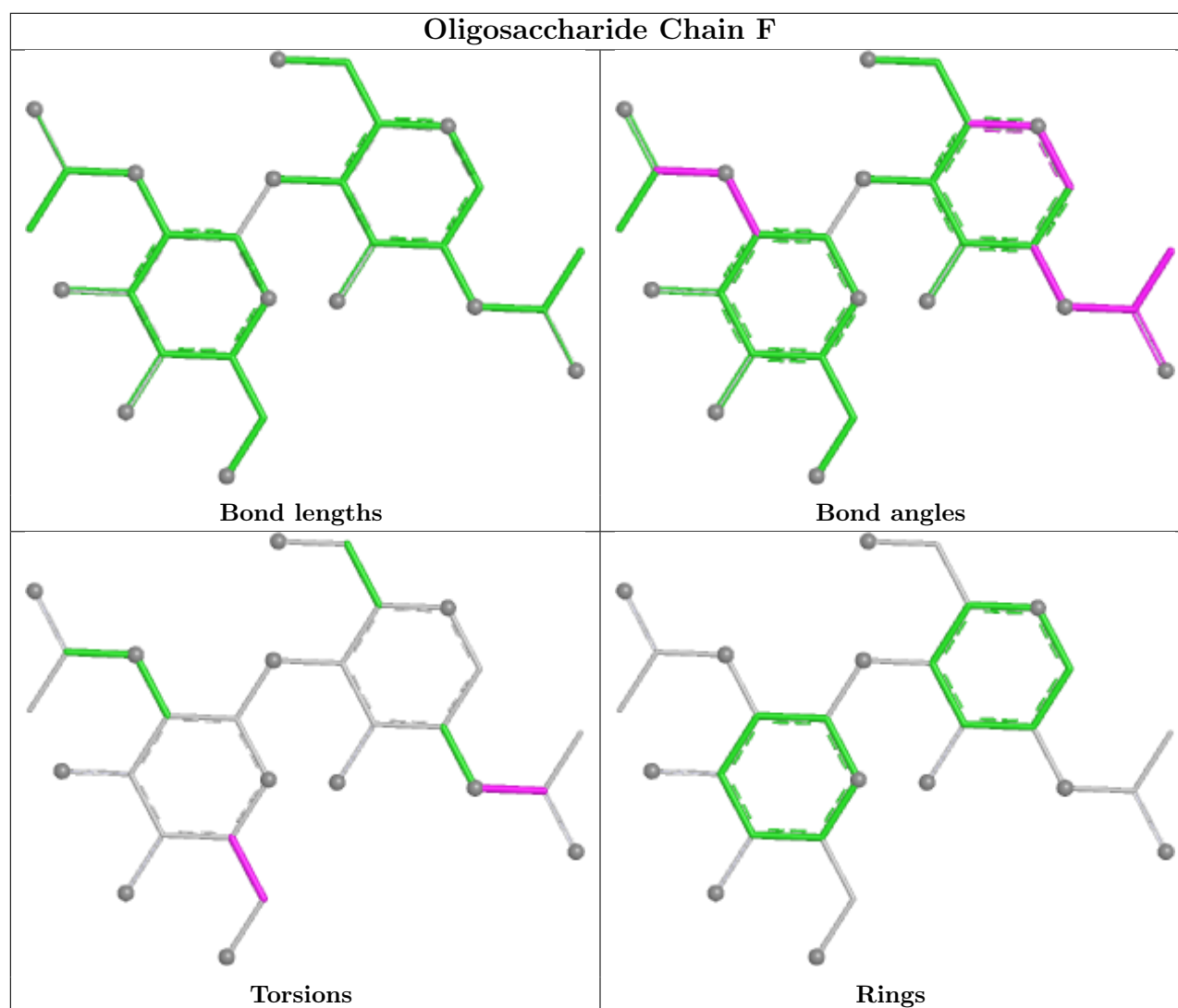
There are no ring outliers.

3 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	F	2	NAG	1	0
2	E	1	NAG	2	0
2	F	1	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 oligosaccharide.





5.6 Ligand geometry [i](#)

Of 12 ligands modelled in this entry, 2 are monoatomic - leaving 10 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	NAG	A	1674	1	14,14,15	0.51	0	17,19,21	0.61	0
3	NAG	A	1141	1	14,14,15	0.50	0	17,19,21	0.63	0
3	NAG	C	1430	1	14,14,15	0.48	0	17,19,21	1.15	1 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	NAG	B	1430	1	14,14,15	0.60	0	17,19,21	1.29	2 (11%)
3	NAG	C	1141	1	14,14,15	0.52	0	17,19,21	0.58	0
5	MRY	C	2001	-	7,7,7	0.41	0	8,8,8	0.35	0
3	NAG	A	1430	1	14,14,15	0.56	0	17,19,21	0.87	1 (5%)
3	NAG	D	1398	1	14,14,15	0.54	0	17,19,21	0.55	0
3	NAG	B	1398	1	14,14,15	0.54	0	17,19,21	0.60	0
3	NAG	C	1674	1	14,14,15	0.52	0	17,19,21	0.59	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	A	1674	1	-	0/6/23/26	0/1/1/1
3	NAG	A	1141	1	-	2/6/23/26	0/1/1/1
3	NAG	C	1430	1	-	0/6/23/26	0/1/1/1
3	NAG	B	1430	1	-	4/6/23/26	0/1/1/1
3	NAG	C	1141	1	-	4/6/23/26	0/1/1/1
5	MRY	C	2001	-	-	0/8/8/8	-
3	NAG	A	1430	1	-	2/6/23/26	0/1/1/1
3	NAG	D	1398	1	-	3/6/23/26	0/1/1/1
3	NAG	B	1398	1	-	0/6/23/26	0/1/1/1
3	NAG	C	1674	1	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	1430	NAG	C1-O5-C5	3.32	116.63	112.19
3	B	1430	NAG	O5-C1-C2	3.25	116.33	111.29
3	B	1430	NAG	C1-O5-C5	2.86	116.02	112.19
3	A	1430	NAG	C1-O5-C5	2.39	115.39	112.19

There are no chirality outliers.

All (15) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	1430	NAG	C3-C2-N2-C7

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Mol	Chain	Res	Type	Atoms
3	B	1430	NAG	C8-C7-N2-C2
3	B	1430	NAG	O7-C7-N2-C2
3	C	1141	NAG	C3-C2-N2-C7
3	C	1141	NAG	C8-C7-N2-C2
3	C	1141	NAG	O7-C7-N2-C2
3	D	1398	NAG	C3-C2-N2-C7
3	D	1398	NAG	C8-C7-N2-C2
3	D	1398	NAG	O7-C7-N2-C2
3	A	1141	NAG	C4-C5-C6-O6
3	A	1141	NAG	O5-C5-C6-O6
3	A	1430	NAG	C8-C7-N2-C2
3	A	1430	NAG	O7-C7-N2-C2
3	B	1430	NAG	O5-C5-C6-O6
3	C	1141	NAG	O5-C5-C6-O6

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	C	2001	MRY	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

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

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	606/703 (86%)	-1.48	2 (0%) 90 87	20, 66, 145, 550	1 (0%)
1	B	597/703 (84%)	-1.33	7 (1%) 76 70	23, 71, 174, 550	1 (0%)
1	C	601/703 (85%)	-1.39	2 (0%) 90 87	20, 76, 176, 550	0
1	D	606/703 (86%)	-1.42	3 (0%) 87 83	22, 64, 151, 550	1 (0%)
All	All	2410/2812 (85%)	-1.40	14 (0%) 85 81	20, 69, 164, 550	3 (0%)

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	612	LEU	4.9
1	B	612	LEU	4.6
1	D	603	SER	4.4
1	B	581	VAL	3.3
1	C	594	GLY	2.8
1	B	600	PRO	2.7
1	B	613	VAL	2.7
1	D	613	VAL	2.7
1	B	578	ALA	2.6
1	B	599	ARG	2.4
1	C	595	ALA	2.3
1	A	594	GLY	2.1
1	B	402	TYR	2.1
1	A	595	ALA	2.0

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

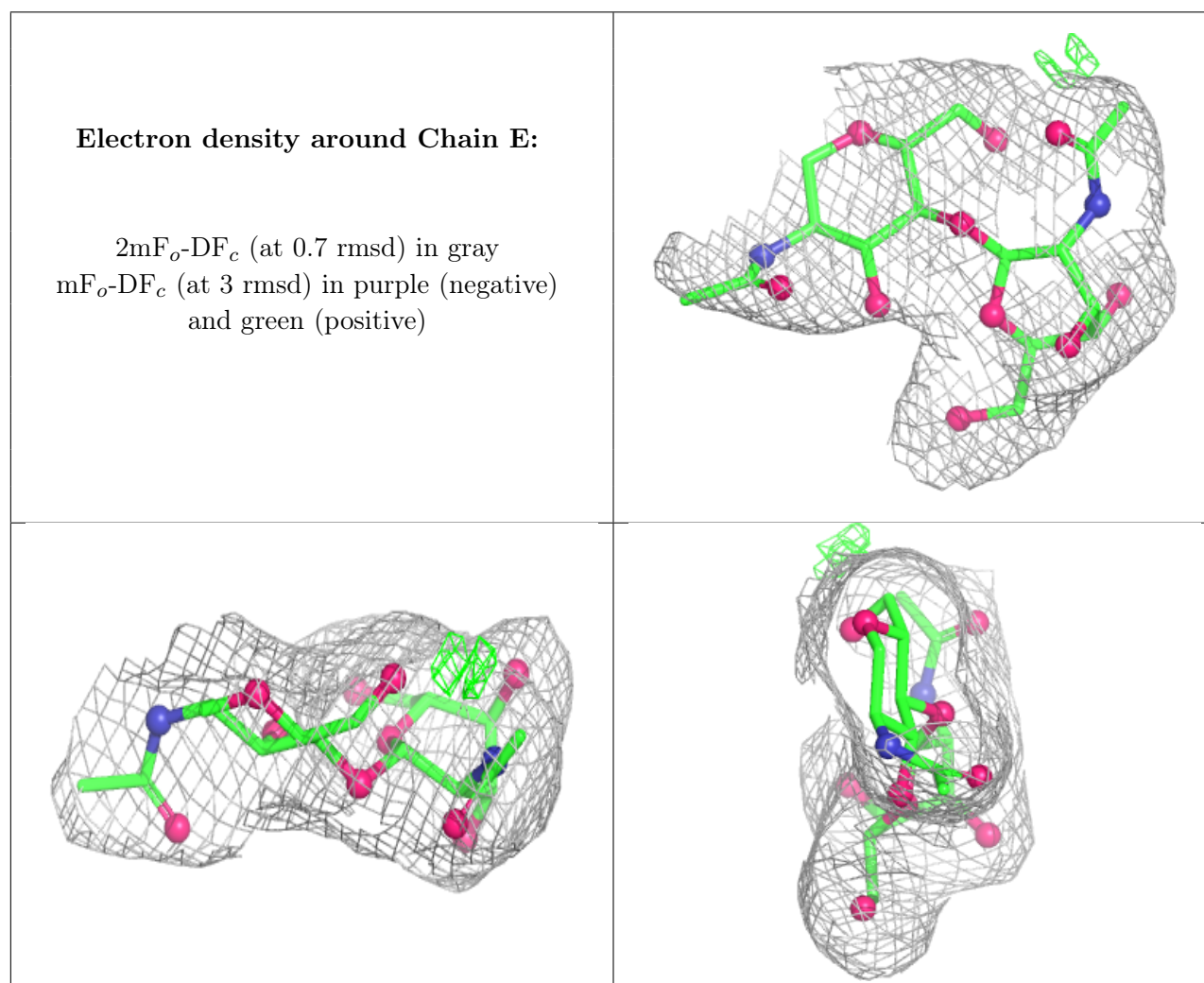
There are no non-standard protein/DNA/RNA residues in this entry.

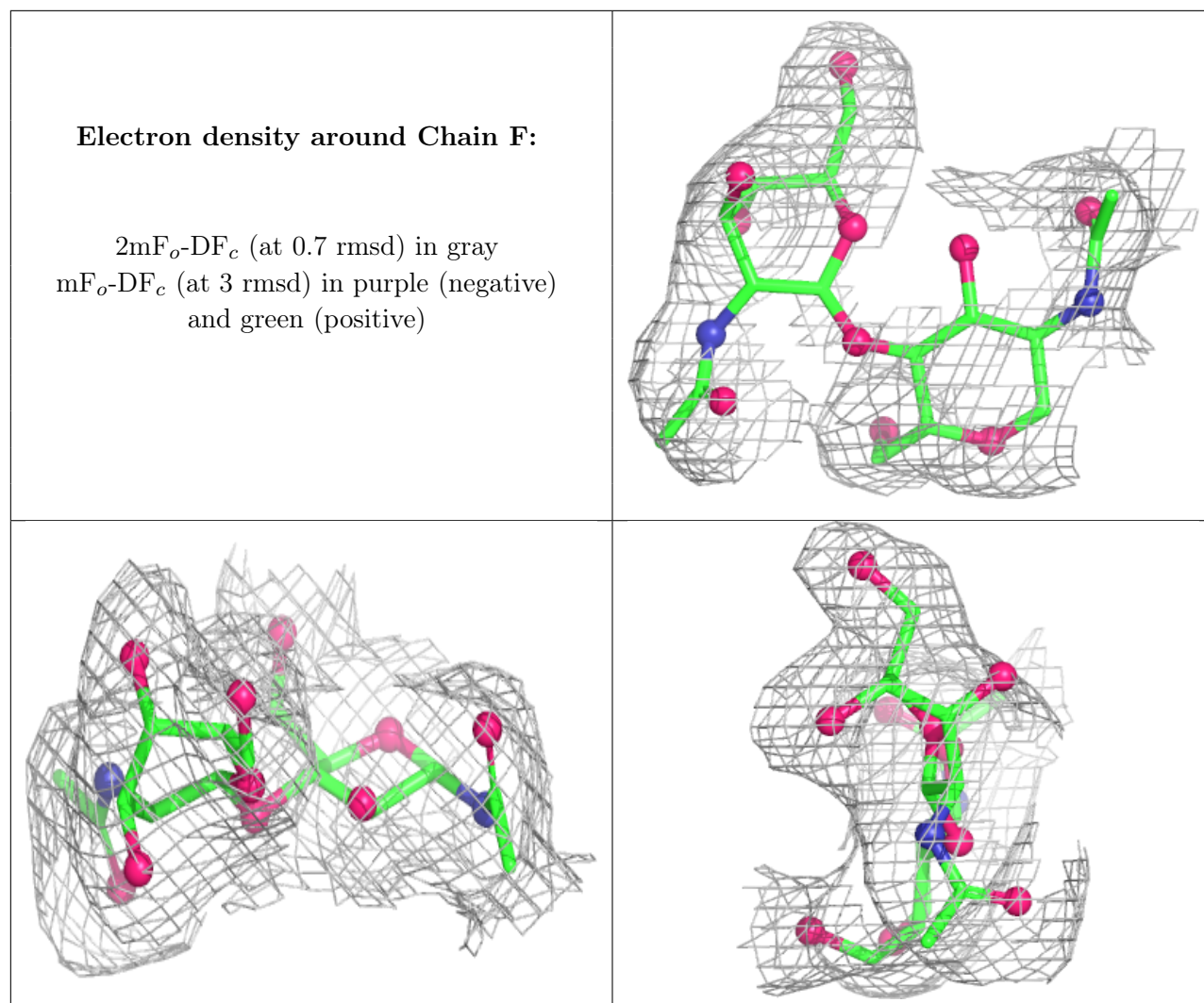
6.3 Carbohydrates [i](#)

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(Å ²)	Q<0.9
2	NAG	E	1	14/15	0.98	0.04	104,104,104,104	0
2	NAG	E	2	14/15	0.98	0.03	89,89,89,89	0
2	NAG	F	2	14/15	0.98	0.04	91,91,91,91	0
2	NAG	F	1	14/15	0.99	0.03	77,77,77,77	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 ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	NAG	C	1430	14/15	0.94	0.05	117,117,117,117	0
3	NAG	A	1430	14/15	0.96	0.05	118,118,118,118	0
3	NAG	B	1430	14/15	0.96	0.05	117,117,117,117	0
3	NAG	A	1674	14/15	0.98	0.04	89,89,89,89	0
3	NAG	A	1141	14/15	0.98	0.05	120,120,120,120	0
3	NAG	C	1674	14/15	0.98	0.04	77,77,77,77	0
3	NAG	D	1398	14/15	0.98	0.05	87,87,87,87	0
3	NAG	C	1141	14/15	0.99	0.04	113,113,113,113	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	NAG	B	1398	14/15	0.99	0.03	92,92,92,92	0
4	CL	A	2	1/1	0.99	0.04	78,78,78,78	0
4	CL	D	1	1/1	0.99	0.04	59,59,59,59	0
5	MRY	C	2001	8/8	0.99	0.04	69,69,69,69	0

6.5 Other polymers [i](#)

There are no such residues in this entry.