



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 08:13 AM UTC

PDB ID : 2Y7Q / pdb_00002y7q
Title : THE HIGH-AFFINITY COMPLEX BETWEEN IGE AND ITS RECEPTOR
FC EPSILON RI
Authors : Davies, A.M.; Holdom, M.D.; Nettleship, J.E.; Beavil, A.J.; Owens, R.J.;
Sutton, B.J.
Deposited on : 2011-02-01
Resolution : 3.40 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

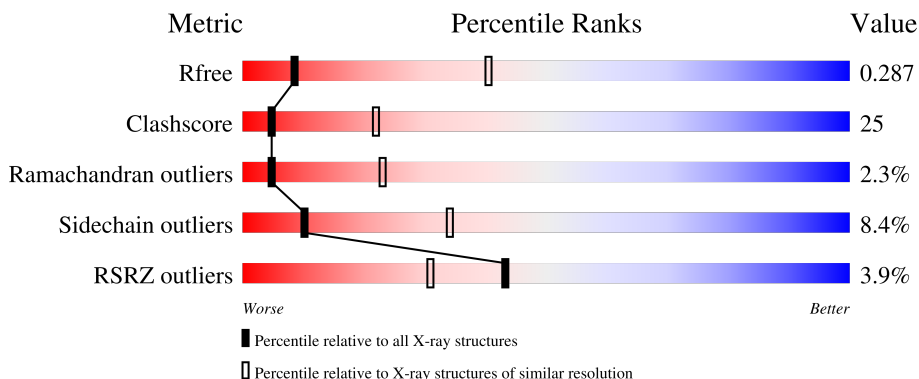
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.40 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	1001 (3.44-3.36)
Clashscore	190562	1022 (3.44-3.36)
Ramachandran outliers	187476	1012 (3.44-3.36)
Sidechain outliers	187428	1012 (3.44-3.36)
RSRZ outliers	180081	1001 (3.44-3.36)

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	188	
2	B	327	
2	D	327	
3	C	4	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5355 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 HIGH AFFINITY IMMUNOGLOBULIN EPSILON RECEPTOR SUBUNIT ALPHA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	158	Total	C	N	O	S	0	0	0
			1188	773	189	222	4			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLU	-	cloning artifact	UNP P12319
A	-1	THR	-	cloning artifact	UNP P12319
A	0	GLY	-	cloning artifact	UNP P12319
A	74	ALA	ASN	engineered mutation	UNP P12319
A	135	ALA	ASN	engineered mutation	UNP P12319
A	142	ALA	THR	engineered mutation	UNP P12319
A	143	ALA	VAL	cloning artifact	UNP P12319
A	177	GLY	-	cloning artifact	UNP P12319
A	178	THR	-	cloning artifact	UNP P12319
A	179	LYS	-	cloning artifact	UNP P12319
A	180	HIS	-	expression tag	UNP P12319
A	181	HIS	-	expression tag	UNP P12319
A	182	HIS	-	expression tag	UNP P12319
A	183	HIS	-	expression tag	UNP P12319
A	184	HIS	-	expression tag	UNP P12319
A	185	HIS	-	expression tag	UNP P12319

- Molecule 2 is a protein called IG EPSILON CHAIN C REGION.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	287	Total	C	N	O	S	0	0	0
			2043	1290	351	391	11			
2	D	290	Total	C	N	O	S	0	0	0
			2032	1287	350	385	10			

There are 10 discrepancies between the modelled and reference sequences:

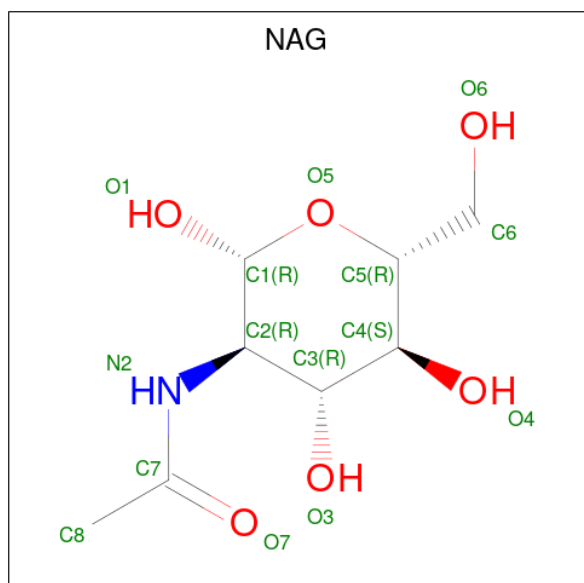
Chain	Residue	Modelled	Actual	Comment	Reference
B	222	ASP	-	expression tag	UNP P01854
B	223	ILE	-	expression tag	UNP P01854
B	225	ALA	CYS	engineered mutation	UNP P01854
B	265	GLN	ASN	engineered mutation	UNP P01854
B	371	GLN	ASN	engineered mutation	UNP P01854
D	222	ASP	-	expression tag	UNP P01854
D	223	ILE	-	expression tag	UNP P01854
D	225	ALA	CYS	engineered mutation	UNP P01854
D	265	GLN	ASN	engineered mutation	UNP P01854
D	371	GLN	ASN	engineered mutation	UNP P01854

- Molecule 3 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-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
3	C	4	Total	C	N	O	0	0	0
			50	28	2	20			

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

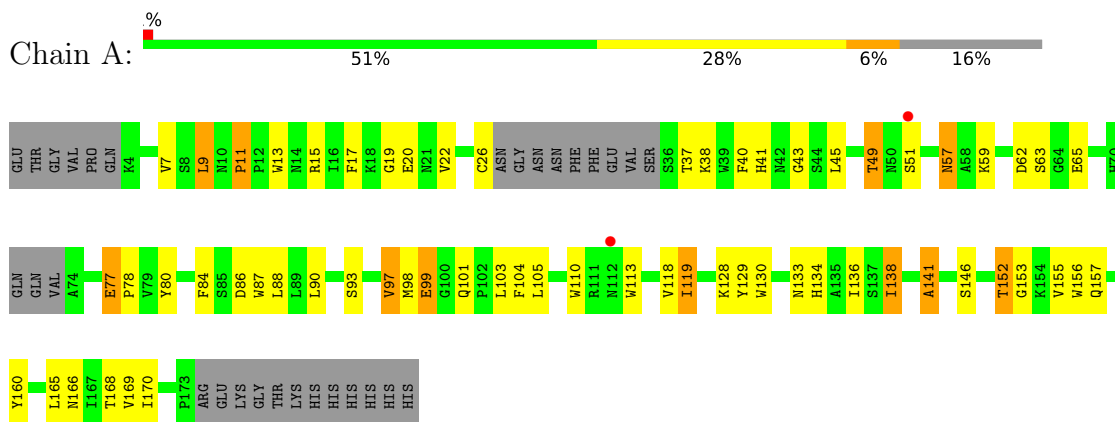


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total 14	C 8	N 1	O 5	0	0
4	B	1	Total 14	C 8	N 1	O 5	0	0
4	D	1	Total 14	C 8	N 1	O 5	0	0

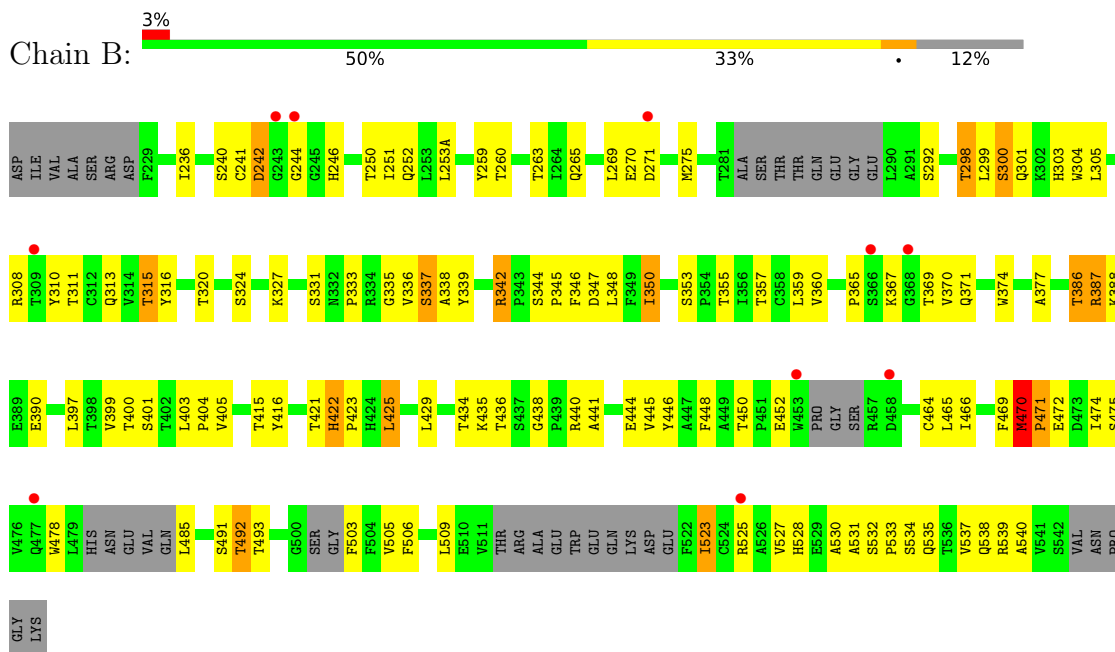
3 Residue-property plots [i](#)

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

- Molecule 1: HIGH AFFINITY IMMUNOGLOBULIN EPSILON RECEPTOR SUBUNIT ALPHA

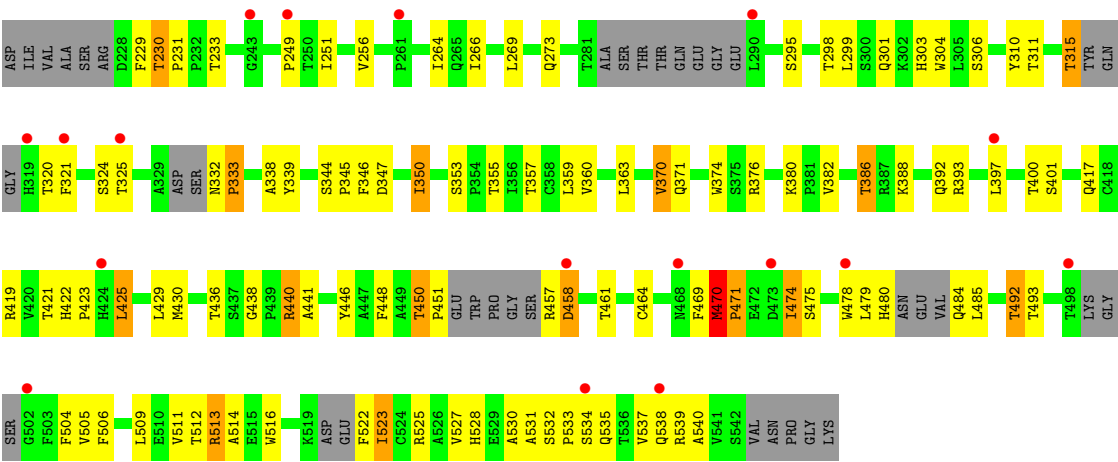


- Molecule 2: IG EPSILON CHAIN C REGION



- Molecule 2: IG EPSILON CHAIN C REGION





● Molecule 3: alpha-D-mannopyranose-(1-3)-beta-D-mannopyranose-(1-4)-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 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	99.49Å 103.34Å 110.12Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	37.68 – 3.40 37.68 – 3.40	Depositor EDS
% Data completeness (in resolution range)	99.7 (37.68-3.40) 99.7 (37.68-3.40)	Depositor EDS
R_{merge}	0.25	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.81 (at 3.40Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.245 , 0.292 0.238 , 0.287	Depositor DCC
R_{free} test set	812 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	85.6	Xtriage
Anisotropy	0.275	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 107.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	0.047 for k,h,-l	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	5355	wwPDB-VP
Average B, all atoms (Å ²)	77.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.51% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NAG, BMA, 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	0.41	0/1225	0.83	2/1679 (0.1%)
2	B	0.38	0/2089	0.99	19/2847 (0.7%)
2	D	0.37	0/2077	1.09	14/2829 (0.5%)
All	All	0.38	0/5391	1.00	35/7355 (0.5%)

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
2	B	0	1
2	D	0	1
All	All	0	2

There are no bond length outliers.

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	532	SER	CA-C-N	21.98	147.32	119.84
2	D	532	SER	C-N-CA	21.98	147.32	119.84
2	B	532	SER	CA-C-N	8.66	147.79	127.00
2	B	532	SER	C-N-CA	8.66	147.79	127.00
1	A	11	PRO	CA-C-N	7.38	127.08	119.56
1	A	11	PRO	C-N-CA	7.38	127.08	119.56
2	B	438	GLY	CA-C-N	6.60	127.12	119.99
2	B	438	GLY	C-N-CA	6.60	127.12	119.99
2	B	470	MET	CA-C-N	-6.50	111.72	119.84
2	B	470	MET	C-N-CA	-6.50	111.72	119.84
2	B	450	THR	CA-C-N	6.28	127.69	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	450	THR	C-N-CA	6.28	127.69	119.84
2	B	425	LEU	CA-C-N	6.07	125.62	119.19
2	B	425	LEU	C-N-CA	6.07	125.62	119.19
2	D	353	SER	CA-C-N	5.95	125.71	119.76
2	D	353	SER	C-N-CA	5.95	125.71	119.76
2	B	523	ILE	N-CA-C	5.81	116.25	108.11
2	D	438	GLY	CA-C-N	5.72	126.17	119.99
2	D	438	GLY	C-N-CA	5.72	126.17	119.99
2	B	485	LEU	CA-C-N	5.71	125.66	119.78
2	B	485	LEU	C-N-CA	5.71	125.66	119.78
2	D	425	LEU	CA-C-N	5.54	125.15	119.56
2	D	425	LEU	C-N-CA	5.54	125.15	119.56
2	B	452	GLU	N-CA-C	5.53	118.27	109.81
2	D	523	ILE	N-CA-C	5.50	115.81	108.11
2	D	332	ASN	CA-C-N	5.46	126.67	119.84
2	D	332	ASN	C-N-CA	5.46	126.67	119.84
2	B	353	SER	CA-C-N	5.28	125.04	119.76
2	B	353	SER	C-N-CA	5.28	125.04	119.76
2	D	485	LEU	CA-C-N	5.27	125.21	119.78
2	D	485	LEU	C-N-CA	5.27	125.21	119.78
2	B	422	HIS	CA-C-N	5.23	124.89	119.56
2	B	422	HIS	C-N-CA	5.23	124.89	119.56
2	D	470	MET	N-CA-C	5.09	121.06	109.81
2	B	470	MET	N-CA-C	5.04	120.95	109.81

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	470	MET	Peptide
2	D	470	MET	Peptide

5.2 Too-close contacts ⓘ

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	1188	0	1015	53	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	2043	0	1789	114	0
2	D	2032	0	1769	100	0
3	C	50	0	43	2	0
4	A	14	0	13	0	0
4	B	14	0	13	0	0
4	D	14	0	13	0	0
All	All	5355	0	4655	255	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.

All (255) 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)
2:B:342:ARG:HG3	2:B:342:ARG:HH11	1.12	1.06
2:B:471:PRO:HD2	2:B:528:HIS:CE1	1.99	0.97
2:D:471:PRO:HD2	2:D:528:HIS:CE1	1.99	0.97
2:D:344:SER:HB3	2:D:347:ASP:HB2	1.48	0.95
2:B:344:SER:HB3	2:B:347:ASP:HB2	1.50	0.93
2:D:363:LEU:HD13	2:D:370:VAL:HG21	1.55	0.89
1:A:65:GLU:HB2	1:A:80:TYR:CZ	2.11	0.86
2:B:333:PRO:HG2	2:B:429:LEU:HD11	1.57	0.85
2:B:342:ARG:HG3	2:B:342:ARG:NH1	1.90	0.85
2:B:251:ILE:HD13	2:B:304:TRP:CD1	2.12	0.85
1:A:156:TRP:CZ3	2:D:333:PRO:HG2	2.12	0.84
2:B:539:ARG:HG2	2:B:540:ALA:H	1.46	0.80
1:A:65:GLU:HB2	1:A:80:TYR:CE2	2.19	0.78
2:B:251:ILE:HD13	2:B:304:TRP:CG	2.20	0.77
1:A:156:TRP:HZ3	2:D:333:PRO:HG2	1.48	0.76
2:D:539:ARG:HG2	2:D:540:ALA:H	1.51	0.75
2:D:513:ARG:HA	2:D:516:TRP:CG	2.22	0.74
1:A:41:HIS:HE1	1:A:63:SER:O	1.70	0.74
2:D:471:PRO:HD2	2:D:528:HIS:NE2	2.02	0.74
2:B:528:HIS:CD2	2:B:530:ALA:H	2.08	0.72
2:B:252:GLN:HG2	2:B:298:THR:HG23	1.71	0.71
2:B:242:ASP:HB3	2:B:246:HIS:H	1.54	0.71
2:B:471:PRO:HD2	2:B:528:HIS:NE2	2.04	0.71
1:A:77:GLU:HB3	1:A:78:PRO:CD	2.20	0.70
2:D:528:HIS:CD2	2:D:530:ALA:H	2.09	0.70
2:B:365:PRO:HA	2:B:397:LEU:HD13	1.75	0.69
2:B:259:TYR:HE1	2:B:292:SER:HA	1.57	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:1:NAG:H62	3:C:2:NAG:H82	1.75	0.69
2:D:440:ARG:HD3	2:D:440:ARG:N	2.07	0.68
2:D:346:PHE:HA	2:D:527:VAL:HG11	1.76	0.67
1:A:105:LEU:HB2	1:A:136:ILE:HB	1.76	0.67
1:A:130:TRP:HB3	1:A:134:HIS:CD2	2.30	0.66
2:B:492:THR:HG23	2:B:505:VAL:CG1	2.25	0.66
2:D:457:ARG:O	2:D:458:ASP:HB2	1.95	0.65
2:D:492:THR:HG23	2:D:505:VAL:CG1	2.26	0.64
2:B:475:SER:HB2	2:B:527:VAL:HB	1.79	0.64
2:B:342:ARG:HH11	2:B:342:ARG:CG	2.00	0.63
2:D:269:LEU:HB2	2:D:311:THR:HB	1.81	0.63
2:D:363:LEU:CD1	2:D:370:VAL:HG21	2.26	0.63
2:B:377:ALA:HB2	2:B:415:THR:HB	1.81	0.62
2:B:422:HIS:CG	2:B:423:PRO:HD2	2.34	0.62
2:B:448:PHE:CD2	2:D:448:PHE:CD2	2.86	0.62
2:B:425:LEU:HD13	2:B:429:LEU:HD13	1.81	0.61
2:D:371:GLN:HB2	2:D:421:THR:OG1	2.01	0.61
2:B:448:PHE:HD2	2:D:448:PHE:HD2	1.49	0.61
2:D:513:ARG:HD2	2:D:514:ALA:N	2.15	0.60
1:A:155:VAL:HG12	1:A:156:TRP:CD1	2.36	0.60
2:D:475:SER:HB2	2:D:527:VAL:HB	1.83	0.60
2:D:523:ILE:HD13	2:D:540:ALA:HB2	1.81	0.60
2:B:315:THR:HG23	2:B:320:THR:HG22	1.82	0.60
2:B:441:ALA:O	2:B:469:PHE:HA	2.02	0.60
1:A:133:ASN:HD22	1:A:134:HIS:H	1.47	0.60
2:B:374:TRP:HZ2	2:B:401:SER:O	1.85	0.59
2:B:493:THR:HG21	2:D:493:THR:HG21	1.84	0.59
1:A:77:GLU:HB3	1:A:78:PRO:HD3	1.82	0.59
2:B:464:CYS:HB2	2:B:478:TRP:CH2	2.38	0.59
2:B:446:TYR:CE1	2:D:451:PRO:HD2	2.37	0.59
2:D:425:LEU:HD13	2:D:429:LEU:HD13	1.84	0.59
2:D:479:LEU:O	2:D:480:HIS:HB2	2.02	0.58
2:B:448:PHE:CD2	2:D:448:PHE:HD2	2.20	0.58
2:B:478:TRP:CG	2:B:509:LEU:HD22	2.37	0.58
2:B:370:VAL:HG22	2:B:422:HIS:CD2	2.38	0.58
2:D:464:CYS:HB2	2:D:478:TRP:CH2	2.39	0.58
2:D:425:LEU:HD13	2:D:429:LEU:CD1	2.34	0.57
2:B:446:TYR:HB3	2:B:448:PHE:CE1	2.40	0.57
1:A:37:THR:O	1:A:49:THR:HA	2.04	0.57
2:B:331:SER:C	2:B:333:PRO:HD3	2.30	0.57
2:D:229:PHE:CD2	2:D:230:THR:N	2.72	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:315:THR:CG2	2:D:320:THR:HG22	2.34	0.56
2:D:450:THR:HG21	2:D:461:THR:HB	1.86	0.56
2:D:531:ALA:O	2:D:535:GLN:HA	2.06	0.56
1:A:11:PRO:HB2	1:A:13:TRP:CD1	2.41	0.56
2:D:363:LEU:HB2	2:D:397:LEU:HB3	1.87	0.55
1:A:40:PHE:CE1	1:A:45:LEU:HD12	2.41	0.55
2:D:266:ILE:HD12	2:D:295:SER:HB2	1.88	0.55
2:D:478:TRP:CG	2:D:509:LEU:HD22	2.42	0.55
2:B:444:GLU:O	2:B:466:ILE:HA	2.06	0.55
2:B:531:ALA:O	2:B:535:GLN:HA	2.06	0.55
2:D:422:HIS:CG	2:D:423:PRO:HD2	2.41	0.55
2:D:492:THR:HG23	2:D:505:VAL:HG11	1.88	0.55
1:A:110:TRP:O	1:A:113:TRP:HB2	2.07	0.55
1:A:77:GLU:CB	1:A:78:PRO:HD3	2.38	0.55
2:B:250:THR:CG2	2:B:298:THR:HG22	2.36	0.55
2:B:333:PRO:HG2	2:B:429:LEU:CD1	2.32	0.55
2:B:448:PHE:HD2	2:D:448:PHE:CD2	2.25	0.55
1:A:104:PHE:O	1:A:105:LEU:HD23	2.07	0.54
1:A:156:TRP:O	1:A:157:GLN:HB2	2.07	0.54
2:D:311:THR:HG23	2:D:324:SER:OG	2.07	0.54
2:D:513:ARG:O	2:D:516:TRP:HB2	2.07	0.54
2:B:492:THR:HG23	2:B:505:VAL:HG11	1.88	0.54
2:B:338:ALA:HA	2:B:359:LEU:O	2.07	0.54
2:B:387:ARG:HG3	2:B:387:ARG:O	2.07	0.54
2:B:269:LEU:HB2	2:B:311:THR:HB	1.90	0.54
2:B:539:ARG:HG2	2:B:540:ALA:N	2.21	0.54
2:B:425:LEU:HD13	2:B:429:LEU:CD1	2.38	0.54
2:D:446:TYR:HB3	2:D:448:PHE:CE1	2.43	0.53
1:A:133:ASN:ND2	1:A:134:HIS:H	2.06	0.53
2:D:359:LEU:HD12	2:D:360:VAL:N	2.23	0.53
2:B:446:TYR:CD1	2:D:451:PRO:HD2	2.43	0.53
2:B:434:THR:HG22	2:B:435:LYS:N	2.22	0.53
2:D:339:TYR:CD1	2:D:339:TYR:N	2.77	0.52
1:A:119:ILE:HD12	1:A:152:THR:CG2	2.40	0.52
2:B:311:THR:HG23	2:B:324:SER:OG	2.10	0.51
2:D:251:ILE:HD13	2:D:304:TRP:CD1	2.45	0.51
2:B:301:GLN:O	2:B:305:LEU:HG	2.11	0.51
2:B:465:LEU:HG	2:B:465:LEU:O	2.09	0.51
2:B:304:TRP:CZ2	2:B:327:LYS:HA	2.46	0.51
2:B:251:ILE:CD1	2:B:304:TRP:CD1	2.90	0.50
2:D:229:PHE:HD2	2:D:230:THR:H	1.59	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:367:LYS:C	2:B:369:THR:H	2.19	0.50
2:D:376:ARG:HD3	2:D:380:LYS:O	2.12	0.50
2:B:434:THR:CG2	2:B:435:LYS:N	2.74	0.50
2:D:299:LEU:HD21	2:D:310:TYR:CE2	2.46	0.50
2:B:250:THR:HG22	2:B:298:THR:HG22	1.94	0.49
2:B:445:VAL:CG1	2:B:464:CYS:SG	3.00	0.49
2:D:345:PRO:HD2	2:D:475:SER:HG	1.78	0.49
2:D:479:LEU:HA	2:D:484:GLN:HA	1.94	0.49
1:A:41:HIS:NE2	3:C:1:NAG:H82	2.28	0.49
2:B:333:PRO:C	2:B:335:GLY:H	2.19	0.49
2:B:342:ARG:NH1	2:B:342:ARG:CG	2.67	0.49
2:D:422:HIS:HB3	2:D:425:LEU:HG	1.95	0.49
2:B:344:SER:HB3	2:B:347:ASP:CB	2.31	0.49
2:B:371:GLN:HB2	2:B:421:THR:OG1	2.13	0.49
1:A:15:ARG:HB3	1:A:84:PHE:HE1	1.78	0.48
2:B:331:SER:O	2:B:333:PRO:HD3	2.13	0.48
2:B:236:ILE:HA	2:B:253(A):LEU:O	2.13	0.48
2:D:338:ALA:HA	2:D:359:LEU:O	2.13	0.48
2:D:345:PRO:HB2	2:D:527:VAL:O	2.13	0.48
1:A:101:GLN:O	1:A:141:ALA:HB2	2.13	0.48
2:D:256:VAL:HG11	2:D:264:ILE:CD1	2.43	0.48
2:B:355:THR:HG22	2:B:404:PRO:HA	1.96	0.48
2:B:377:ALA:CB	2:B:415:THR:HB	2.44	0.48
2:B:240:SER:HB2	2:D:325:THR:CB	2.44	0.48
2:B:303:HIS:HD2	2:B:310:TYR:OH	1.96	0.48
1:A:133:ASN:HD22	1:A:134:HIS:N	2.12	0.47
2:B:386:THR:CG2	2:B:387:ARG:N	2.77	0.47
2:B:478:TRP:CB	2:B:509:LEU:HD22	2.45	0.47
2:B:446:TYR:O	2:B:464:CYS:HA	2.14	0.47
2:D:231:PRO:HB3	2:D:321:PHE:CZ	2.49	0.47
1:A:19:GLY:O	1:A:57:ASN:HA	2.14	0.47
2:B:333:PRO:C	2:B:335:GLY:N	2.72	0.47
2:B:471:PRO:HD2	2:B:528:HIS:HE1	1.68	0.47
2:D:478:TRP:CB	2:D:509:LEU:HD22	2.45	0.47
2:B:339:TYR:CD1	2:B:339:TYR:N	2.82	0.47
1:A:37:THR:HG22	1:A:38:LYS:N	2.30	0.47
2:D:345:PRO:HD2	2:D:475:SER:OG	2.14	0.47
1:A:128:LYS:HG2	1:A:129:TYR:N	2.29	0.46
2:B:446:TYR:CB	2:B:448:PHE:HE1	2.29	0.46
2:D:370:VAL:HG22	2:D:422:HIS:HD2	1.80	0.46
1:A:86:ASP:OD1	1:A:87:TRP:N	2.44	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:300:SER:HB2	2:B:303:HIS:H	1.80	0.46
1:A:146:SER:HB3	1:A:168:THR:HG22	1.97	0.46
2:D:474:ILE:HG12	2:D:475:SER:N	2.31	0.46
2:B:370:VAL:HG22	2:B:422:HIS:HD2	1.79	0.45
2:D:346:PHE:CE1	2:D:350:ILE:HG21	2.51	0.45
2:D:231:PRO:HB3	2:D:321:PHE:CE2	2.51	0.45
1:A:98:MET:O	1:A:99:GLU:C	2.60	0.45
1:A:155:VAL:HB	1:A:160:TYR:CE2	2.52	0.45
2:D:374:TRP:HZ2	2:D:401:SER:O	1.98	0.45
1:A:93:SER:CB	1:A:104:PHE:H	2.29	0.45
2:D:419:ARG:HD2	2:D:430:MET:SD	2.57	0.45
2:B:336:VAL:HG12	2:B:337:SER:N	2.32	0.45
2:B:422:HIS:HB3	2:B:425:LEU:HG	1.99	0.45
1:A:133:ASN:ND2	1:A:134:HIS:N	2.65	0.45
2:B:388:LYS:HB2	2:B:400:THR:HB	1.98	0.45
2:B:472:GLU:HG3	2:B:503:PHE:CE2	2.51	0.45
2:D:513:ARG:HD2	2:D:513:ARG:C	2.42	0.45
2:B:259:TYR:HE1	2:B:292:SER:CA	2.28	0.44
2:B:259:TYR:CE1	2:B:292:SER:HA	2.43	0.44
1:A:113:TRP:CD1	1:A:156:TRP:HZ2	2.35	0.44
2:B:346:PHE:CE1	2:B:350:ILE:HG21	2.52	0.44
2:B:405:VAL:HG13	2:B:416:TYR:OH	2.16	0.44
2:B:448:PHE:HB3	2:D:448:PHE:CE2	2.52	0.44
2:D:471:PRO:HD2	2:D:528:HIS:HE1	1.71	0.44
2:B:270:GLU:HB3	2:B:275:MET:HG2	2.00	0.44
2:D:363:LEU:CB	2:D:397:LEU:HB3	2.47	0.44
1:A:103:LEU:HD23	1:A:138:ILE:HD13	1.99	0.44
2:B:359:LEU:HD12	2:B:360:VAL:H	1.82	0.44
1:A:59:LYS:O	1:A:62:ASP:HB2	2.18	0.44
2:B:299:LEU:HD21	2:B:310:TYR:CE2	2.53	0.44
2:D:370:VAL:HG22	2:D:422:HIS:CD2	2.52	0.44
2:B:506:PHE:CD1	2:D:506:PHE:CD1	3.05	0.44
2:D:251:ILE:CD1	2:D:304:TRP:CD1	3.01	0.44
2:B:345:PRO:HB2	2:B:475:SER:OG	2.18	0.44
2:D:315:THR:HG23	2:D:320:THR:HG22	1.98	0.44
2:D:469:PHE:HB2	2:D:528:HIS:CE1	2.53	0.43
2:D:474:ILE:HG12	2:D:475:SER:H	1.83	0.43
2:D:441:ALA:O	2:D:469:PHE:HA	2.18	0.43
2:B:478:TRP:HA	2:B:523:ILE:O	2.18	0.43
2:D:359:LEU:HD12	2:D:360:VAL:H	1.83	0.43
1:A:77:GLU:CB	1:A:78:PRO:CD	2.91	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:440:ARG:HD2	2:B:471:PRO:HG3	2.00	0.43
2:D:249:PRO:O	2:D:301:GLN:HB2	2.18	0.43
1:A:169:VAL:HG12	1:A:170:ILE:N	2.33	0.43
2:D:446:TYR:CB	2:D:448:PHE:HE1	2.32	0.43
1:A:41:HIS:C	1:A:43:GLY:H	2.27	0.43
1:A:93:SER:HB2	1:A:104:PHE:H	1.83	0.43
2:B:525:ARG:HA	2:B:538:GLN:HB2	2.00	0.43
2:B:534:SER:O	2:B:535:GLN:HB2	2.19	0.43
2:D:534:SER:O	2:D:535:GLN:HB2	2.19	0.43
2:B:346:PHE:N	2:B:475:SER:OG	2.51	0.43
2:D:256:VAL:HG11	2:D:264:ILE:HD13	2.00	0.43
1:A:41:HIS:CE1	1:A:63:SER:O	2.60	0.43
2:B:539:ARG:CG	2:B:540:ALA:H	2.24	0.43
2:D:269:LEU:HA	2:D:273:GLN:O	2.19	0.43
2:B:346:PHE:C	2:B:348:LEU:H	2.26	0.42
2:D:344:SER:HB3	2:D:347:ASP:CB	2.33	0.42
2:D:400:THR:HG22	2:D:401:SER:N	2.33	0.42
1:A:136:ILE:N	1:A:136:ILE:HD12	2.34	0.42
2:B:333:PRO:CG	2:B:429:LEU:HD11	2.39	0.42
2:B:242:ASP:CB	2:B:246:HIS:HB2	2.48	0.42
2:B:303:HIS:CE1	2:B:308:ARG:NH1	2.87	0.42
2:B:422:HIS:CE1	2:B:423:PRO:HG2	2.54	0.42
2:D:386:THR:O	2:D:386:THR:HG22	2.19	0.42
2:D:523:ILE:CD1	2:D:540:ALA:HB2	2.48	0.42
1:A:17:PHE:O	1:A:20:GLU:HB2	2.19	0.42
1:A:88:LEU:HD21	1:A:153:GLY:HA3	2.00	0.42
1:A:118:VAL:HA	1:A:152:THR:O	2.20	0.42
1:A:119:ILE:HD12	1:A:152:THR:HG21	2.01	0.42
1:A:165:LEU:HD12	1:A:166:ASN:N	2.35	0.42
2:B:469:PHE:HB2	2:B:528:HIS:CE1	2.55	0.42
2:D:229:PHE:CE2	2:D:231:PRO:HD3	2.55	0.42
2:B:241:CYS:O	2:D:325:THR:HA	2.19	0.42
2:B:345:PRO:HG2	2:B:474:ILE:C	2.44	0.42
2:B:359:LEU:HD12	2:B:360:VAL:N	2.35	0.42
2:D:392:GLN:HE21	2:D:393:ARG:H	1.67	0.42
2:B:446:TYR:CB	2:B:448:PHE:CE1	3.02	0.41
2:D:512:THR:C	2:D:514:ALA:N	2.78	0.41
1:A:9:LEU:HD23	1:A:22:VAL:HG21	2.01	0.41
1:A:119:ILE:HA	1:A:128:LYS:O	2.20	0.41
2:B:269:LEU:HD11	2:B:313:GLN:NE2	2.35	0.41
2:D:450:THR:CG2	2:D:461:THR:HB	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:97:VAL:HG23	1:A:169:VAL:HG22	2.02	0.41
2:D:479:LEU:HD21	2:D:525:ARG:NH2	2.35	0.41
2:B:370:VAL:HG11	2:B:399:VAL:HG21	2.02	0.41
2:D:251:ILE:HD13	2:D:304:TRP:CG	2.56	0.41
2:D:511:VAL:HG21	2:D:522:PHE:CE2	2.55	0.41
1:A:9:LEU:HD23	1:A:22:VAL:CG2	2.51	0.41
2:D:388:LYS:O	2:D:400:THR:HB	2.21	0.41
2:D:478:TRP:HA	2:D:523:ILE:O	2.21	0.41
2:B:260:THR:HB	2:B:316:TYR:OH	2.20	0.41
1:A:90:LEU:HD23	1:A:165:LEU:HB3	2.03	0.40
2:B:242:ASP:HB3	2:B:246:HIS:HB2	2.03	0.40
2:B:300:SER:HB2	2:B:303:HIS:HB2	2.03	0.40
2:B:464:CYS:HB2	2:B:478:TRP:CZ2	2.56	0.40
2:D:539:ARG:HG2	2:D:540:ALA:N	2.27	0.40
2:B:242:ASP:C	2:B:244:GLY:N	2.76	0.40
2:D:537:VAL:HG12	2:D:538:GLN:N	2.35	0.40
1:A:15:ARG:HB3	1:A:84:PHE:CE1	2.55	0.40
2:B:400:THR:HG22	2:B:401:SER:N	2.35	0.40
2:D:504:PHE:CD1	2:D:504:PHE:C	3.00	0.40
2:D:528:HIS:HD2	2:D:530:ALA:H	1.64	0.40
2:B:537:VAL:O	2:B:538:GLN:HB3	2.21	0.40
2:B:367:LYS:O	2:B:369:THR:N	2.53	0.40
2:D:303:HIS:O	2:D:306:SER:HB2	2.20	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	152/188 (81%)	135 (89%)	13 (9%)	4 (3%)	4	21
2	B	275/327 (84%)	246 (90%)	24 (9%)	5 (2%)	6	26

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	D	274/327 (84%)	244 (89%)	23 (8%)	7 (3%)	4	21
All	All	701/842 (83%)	625 (89%)	60 (9%)	16 (2%)	5	23

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	533	PRO
1	A	57	ASN
1	A	99	GLU
2	B	350	ILE
2	D	458	ASP
2	B	242	ASP
2	D	350	ILE
1	A	77	GLU
2	D	230	THR
2	D	333	PRO
1	A	141	ALA
2	B	471	PRO
2	D	471	PRO
2	B	533	PRO
2	B	470	MET
2	D	470	MET

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	109/167 (65%)	100 (92%)	9 (8%)	10	35
2	B	188/289 (65%)	172 (92%)	16 (8%)	10	33
2	D	182/289 (63%)	167 (92%)	15 (8%)	10	35
All	All	479/745 (64%)	439 (92%)	40 (8%)	10	34

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	VAL
1	A	9	LEU
1	A	26	CYS
1	A	49	THR
1	A	51	SER
1	A	97	VAL
1	A	119	ILE
1	A	138	ILE
1	A	152	THR
2	B	263	THR
2	B	265	GLN
2	B	271	ASP
2	B	298	THR
2	B	300	SER
2	B	315	THR
2	B	337	SER
2	B	342	ARG
2	B	357	THR
2	B	386	THR
2	B	387	ARG
2	B	390	GLU
2	B	403	LEU
2	B	436	THR
2	B	491	SER
2	B	492	THR
2	D	233	THR
2	D	298	THR
2	D	315	THR
2	D	355	THR
2	D	357	THR
2	D	370	VAL
2	D	382	VAL
2	D	386	THR
2	D	417	GLN
2	D	436	THR
2	D	440	ARG
2	D	450	THR
2	D	474	ILE
2	D	492	THR
2	D	513	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	10	ASN
1	A	14	ASN
1	A	41	HIS
1	A	133	ASN
2	B	303	HIS
2	B	313	GLN
2	B	371	GLN
2	B	384	HIS
2	B	468	ASN
2	B	490	HIS
2	B	528	HIS
2	D	246	HIS
2	D	392	GLN
2	D	477	GLN
2	D	528	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	$\# Z > 2$
3	NAG	C	1	1,3	14,14,15	0.60	0	17,19,21	0.93	1 (5%)
3	NAG	C	2	3	14,14,15	0.56	0	17,19,21	1.34	1 (5%)
3	BMA	C	3	3	11,11,12	0.89	1 (9%)	15,15,17	0.79	0
3	MAN	C	4	3	11,11,12	0.74	0	15,15,17	0.97	1 (6%)

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	C	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	C	2	3	-	0/6/23/26	0/1/1/1
3	BMA	C	3	3	-	1/2/19/22	0/1/1/1
3	MAN	C	4	3	-	2/2/19/22	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	3	BMA	O5-C5	-2.03	1.39	1.43

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	2	NAG	C2-N2-C7	-4.00	117.54	122.90
3	C	1	NAG	C2-N2-C7	-2.47	119.59	122.90
3	C	4	MAN	C1-C2-C3	2.08	112.68	109.64

There are no chirality outliers.

All (3) torsion outliers are listed below:

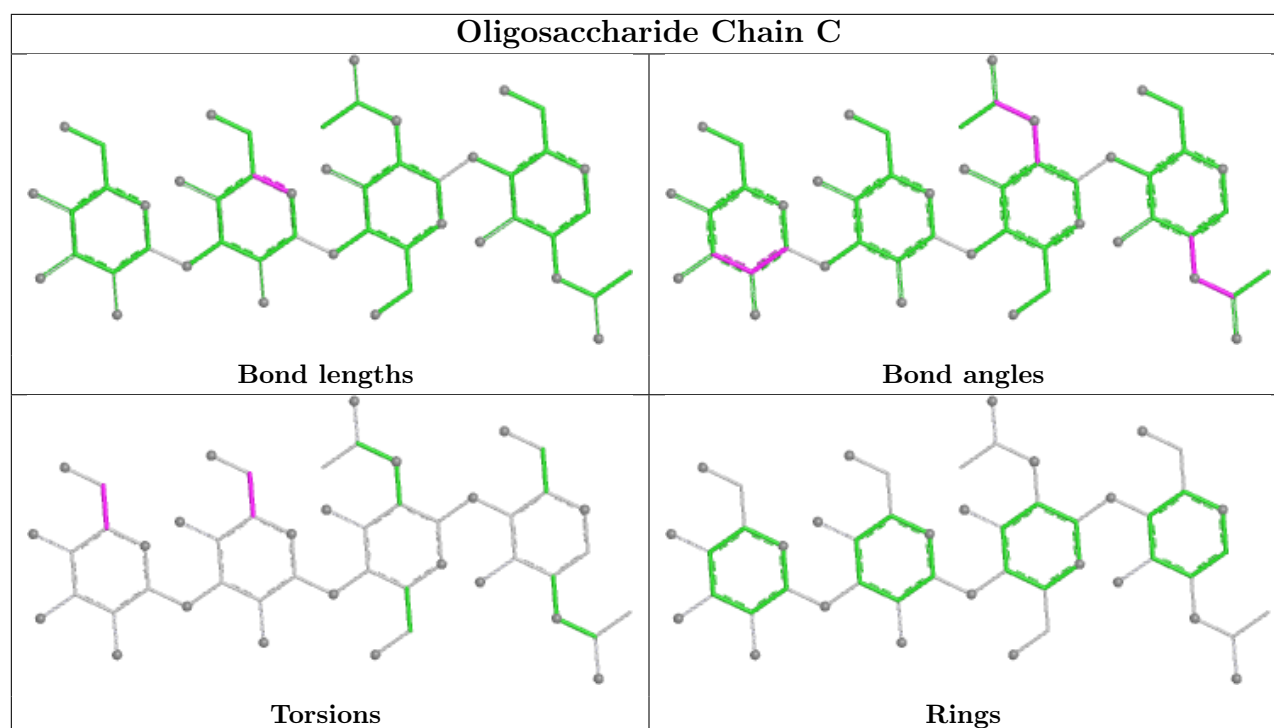
Mol	Chain	Res	Type	Atoms
3	C	4	MAN	O5-C5-C6-O6
3	C	4	MAN	C4-C5-C6-O6
3	C	3	BMA	O5-C5-C6-O6

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	2	NAG	1	0
3	C	1	NAG	2	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](#)

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	$\# Z > 2$
4	NAG	B	601	2	14,14,15	0.61	0	17,19,21	0.76	0
4	NAG	A	201	1	14,14,15	0.48	0	17,19,21	1.65	3 (17%)
4	NAG	D	601	2	14,14,15	0.43	0	17,19,21	1.84	3 (17%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	B	601	2	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	A	201	1	-	2/6/23/26	0/1/1/1
4	NAG	D	601	2	-	3/6/23/26	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(°)
4	D	601	NAG	C1-O5-C5	5.18	119.12	112.19
4	A	201	NAG	C1-O5-C5	4.86	118.69	112.19
4	D	601	NAG	C4-C3-C2	-3.69	105.61	111.02
4	A	201	NAG	C3-C4-C5	2.71	115.15	110.23
4	D	601	NAG	C2-N2-C7	-2.19	119.97	122.90
4	A	201	NAG	O5-C5-C4	2.17	116.10	110.83

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	201	NAG	C8-C7-N2-C2
4	A	201	NAG	O7-C7-N2-C2
4	B	601	NAG	C8-C7-N2-C2
4	B	601	NAG	O7-C7-N2-C2
4	D	601	NAG	C8-C7-N2-C2
4	D	601	NAG	O7-C7-N2-C2
4	D	601	NAG	O5-C5-C6-O6

There are no ring outliers.

No monomer is involved in short contacts.

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

6.1 Protein, DNA and RNA chains

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	158/188 (84%)	-0.02	2 (1%) 75 61	28, 64, 114, 146	0
2	B	287/327 (87%)	0.26	10 (3%) 47 34	33, 76, 120, 159	0
2	D	290/327 (88%)	0.40	17 (5%) 28 22	32, 81, 127, 154	0
All	All	735/842 (87%)	0.26	29 (3%) 43 31	28, 76, 124, 159	0

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	473	ASP	4.7
2	B	368	GLY	3.2
2	D	243	GLY	3.2
2	D	478	TRP	3.0
2	D	424	HIS	2.9
2	D	534	SER	2.8
2	D	502	GLY	2.7
2	D	321	PHE	2.7
1	A	51	SER	2.6
2	B	309	THR	2.6
2	D	538	GLN	2.6
2	B	243	GLY	2.6
2	D	319	HIS	2.5
1	A	112	ASN	2.5
2	B	458	ASP	2.5
2	D	498	THR	2.5
2	D	468	ASN	2.4
2	D	249	PRO	2.4
2	B	366	SER	2.4
2	B	477	GLN	2.3
2	D	458	ASP	2.2
2	D	325	THR	2.2
2	D	261	PRO	2.2

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Mol	Chain	Res	Type	RSRZ
2	B	453	TRP	2.1
2	D	397	LEU	2.1
2	B	525	ARG	2.1
2	D	290	LEU	2.1
2	B	271	ASP	2.0
2	B	244	GLY	2.0

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

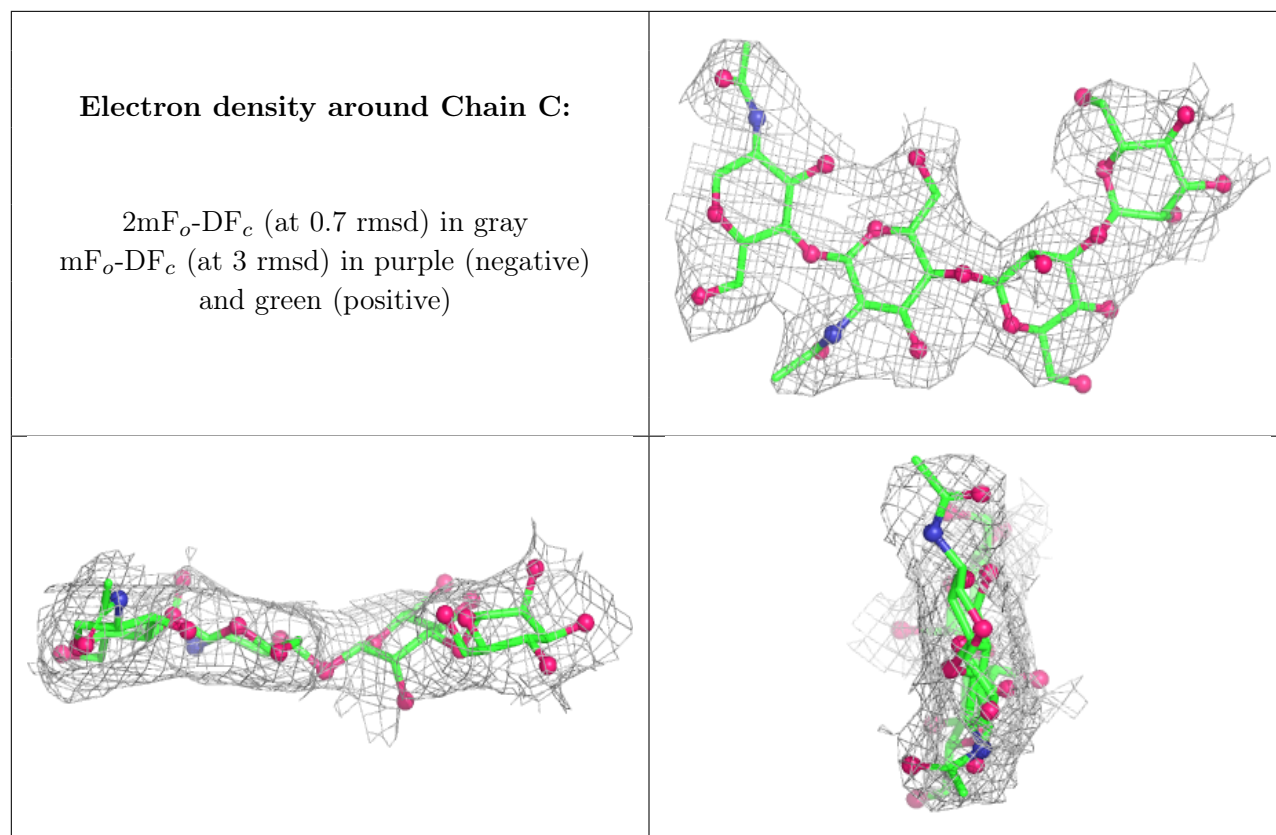
There are no non-standard protein/DNA/RNA residues in this entry.

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
3	NAG	C	1	14/15	-	-	44,44,44,44	0
3	NAG	C	2	14/15	-	-	56,56,56,56	0
3	BMA	C	3	11/12	-	-	104,104,104,104	0
3	MAN	C	4	11/12	-	-	101,101,101,101	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($\text{\AA}^2$)	Q<0.9
4	NAG	D	601	14/15	0.83	0.13	90,90,90,90	0
4	NAG	A	201	14/15	0.88	0.07	83,83,83,83	0
4	NAG	B	601	14/15	0.90	0.10	65,65,65,65	0

6.5 Other polymers [i](#)

There are no such residues in this entry.