



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 09:47 PM UTC

PDB ID : 8SDH / pdb_00008sdh
Title : Crystal structure of SARS-CoV-2 receptor binding domain in complex with neutralizing antibody CC25.56
Authors : Yuan, M.; Wilson, I.A.
Deposited on : 2023-04-06
Resolution : 2.84 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

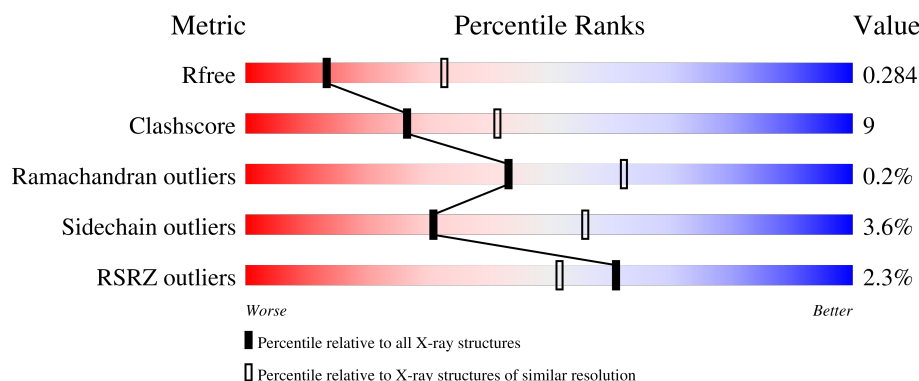
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.84 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	1520 (2.86-2.82)
Clashscore	190562	1559 (2.86-2.82)
Ramachandran outliers	187476	1517 (2.86-2.82)
Sidechain outliers	187428	1518 (2.86-2.82)
RSRZ outliers	180081	1521 (2.86-2.82)

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	205	6% 73% 21% 5%
1	B	205	3% 71% 24% 5%
2	C	223	74% 24% .
2	H	223	% 83% 13% ..
3	D	214	2% 78% 17% ..

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Mol	Chain	Length	Quality of chain
3	L	214	2% 79% 19%
4	E	2	100%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 9606 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 Spike protein S1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	195	Total	C	N	O	S	0	0	0
			1541	988	257	288	8			
1	B	195	Total	C	N	O	S	0	0	0
			1535	984	256	287	8			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	531	GLY	-	expression tag	UNP P0DTC2
A	532	HIS	-	expression tag	UNP P0DTC2
A	533	HIS	-	expression tag	UNP P0DTC2
A	534	HIS	-	expression tag	UNP P0DTC2
A	535	HIS	-	expression tag	UNP P0DTC2
A	536	HIS	-	expression tag	UNP P0DTC2
A	537	HIS	-	expression tag	UNP P0DTC2
B	531	GLY	-	expression tag	UNP P0DTC2
B	532	HIS	-	expression tag	UNP P0DTC2
B	533	HIS	-	expression tag	UNP P0DTC2
B	534	HIS	-	expression tag	UNP P0DTC2
B	535	HIS	-	expression tag	UNP P0DTC2
B	536	HIS	-	expression tag	UNP P0DTC2
B	537	HIS	-	expression tag	UNP P0DTC2

- Molecule 2 is a protein called Neutralizing antibody CC25.56 Heavy Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	216	Total	C	N	O	S	0	0	0
			1626	1037	271	310	8			
2	C	222	Total	C	N	O	S	0	0	0
			1664	1058	278	320	8			

- Molecule 3 is a protein called Neutralizing antibody CC25.56 Light Chain.

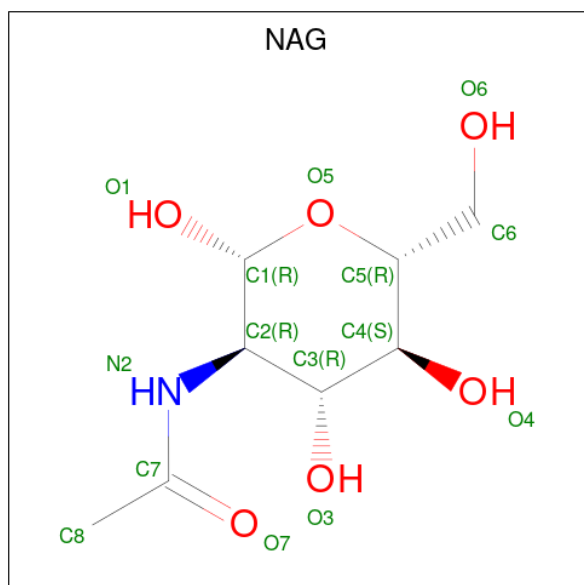
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	L	210	Total	C	N	O	S	0	0	0
			1573	986	265	318	4			
3	D	210	Total	C	N	O	S	0	1	0
			1584	995	266	319	4			

- Molecule 4 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
4	E	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



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

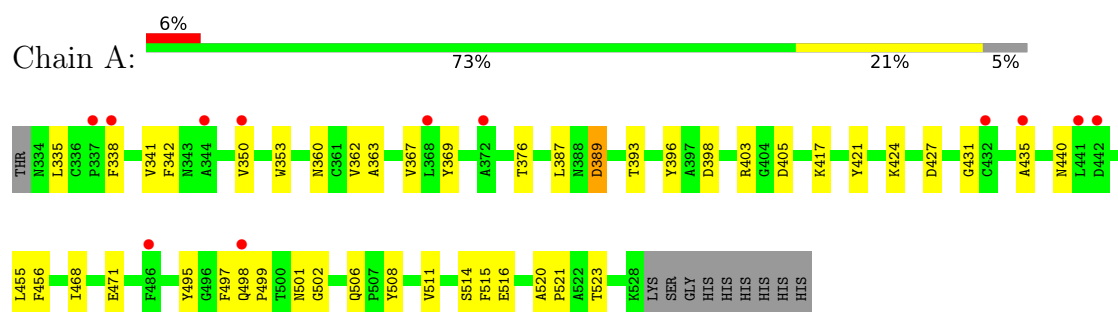
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	7	Total 7	O 7	0	0
6	B	3	Total 3	O 3	0	0
6	H	6	Total 6	O 6	0	0
6	L	7	Total 7	O 7	0	0
6	C	10	Total 10	O 10	0	0
6	D	8	Total 8	O 8	0	0

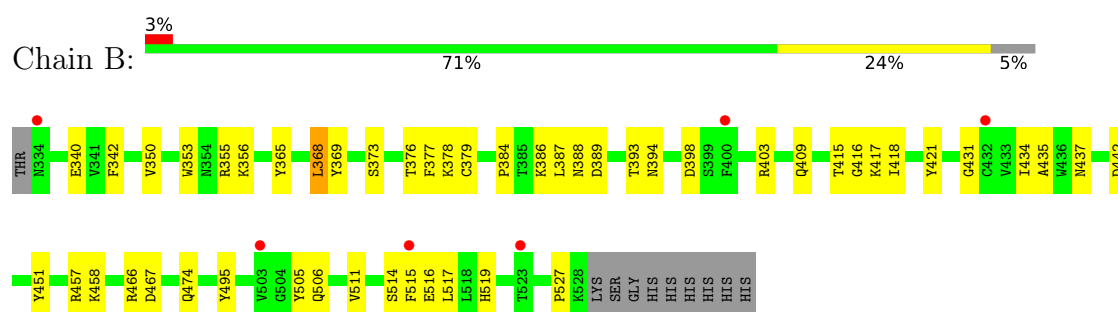
3 Residue-property plots [i](#)

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.

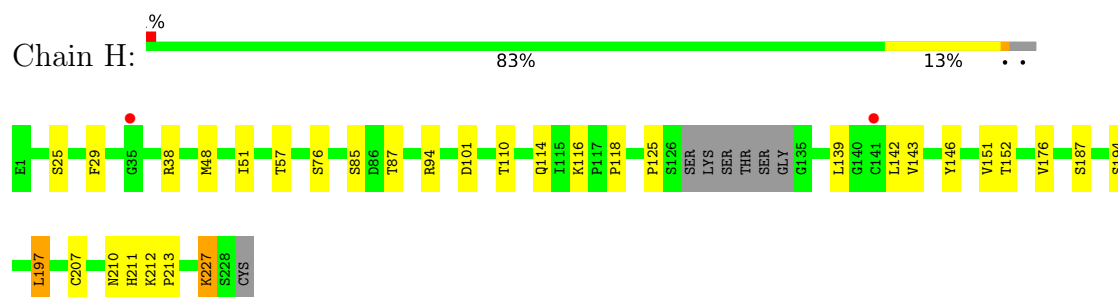
- Molecule 1: Spike protein S1



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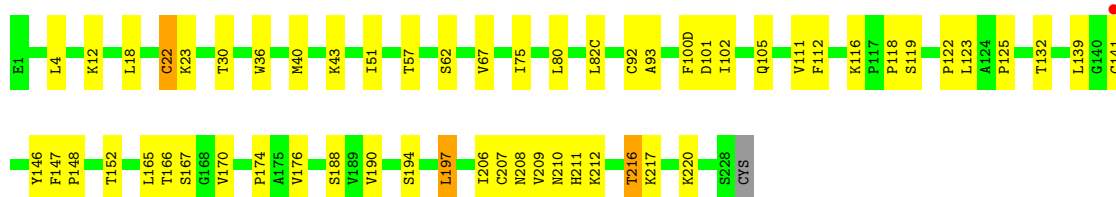


- Molecule 2: Neutralizing antibody CC25.56 Heavy Chain

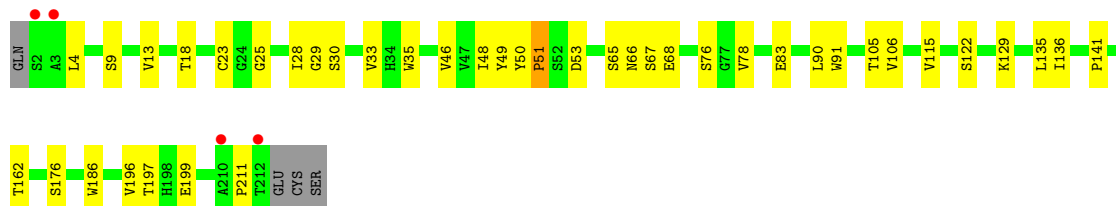
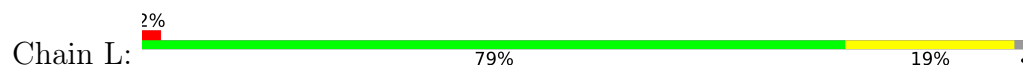


- Molecule 2: Neutralizing antibody CC25.56 Heavy Chain

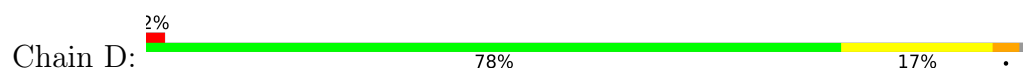




• Molecule 3: Neutralizing antibody CC25.56 Light Chain



• Molecule 3: Neutralizing antibody CC25.56 Light Chain



• Molecule 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	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	171.05Å 105.59Å 98.99Å 90.00° 117.31° 90.00°	Depositor
Resolution (Å)	47.69 – 2.84 47.69 – 2.84	Depositor EDS
% Data completeness (in resolution range)	98.8 (47.69-2.84) 98.8 (47.69-2.84)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.48 (at 2.86Å)	Xtriage
Refinement program	PHENIX (1.19.2_4158: ???)	Depositor
R, R_{free}	0.246 , 0.283 0.246 , 0.284	Depositor DCC
R_{free} test set	1915 reflections (5.17%)	wwPDB-VP
Wilson B-factor (Å ²)	54.8	Xtriage
Anisotropy	0.635	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 35.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	9606	wwPDB-VP
Average B, all atoms (Å ²)	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.15% 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

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	0.23	1/1585 (0.1%)	0.73	3/2158 (0.1%)
1	B	0.13	0/1579	0.33	0/2150
2	C	0.27	0/1708	0.58	5/2325 (0.2%)
2	H	0.19	0/1669	0.43	1/2272 (0.0%)
3	D	0.15	0/1627	0.43	2/2227 (0.1%)
3	L	0.15	0/1615	0.39	0/2211
All	All	0.19	1/9783 (0.0%)	0.50	11/13343 (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
3	D	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	497	PHE	C-N	7.14	1.45	1.33

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	497	PHE	CA-C-N	19.11	148.17	121.61
1	A	497	PHE	C-N-CA	19.11	148.17	121.61
1	A	497	PHE	O-C-N	-12.98	107.06	123.02
2	C	206	ILE	CA-CB-CG1	9.57	126.66	110.40
3	D	50	TYR	C-N-CD	-6.39	98.82	125.00
2	H	227	LYS	CD-CE-NZ	-6.29	91.78	111.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	206	ILE	CG1-CB-CG2	-6.09	92.42	110.70
3	D	129	LYS	CA-CB-CG	5.28	124.65	114.10
2	C	216	THR	CA-C-N	-5.17	113.47	122.37
2	C	216	THR	C-N-CA	-5.17	113.47	122.37
2	C	206	ILE	CB-CA-C	-5.00	103.66	110.96

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	D	50	TYR	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	1541	0	1454	30	0
1	B	1535	0	1441	37	0
2	C	1664	0	1642	41	0
2	H	1626	0	1603	16	0
3	D	1584	0	1521	29	0
3	L	1573	0	1513	32	0
4	E	28	0	25	0	0
5	B	14	0	13	0	0
6	A	7	0	0	0	0
6	B	3	0	0	0	0
6	C	10	0	0	0	0
6	D	8	0	0	0	0
6	H	6	0	0	0	0
6	L	7	0	0	0	0
All	All	9606	0	9212	174	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 9.

All (174) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:29:GLY:HA2	3:L:66:ASN:HD21	1.36	0.90
2:C:210:ASN:HB2	2:C:217:LYS:HZ3	1.43	0.82
1:A:521:PRO:HG3	1:B:409:GLN:HG3	1.61	0.80
3:L:66:ASN:HD22	3:L:67:SER:N	1.83	0.75
2:C:210:ASN:HB2	2:C:217:LYS:HD2	1.69	0.74
3:D:181:LEU:HD12	3:D:185:GLN:HG3	1.71	0.72
1:B:340:GLU:OE1	1:B:356:LYS:NZ	2.22	0.72
3:D:130:ALA:HB3	3:D:181:LEU:O	1.89	0.71
2:C:51:ILE:HD12	2:C:57:THR:HG22	1.71	0.70
2:C:139:LEU:HD11	2:C:197:LEU:HD11	1.75	0.69
2:H:227:LYS:HZ1	3:L:122:SER:CB	2.06	0.67
3:L:29:GLY:HA2	3:L:66:ASN:ND2	2.09	0.66
1:A:341:VAL:HG23	1:A:342:PHE:HD1	1.61	0.66
2:H:152:THR:HG22	2:H:210:ASN:HB3	1.78	0.66
2:C:125:PRO:HB3	2:C:139:LEU:HD13	1.78	0.66
1:B:457:ARG:NH1	1:B:467:ASP:OD2	2.30	0.65
1:B:458:LYS:NZ	1:B:474:GLN:O	2.29	0.65
2:C:18:LEU:HD12	2:C:82(C):LEU:HD11	1.79	0.63
2:C:209:VAL:O	2:C:217:LYS:HG3	1.98	0.63
1:A:360:ASN:H	1:A:523:THR:HB	1.63	0.62
1:B:434:ILE:HB	1:B:511:VAL:HG12	1.80	0.62
3:L:105:THR:HG21	3:L:141:PRO:HB3	1.81	0.62
3:D:189:HIS:HD1	3:D:192:TYR:HH	1.47	0.61
1:B:365:TYR:HA	1:B:368:LEU:HD12	1.83	0.61
3:L:136:ILE:HG21	3:L:196:VAL:HG11	1.84	0.60
1:A:369:TYR:O	1:A:369:TYR:HD1	1.85	0.59
3:L:66:ASN:HD22	3:L:67:SER:H	1.49	0.58
2:C:112:PHE:HB2	2:C:147:PHE:CZ	2.38	0.58
3:L:83:GLU:HG3	3:L:106:VAL:HG23	1.84	0.58
2:H:212:LYS:HG2	2:H:213:PRO:HD3	1.84	0.58
2:C:23:LYS:NZ	2:C:75:ILE:O	2.36	0.58
3:D:128:ASN:O	3:D:128:ASN:ND2	2.37	0.57
3:D:195:GLN:NE2	3:D:206:GLU:OE2	2.19	0.57
3:D:127:ALA:HB1	3:D:129:LYS:NZ	2.19	0.57
3:L:90:LEU:HD12	3:L:91:TRP:H	1.69	0.57
2:C:194:SER:HA	2:C:197:LEU:HD23	1.85	0.57
2:C:176:VAL:HB	3:D:162:THR:HG22	1.86	0.57
3:L:90:LEU:HD12	3:L:91:TRP:N	2.20	0.57
1:B:355:ARG:HD3	1:B:398:ASP:OD1	2.05	0.57
1:B:342:PHE:CZ	1:B:511:VAL:HG11	2.40	0.56
2:H:114:GLN:HG2	2:H:116:LYS:HB2	1.86	0.56
2:C:101:ASP:HA	3:D:46:VAL:HG21	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:125:PRO:HD3	2:H:139:LEU:HD13	1.86	0.56
3:L:4:LEU:HD22	3:L:23:CYS:SG	2.45	0.56
1:B:431:GLY:HA2	1:B:515:PHE:CD2	2.40	0.56
3:L:48:ILE:HG22	3:L:49:TYR:O	2.06	0.55
2:C:118:PRO:HB3	2:C:146:TYR:HB3	1.86	0.55
2:H:227:LYS:NZ	3:L:122:SER:HB3	2.21	0.55
3:D:127:ALA:HB1	3:D:129:LYS:HZ2	1.72	0.55
1:B:388:ASN:HB2	1:B:527:PRO:HD2	1.89	0.54
3:D:119:PRO:HA	3:D:132:LEU:HD13	1.88	0.54
1:B:365:TYR:CD2	1:B:387:LEU:HD22	2.42	0.54
1:B:393:THR:N	1:B:516:GLU:O	2.40	0.54
2:H:227:LYS:NZ	3:L:122:SER:CB	2.69	0.54
3:L:18:THR:HG23	3:L:76:SER:HA	1.90	0.54
3:D:128:ASN:C	3:D:129:LYS:HD3	2.31	0.54
1:B:369:TYR:HA	1:B:377:PHE:CE2	2.43	0.54
2:C:40:MET:HB2	2:C:43:LYS:HG3	1.90	0.53
1:A:455:LEU:HD12	1:A:456:PHE:CZ	2.44	0.53
3:L:68:GLU:H	3:L:68:GLU:CD	2.16	0.53
2:C:209:VAL:C	2:C:217:LYS:HG3	2.34	0.53
3:D:149:LYS:HZ1	3:D:154:PRO:HB3	1.74	0.52
2:C:166:THR:O	2:C:170:VAL:HG22	2.10	0.52
2:H:118:PRO:HB3	2:H:146:TYR:HB3	1.91	0.51
1:B:437:ASN:ND2	1:B:506:GLN:OE1	2.35	0.51
3:L:28:ILE:N	3:L:28:ILE:HD12	2.26	0.51
3:L:29:GLY:CA	3:L:66:ASN:HD21	2.17	0.51
2:C:152:THR:OG1	2:C:210:ASN:HB3	2.11	0.51
3:D:185:GLN:O	3:D:189:HIS:ND1	2.44	0.51
2:C:4:LEU:HD12	2:C:92:CYS:SG	2.51	0.50
3:D:151:ASP:OD1	3:D:190:ARG:HG2	2.11	0.50
1:A:398:ASP:O	1:A:511:VAL:HA	2.12	0.50
1:B:417:LYS:O	1:B:421:TYR:HB2	2.12	0.49
3:D:135:LEU:HD23	3:D:176:SER:HB3	1.94	0.49
1:B:353:TRP:CZ2	1:B:466:ARG:HB3	2.47	0.49
2:H:176:VAL:HB	3:L:162:THR:HG22	1.95	0.49
3:D:68:GLU:HG2	3:D:69:ASN:OD1	2.13	0.49
3:L:50:TYR:HB3	3:L:51:PRO:HD3	1.95	0.48
1:A:471:GLU:OE2	3:D:31:LYS:NZ	2.46	0.48
2:C:147:PHE:HD1	2:C:148:PRO:CA	2.27	0.48
3:D:28:ILE:HG22	3:D:33:VAL:HG21	1.96	0.48
1:B:369:TYR:CE2	1:B:384:PRO:HB2	2.49	0.48
2:H:94:ARG:NH2	2:H:101:ASP:OD2	2.39	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:4:LEU:HD12	2:C:22:CYS:SG	2.53	0.48
1:A:417:LYS:O	1:A:421:TYR:HB2	2.14	0.47
3:D:189:HIS:CE1	3:D:192:TYR:HH	2.31	0.47
1:B:353:TRP:O	1:B:466:ARG:NH1	2.47	0.47
1:B:398:ASP:O	1:B:511:VAL:HA	2.14	0.47
1:A:506:GLN:HB2	1:A:508:TYR:HE1	1.79	0.47
1:B:394:ASN:HB2	1:B:516:GLU:OE2	2.14	0.47
2:C:36:TRP:CZ3	2:C:92:CYS:HB3	2.49	0.47
1:B:365:TYR:CE2	1:B:387:LEU:HD22	2.49	0.47
1:B:409:GLN:NE2	1:B:418:ILE:HB	2.30	0.47
1:B:517:LEU:HD13	1:B:519:HIS:H	1.80	0.46
1:A:338:PHE:CE2	1:A:363:ALA:HB1	2.51	0.46
2:C:116:LYS:HB2	2:C:116:LYS:HE2	1.65	0.46
3:L:49:TYR:HB2	3:L:53:ASP:HB3	1.97	0.46
2:C:101:ASP:OD1	2:C:101:ASP:N	2.48	0.46
3:D:35:TRP:HB2	3:D:48:ILE:HG22	1.97	0.46
2:C:210:ASN:CB	2:C:217:LYS:HD2	2.40	0.45
1:A:506:GLN:HB2	1:A:508:TYR:CE1	2.51	0.45
2:C:105:GLN:H	2:C:105:GLN:CD	2.24	0.45
1:B:377:PHE:CE1	1:B:434:ILE:HG12	2.51	0.45
3:L:135:LEU:HD23	3:L:176:SER:HB3	1.98	0.45
3:L:13:VAL:HG23	3:L:78:VAL:HG21	1.99	0.45
3:L:28:ILE:HG23	3:L:33:VAL:HG21	1.99	0.45
2:C:211:HIS:HB3	2:C:216:THR:HB	1.98	0.45
1:A:376:THR:HB	1:A:435:ALA:HB3	1.98	0.45
1:B:409:GLN:HE22	1:B:418:ILE:H	1.65	0.45
1:A:342:PHE:CZ	1:A:511:VAL:HG21	2.51	0.45
3:L:141:PRO:HD2	3:L:199:GLU:OE1	2.17	0.45
2:C:165:LEU:HD23	2:C:190:VAL:HG21	1.99	0.45
2:C:210:ASN:HB2	2:C:217:LYS:NZ	2.22	0.45
1:B:376:THR:HB	1:B:435:ALA:HB3	1.97	0.44
2:H:87:THR:HG23	2:H:110:THR:HA	1.99	0.44
1:B:409:GLN:NE2	1:B:418:ILE:H	2.15	0.44
3:D:186:TRP:CZ2	3:D:211:PRO:HA	2.53	0.44
1:A:502:GLY:O	1:A:506:GLN:HG2	2.18	0.44
3:L:129:LYS:HA	3:L:129:LYS:HD2	1.67	0.44
2:C:101:ASP:O	2:C:102:ILE:HD13	2.17	0.44
3:L:35:TRP:HB2	3:L:48:ILE:HB	1.99	0.44
2:H:29:PHE:CD2	2:H:76:SER:HA	2.52	0.43
3:L:4:LEU:HD23	3:L:4:LEU:HA	1.77	0.43
2:C:36:TRP:CE2	2:C:80:LEU:HB2	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:403:ARG:HG3	1:A:495:TYR:CE1	2.53	0.43
1:B:350:VAL:O	1:B:353:TRP:HD1	2.01	0.43
1:B:403:ARG:HG3	1:B:495:TYR:CE1	2.53	0.43
2:C:147:PHE:HD1	2:C:148:PRO:N	2.16	0.43
3:L:186:TRP:CZ2	3:L:211:PRO:HA	2.53	0.43
1:B:369:TYR:HA	1:B:377:PHE:HE2	1.82	0.43
2:C:220:LYS:HA	2:C:220:LYS:HD3	1.82	0.43
1:B:342:PHE:HZ	1:B:434:ILE:HD12	1.83	0.43
2:H:194:SER:O	2:H:197:LEU:HB2	2.19	0.43
2:C:123:LEU:HD12	2:C:141:CYS:N	2.33	0.43
1:A:389:ASP:N	1:A:389:ASP:OD1	2.52	0.43
2:C:123:LEU:HB3	3:D:118:PHE:CD1	2.54	0.43
2:C:122:PRO:HD3	2:C:220:LYS:HE3	2.01	0.42
2:C:101:ASP:C	2:C:102:ILE:HD13	2.44	0.42
2:C:174:PRO:HG2	3:D:165:SER:OG	2.18	0.42
1:A:498:GLN:HA	1:A:499:PRO:HD3	1.90	0.42
1:B:378:LYS:HD3	1:B:379:CYS:N	2.33	0.42
3:D:17:LYS:HD3	3:D:17:LYS:HA	1.61	0.42
1:A:431:GLY:HA2	1:A:515:PHE:CD2	2.55	0.42
3:L:13:VAL:HG21	3:L:78:VAL:HG11	2.02	0.42
3:L:50:TYR:CD1	3:L:50:TYR:C	2.98	0.42
1:B:377:PHE:CD1	1:B:434:ILE:HG12	2.54	0.42
2:C:93:ALA:HB1	2:C:100(D):PHE:HB3	2.01	0.42
1:A:468:ILE:HD11	3:D:51:PRO:HD3	2.02	0.42
1:B:415:THR:OG1	1:B:416:GLY:N	2.53	0.42
2:H:38:ARG:HB2	2:H:48:MET:SD	2.60	0.42
1:B:403:ARG:HD2	1:B:505:TYR:HD1	1.84	0.42
3:D:162:THR:HG23	3:D:176:SER:O	2.19	0.42
1:A:341:VAL:HG23	1:A:342:PHE:CD1	2.49	0.42
1:A:350:VAL:O	1:A:353:TRP:HD1	2.02	0.42
1:A:405:ASP:OD1	1:A:405:ASP:N	2.51	0.41
3:D:36:TYR:CE2	3:D:46:VAL:HG22	2.55	0.41
1:A:393:THR:CG2	1:A:520:ALA:HB3	2.50	0.41
1:A:396:TYR:HB2	1:A:514:SER:HB2	2.02	0.41
1:A:424:LYS:HE2	1:A:424:LYS:HB2	1.84	0.41
2:C:12:LYS:O	2:C:111:VAL:HA	2.20	0.41
3:D:13:VAL:HG11	3:D:19:ALA:HB2	2.02	0.41
1:A:393:THR:HG21	1:A:520:ALA:HB3	2.03	0.41
2:C:208:ASN:C	2:C:217:LYS:HE3	2.45	0.41
1:A:335:LEU:HD22	1:A:362:VAL:O	2.21	0.41
1:A:393:THR:N	1:A:516:GLU:O	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:386:LYS:HA	1:B:389:ASP:OD2	2.20	0.41
1:B:442:ASP:HB3	1:B:451:TYR:HE2	1.86	0.41
3:L:25:GLY:CA	3:L:28:ILE:HD11	2.51	0.41
1:A:440:ASN:OD1	1:A:440:ASN:N	2.47	0.40
2:H:211:HIS:CD2	2:H:213:PRO:HD2	2.56	0.40
2:H:51:ILE:HD12	2:H:57:THR:HG22	2.03	0.40
2:C:147:PHE:HA	2:C:148:PRO:HA	1.89	0.40
1:A:427:ASP:OD1	1:A:427:ASP:N	2.53	0.40
3:D:140:TYR:CD1	3:D:141:PRO:HA	2.56	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	193/205 (94%)	190 (98%)	3 (2%)	0	100	100
1	B	193/205 (94%)	189 (98%)	4 (2%)	0	100	100
2	C	220/223 (99%)	215 (98%)	4 (2%)	1 (0%)	24	43
2	H	212/223 (95%)	207 (98%)	5 (2%)	0	100	100
3	D	209/214 (98%)	202 (97%)	6 (3%)	1 (0%)	24	43
3	L	208/214 (97%)	199 (96%)	8 (4%)	1 (0%)	24	43
All	All	1235/1284 (96%)	1202 (97%)	30 (2%)	3 (0%)	43	62

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	C	132	THR
3	D	51	PRO
3	L	51	PRO

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	167/177 (94%)	163 (98%)	4 (2%)	43	67
1	B	165/177 (93%)	162 (98%)	3 (2%)	51	74
2	C	186/187 (100%)	176 (95%)	10 (5%)	20	42
2	H	181/187 (97%)	173 (96%)	8 (4%)	25	49
3	D	176/179 (98%)	169 (96%)	7 (4%)	28	53
3	L	175/179 (98%)	169 (97%)	6 (3%)	32	57
All	All	1050/1086 (97%)	1012 (96%)	38 (4%)	31	56

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	367	VAL
1	A	387	LEU
1	A	389	ASP
1	A	501	ASN
1	B	368	LEU
1	B	373	SER
1	B	514	SER
2	H	25	SER
2	H	85	SER
2	H	142	LEU
2	H	143	VAL
2	H	151	VAL
2	H	187	SER
2	H	197	LEU
2	H	207	CYS
3	L	9	SER
3	L	30	SER
3	L	46	VAL
3	L	65	SER
3	L	115	VAL
3	L	197	THR
2	C	22	CYS

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Mol	Chain	Res	Type
2	C	30	THR
2	C	62	SER
2	C	67	VAL
2	C	119	SER
2	C	167	SER
2	C	188	SER
2	C	197	LEU
2	C	207	CYS
2	C	212	LYS
3	D	11	VAL
3	D	46	VAL
3	D	70	THR
3	D	153	SER
3	D	162	THR
3	D	176	SER
3	D	212	THR

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

Mol	Chain	Res	Type
1	A	501	ASN
1	B	409	GLN
1	B	439	ASN
2	H	3	GLN
2	H	39	GLN
2	H	66	GLN
2	H	178	GLN
3	L	34	HIS
3	L	37	GLN
3	L	38	GLN
3	L	42	GLN
3	L	66	ASN
3	L	195	GLN
2	C	210	ASN
3	D	38	GLN
3	D	94	ASN

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 ⓘ

2 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	$\# Z > 2$
4	NAG	E	1	1,4	14,14,15	0.28	0	17,19,21	0.47	0
4	NAG	E	2	4	14,14,15	0.26	0	17,19,21	0.45	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
4	NAG	E	1	1,4	-	2/6/23/26	0/1/1/1
4	NAG	E	2	4	-	1/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

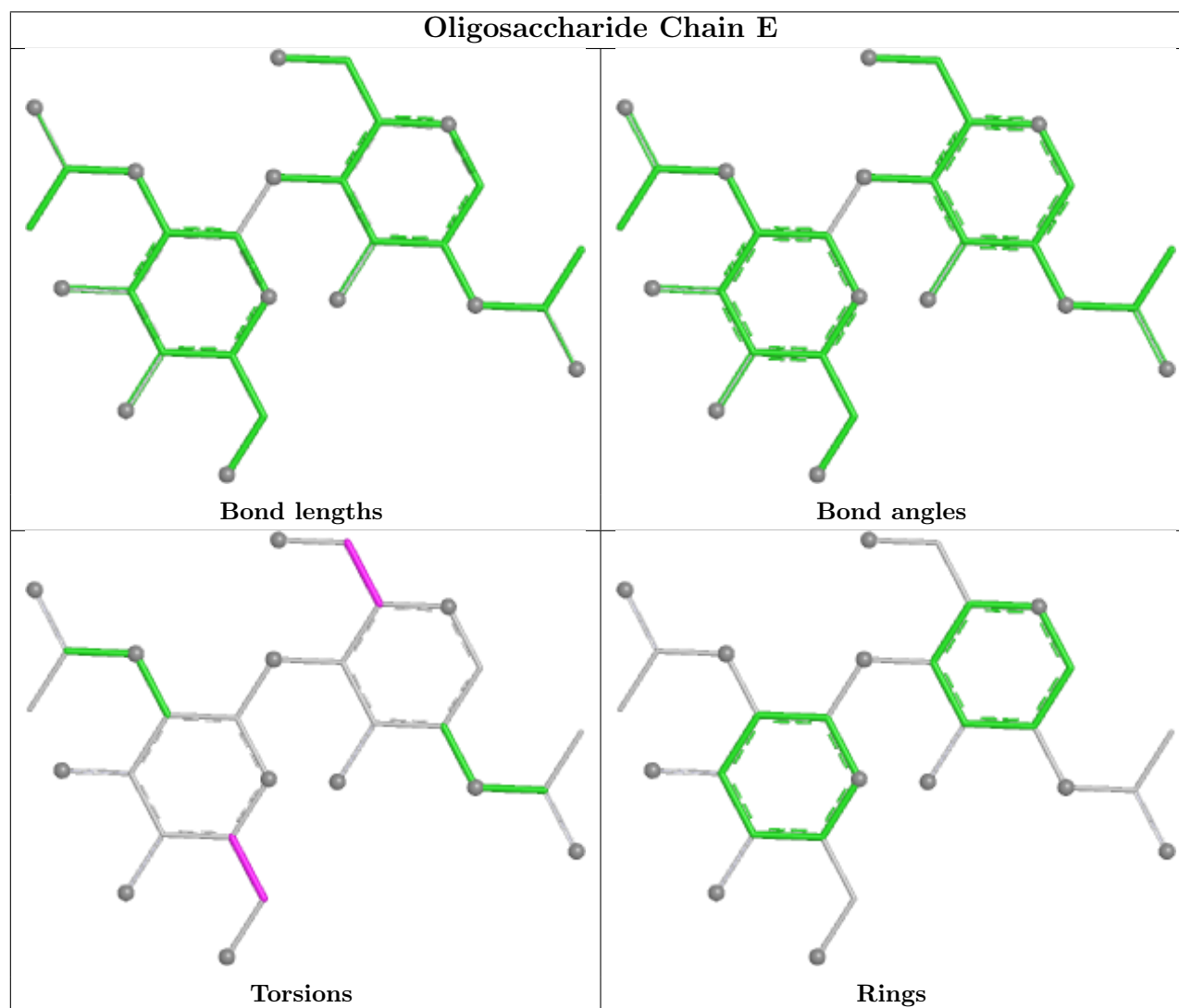
All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	E	1	NAG	C4-C5-C6-O6
4	E	2	NAG	O5-C5-C6-O6
4	E	1	NAG	O5-C5-C6-O6

There are no ring outliers.

No monomer is involved in short contacts.

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](#)

1 ligand is 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	$\# Z > 2$
5	NAG	B	1001	1	14,14,15	0.31	0	17,19,21	0.46	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
5	NAG	B	1001	1	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

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	195/205 (95%)	0.67	12 (6%)	26 20	30, 67, 109, 124	0
1	B	195/205 (95%)	0.58	6 (3%)	51 42	39, 74, 119, 131	0
2	C	222/223 (99%)	0.27	1 (0%)	87 85	28, 52, 87, 106	0
2	H	216/223 (96%)	0.10	2 (0%)	81 76	31, 48, 80, 125	0
3	D	210/214 (98%)	0.20	4 (1%)	66 59	21, 52, 84, 96	1 (0%)
3	L	210/214 (98%)	0.18	4 (1%)	66 59	29, 49, 70, 97	0
All	All	1248/1284 (97%)	0.32	29 (2%)	61 52	21, 54, 102, 131	1 (0%)

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	D	98[A]	PHE	4.8
1	A	338	PHE	3.7
1	B	334	ASN	3.5
1	A	350	VAL	3.1
3	L	212	THR	2.9
3	D	101	GLY	2.9
3	D	25	GLY	2.8
1	B	515	PHE	2.8
1	B	400	PHE	2.8
1	A	337	PRO	2.8
1	A	344	ALA	2.7
3	L	2	SER	2.6
1	B	432	CYS	2.6
1	A	368	LEU	2.4
3	L	3	ALA	2.4
1	A	432	CYS	2.3
2	H	35	GLY	2.3
1	A	498	GLN	2.3
1	A	486	PHE	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	523	THR	2.2
1	A	441	LEU	2.2
1	A	372	ALA	2.1
3	L	210	ALA	2.1
2	C	141	CYS	2.1
3	D	2	SER	2.1
1	A	442	ASP	2.1
1	A	435	ALA	2.1
2	H	141	CYS	2.0
1	B	503	VAL	2.0

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

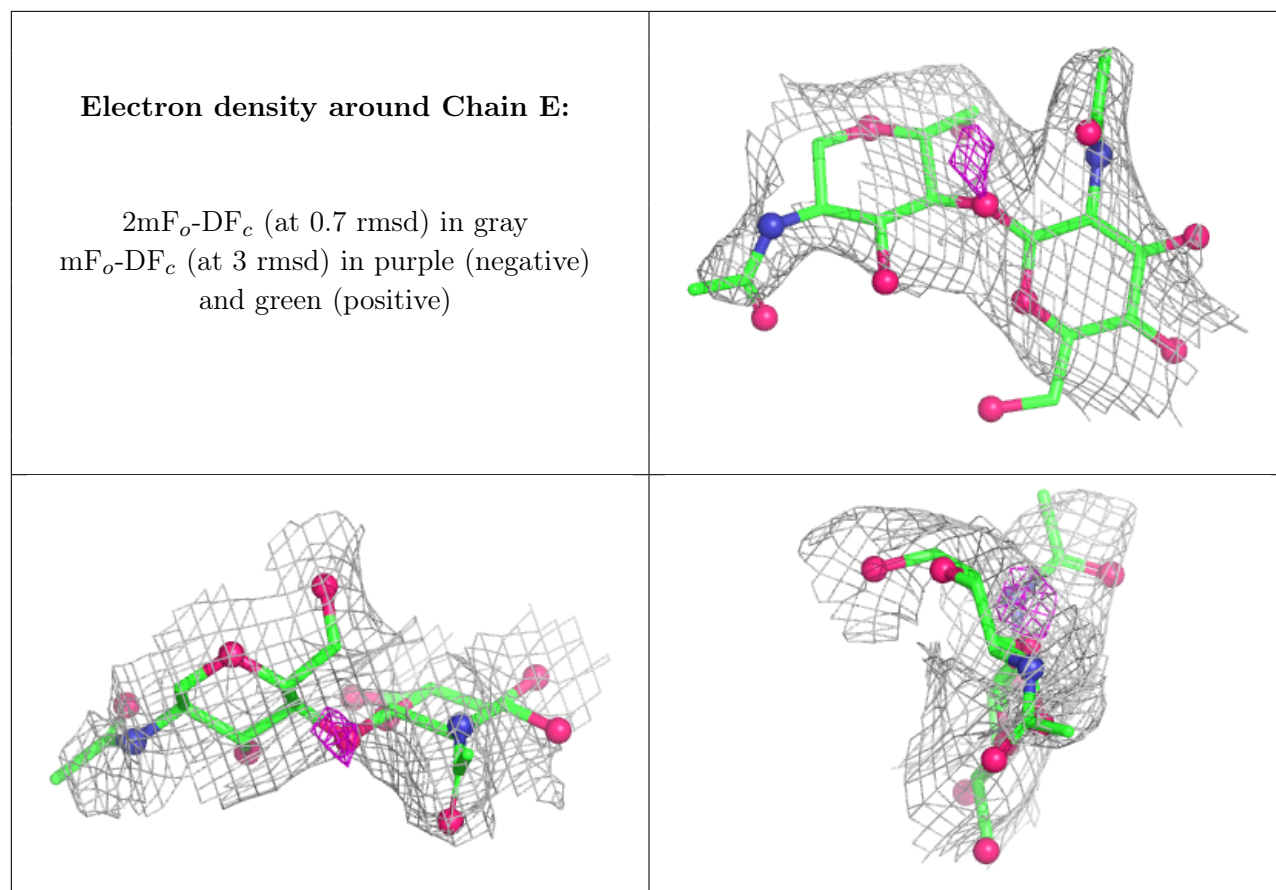
There are no non-standard protein/DNA/RNA residues in this entry.

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
4	NAG	E	1	14/15	-	-	90,108,115,117	0
4	NAG	E	2	14/15	-	-	92,112,117,122	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
5	NAG	B	1001	14/15	0.77	0.11	99,110,111,115	0

6.5 Other polymers [i](#)

There are no such residues in this entry.