



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 11:47 PM UTC

PDB ID : 6BJ5 / pdb_00006bj5
Title : Structure of the Clinically used Myxomaviral Serine Protease Inhibitor 1 (SERP-1)
Authors : Mahon, B.P.; Lomelino, C.L.; McKenna, R.
Deposited on : 2017-11-05
Resolution : 2.50 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

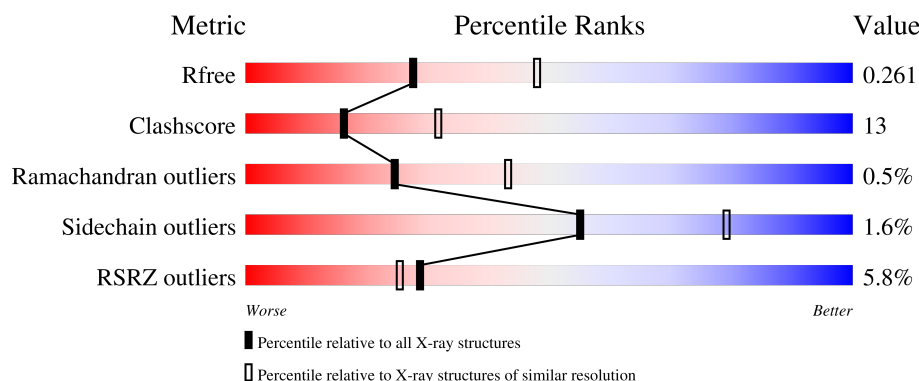
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 Å.

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



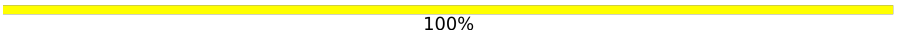
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	315	4% 73% 18% • 7%
1	B	315	5% 68% 24% • 7%
1	C	315	8% 67% 23% • 9%
2	F	54	2% 52% 17% • 30%
2	G	54	4% 57% 15% 28%

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Mol	Chain	Length	Quality of chain
2	H	54	 69% 6% 26%
3	D	2	 50% 50%
3	E	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
3	NAG	D	1	-	-	X	-
3	NAG	E	2	X	-	-	-

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 8023 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 Serine proteinase inhibitor 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	293	Total	C	N	O	S	0	0	0
			2334	1476	399	451	8			
1	A	292	Total	C	N	O	S	0	0	0
			2322	1467	397	450	8			
1	C	288	Total	C	N	O	S	0	0	0
			2297	1451	394	444	8			

- Molecule 2 is a protein called Serine proteinase inhibitor 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	F	38	Total	C	N	O	S	0	0	0
			302	203	45	52	2			
2	G	39	Total	C	N	O	S	0	0	0
			304	205	46	51	2			
2	H	40	Total	C	N	O	S	0	0	0
			316	211	47	56	2			

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



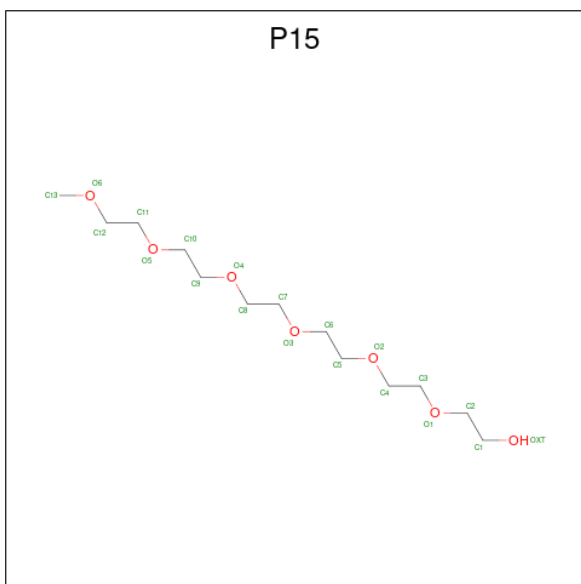
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	D	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	E	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 4 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



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

- Molecule 5 is 2,5,8,11,14,17-HEXAOXANONADECAN-19-OL (CCD ID: P15) (formula: $C_{13}H_{28}O_7$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	C	1	Total	C	O	0	0
			20	13	7		

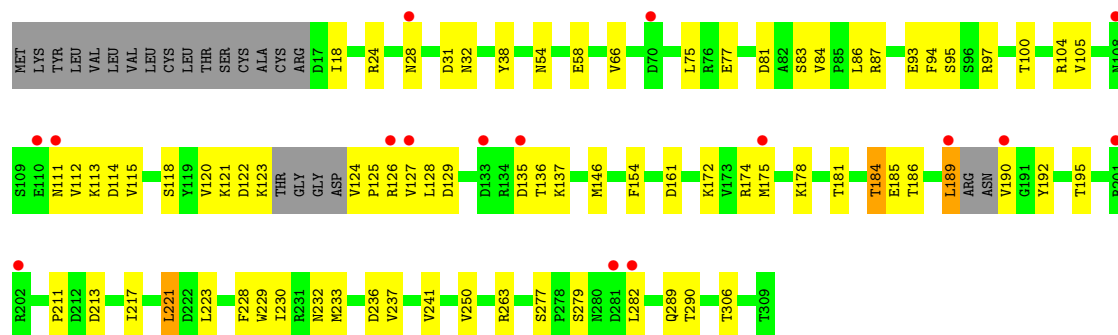
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	14	Total 14	O 14	0	0
6	F	1	Total 1	O 1	0	0
6	A	16	Total 16	O 16	0	0
6	G	7	Total 7	O 7	0	0
6	C	18	Total 18	O 18	0	0
6	H	2	Total 2	O 2	0	0

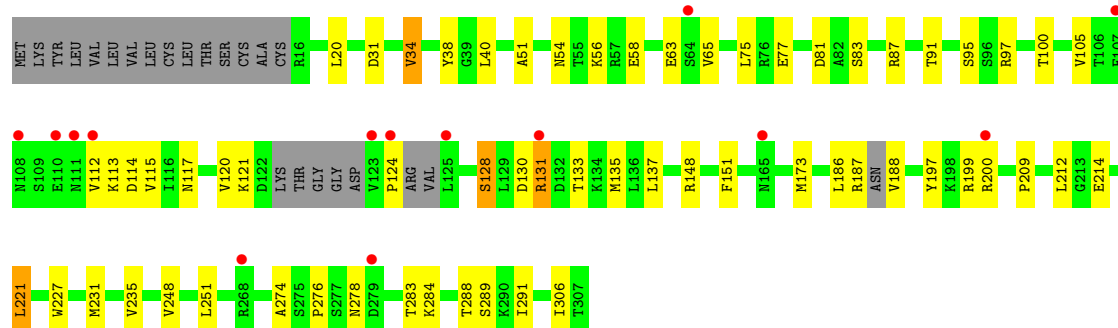
3 Residue-property plots [i](#)

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: Serine proteinase inhibitor 1

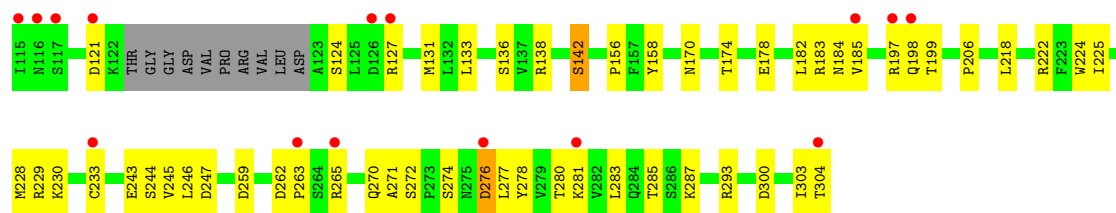


- Molecule 1: Serine proteinase inhibitor 1



- Molecule 1: Serine proteinase inhibitor 1

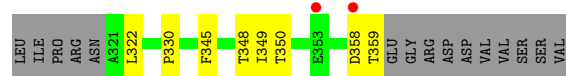




• Molecule 2: Serine proteinase inhibitor 1



• Molecule 2: Serine proteinase inhibitor 1



• Molecule 2: Serine proteinase inhibitor 1



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



• Molecule 3: 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	99.01Å 122.07Å 130.09Å 90.00° 94.88° 90.00°	Depositor
Resolution (Å)	19.83 – 2.50 19.83 – 2.50	Depositor EDS
% Data completeness (in resolution range)	86.8 (19.83-2.50) 86.5 (19.83-2.50)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.97 (at 2.50Å)	Xtriage
Refinement program	PHENIX 1.10.1_2155	Depositor
R, R_{free}	0.207 , 0.257 0.211 , 0.261	Depositor DCC
R_{free} test set	1991 reflections (3.74%)	wwPDB-VP
Wilson B-factor (Å ²)	35.9	Xtriage
Anisotropy	0.060	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 54.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	8023	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.60% 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, P15

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	0.42	0/2362	0.74	5/3201 (0.2%)
1	B	0.44	0/2375	0.73	0/3220
1	C	0.42	0/2337	0.77	1/3166 (0.0%)
2	F	0.47	0/309	0.74	1/419 (0.2%)
2	G	0.43	0/311	0.79	0/422
2	H	0.55	0/323	0.70	0/438
All	All	0.44	0/8017	0.75	7/10866 (0.1%)

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	131	ARG	CG-CD-NE	7.61	128.75	112.00
1	A	131	ARG	CB-CG-CD	-6.15	97.16	111.30
1	A	130	ASP	N-CA-C	-5.42	103.23	110.55
1	C	276	ASP	N-CA-C	-5.37	105.17	112.26
1	A	130	ASP	CA-C-N	-5.20	112.19	122.02
1	A	130	ASP	C-N-CA	-5.20	112.19	122.02
2	F	346	MET	CB-CG-SD	-5.19	97.12	112.70

There are no chirality outliers.

There are no planarity outliers.

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	2322	0	2308	51	0
1	B	2334	0	2331	69	0
1	C	2297	0	2291	80	0
2	F	302	0	320	10	0
2	G	304	0	323	6	0
2	H	316	0	331	5	0
3	D	28	0	25	13	0
3	E	28	0	25	5	0
4	A	14	0	13	0	0
5	C	20	0	28	2	0
6	A	16	0	0	0	0
6	B	14	0	0	2	0
6	C	18	0	0	2	0
6	F	1	0	0	0	0
6	G	7	0	0	0	0
6	H	2	0	0	0	0
All	All	8023	0	7995	209	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 13.

All (209) 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:B:28:ASN:ND2	3:D:1:NAG:O5	1.62	1.31
1:C:28:ASN:HD21	3:E:1:NAG:C1	1.50	1.23
1:C:28:ASN:ND2	3:E:1:NAG:C1	2.00	1.22
1:B:175:MET:SD	1:B:241:VAL:CG1	2.35	1.13
1:B:175:MET:SD	1:B:241:VAL:HG12	1.91	1.09
3:D:2:NAG:H5	3:D:2:NAG:H82	1.15	1.08
1:B:28:ASN:ND2	3:D:1:NAG:C1	2.21	1.04
1:B:28:ASN:CG	3:D:1:NAG:C1	2.38	0.97
1:C:233:CYS:SG	2:H:323:THR:HB	2.13	0.87
3:D:2:NAG:H82	3:D:2:NAG:C5	2.03	0.87
1:B:28:ASN:OD1	3:D:1:NAG:C1	2.23	0.86
1:B:28:ASN:ND2	1:B:31:ASP:CB	2.39	0.84
1:B:24:ARG:NH2	1:B:221:LEU:O	2.09	0.84
1:B:113:LYS:HZ1	1:B:129:ASP:HB2	1.46	0.81
1:A:51:ALA:O	1:A:56:LYS:NZ	2.15	0.80
1:C:107:PHE:HA	1:C:111:VAL:HG21	1.61	0.79
1:B:28:ASN:ND2	1:B:31:ASP:HB2	1.98	0.79
1:B:28:ASN:ND2	1:B:31:ASP:HB3	1.98	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:63:GLU:O	1:A:97:ARG:NH2	2.18	0.77
1:C:272:SER:HG	1:C:274:SER:HG	1.23	0.76
1:C:259:ASP:OD1	1:C:265:ARG:NH2	2.18	0.76
1:A:113:LYS:O	1:A:117:ASN:ND2	2.17	0.76
1:B:113:LYS:NZ	1:B:129:ASP:HB2	2.00	0.76
1:C:170:ASN:HD21	5:C:401:P15:H71	1.48	0.76
1:C:178:GLU:CD	1:C:229:ARG:HH12	1.93	0.75
2:F:358:ASP:OD2	2:F:359:THR:N	2.20	0.74
1:B:146:MET:HE1	2:F:335:ILE:HG21	1.71	0.72
1:A:34:VAL:HG13	1:A:291:ILE:HD12	1.71	0.71
1:A:187:ARG:C	1:A:187:ARG:HD2	2.16	0.71
1:B:31:ASP:HB2	3:D:1:NAG:H62	1.73	0.69
1:B:28:ASN:HD21	3:D:1:NAG:C5	2.01	0.68
1:A:131:ARG:HD2	1:A:131:ARG:C	2.18	0.67
1:A:91:THR:O	1:A:95:SER:OG	2.10	0.67
1:C:93:GLU:N	1:C:93:GLU:OE2	2.28	0.66
1:B:121:LYS:HB2	1:B:124:VAL:HG23	1.78	0.65
1:C:80:VAL:HG11	1:C:86:LEU:HD11	1.78	0.65
1:C:133:LEU:HB3	1:C:303:ILE:HB	1.79	0.64
1:B:87:ARG:NH1	6:B:501:HOH:O	2.27	0.64
1:C:136:SER:HB3	1:C:138:ARG:HH12	1.61	0.64
1:B:122:ASP:O	1:B:123:LYS:HB3	1.96	0.64
1:C:233:CYS:SG	2:H:323:THR:CB	2.85	0.64
1:C:28:ASN:ND2	3:E:1:NAG:O5	2.22	0.63
1:B:32:ASN:ND2	2:F:350:THR:H	1.97	0.63
1:A:81:ASP:OD2	1:A:83:SER:OG	2.11	0.63
1:C:233:CYS:SG	2:H:323:THR:CG2	2.87	0.63
1:B:28:ASN:HD22	1:B:31:ASP:HB3	1.62	0.63
1:B:277:SER:OG	1:B:279:SER:HB3	1.99	0.62
1:B:31:ASP:HB2	3:D:1:NAG:C6	2.30	0.62
1:A:197:TYR:O	1:A:199:ARG:HA	2.00	0.61
1:C:281:LYS:CE	1:C:304:THR:HA	2.30	0.61
3:D:1:NAG:O7	3:D:1:NAG:O3	2.20	0.60
1:A:148:ARG:HE	1:A:199:ARG:HH21	1.48	0.60
1:C:280:THR:H	1:C:304:THR:HG1	1.47	0.59
1:B:75:LEU:HD21	1:B:77:GLU:HG2	1.83	0.59
1:B:175:MET:SD	1:B:241:VAL:HG13	2.40	0.59
2:F:334:LEU:HD23	2:F:346:MET:HG2	1.84	0.59
1:C:281:LYS:HE3	1:C:304:THR:HA	1.84	0.59
1:B:84:VAL:O	1:B:104:ARG:NH1	2.30	0.59
1:A:199:ARG:HB3	1:A:200:ARG:HD3	1.85	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:280:THR:N	1:C:304:THR:OG1	2.25	0.59
1:A:248:VAL:HG22	1:A:288:THR:HG22	1.85	0.58
1:C:170:ASN:ND2	5:C:401:P15:H71	2.18	0.58
1:B:175:MET:SD	1:B:241:VAL:HG11	2.38	0.57
1:A:148:ARG:HE	1:A:199:ARG:NH2	2.01	0.57
1:B:190:VAL:HG21	1:B:213:ASP:HB3	1.87	0.57
1:C:222:ARG:HD3	6:C:513:HOH:O	2.03	0.57
1:A:137:LEU:HB3	1:A:306:ILE:HB	1.87	0.56
1:C:182:LEU:O	1:C:185:VAL:HG22	2.07	0.55
1:B:111:ASN:O	1:B:115:VAL:HG23	2.07	0.55
1:C:26:VAL:HB	1:C:246:LEU:HD23	1.88	0.54
1:A:120:VAL:HG21	1:A:124:PRO:HG2	1.90	0.54
1:A:131:ARG:C	1:A:133:THR:H	2.16	0.54
1:C:124:SER:O	1:C:281:LYS:HE2	2.07	0.54
1:A:114:ASP:HA	1:A:117:ASN:HB2	1.89	0.53
1:C:262:ASP:CG	1:C:265:ARG:HH12	2.16	0.53
1:B:105:VAL:HG22	1:B:112:VAL:HG23	1.89	0.53
1:C:281:LYS:CD	1:C:304:THR:HA	2.38	0.53
1:C:197:ARG:O	1:C:197:ARG:HG2	2.08	0.53
1:B:135:ASP:O	1:B:137:LYS:HE2	2.09	0.53
1:B:120:VAL:O	1:B:121:LYS:HD2	2.09	0.52
1:B:229:TRP:O	1:B:233:MET:HG3	2.10	0.52
1:C:136:SER:HB3	1:C:138:ARG:NH1	2.24	0.52
1:C:281:LYS:NZ	1:C:304:THR:O	2.41	0.52
1:A:235:VAL:HG12	2:G:322:LEU:HD11	1.91	0.52
1:C:272:SER:HB3	1:C:277:LEU:CD1	2.40	0.52
1:C:79:PHE:CD1	1:C:103:GLN:HB2	2.45	0.52
1:C:281:LYS:HD2	1:C:304:THR:HA	1.92	0.52
1:A:75:LEU:HD11	1:A:77:GLU:OE1	2.10	0.52
1:C:142:SER:HB3	1:C:293:ARG:HD3	1.91	0.51
1:A:87:ARG:HG3	1:A:274:ALA:O	2.11	0.51
1:A:34:VAL:HG12	1:A:289:SER:HB2	1.92	0.51
1:B:263:ARG:HD3	6:B:514:HOH:O	2.10	0.51
1:A:283:THR:OG1	1:A:284:LYS:HG3	2.11	0.51
1:B:114:ASP:O	1:B:118:SER:HB3	2.10	0.51
1:B:228:PHE:HE2	1:B:232:ASN:ND2	2.08	0.50
1:C:138:ARG:HH11	1:C:138:ARG:HG3	1.76	0.50
1:A:38:TYR:OH	1:A:65:VAL:HG12	2.12	0.50
1:C:272:SER:HB3	1:C:277:LEU:HD11	1.94	0.50
1:C:86:LEU:HD23	1:C:91:THR:HG22	1.94	0.50
1:C:246:LEU:HD12	1:C:246:LEU:N	2.26	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:197:ARG:C	1:C:198:GLN:OE1	2.54	0.49
1:A:34:VAL:CG1	1:A:291:ILE:HD12	2.40	0.49
1:A:63:GLU:HB3	1:A:97:ARG:HH22	1.77	0.49
1:B:32:ASN:HD22	2:F:350:THR:H	1.59	0.49
1:B:184:THR:HG23	1:B:195:THR:OG1	2.13	0.49
1:C:93:GLU:HA	1:C:96:SER:HB3	1.95	0.48
3:E:2:NAG:O7	3:E:2:NAG:O3	2.31	0.48
1:A:95:SER:HA	1:A:100:THR:O	2.12	0.48
1:A:151:PHE:CD1	1:A:173:MET:HB3	2.49	0.48
1:C:243:GLU:HG3	1:C:287:LYS:HB2	1.95	0.48
1:C:138:ARG:NH1	1:C:138:ARG:HG3	2.28	0.48
1:B:94:PHE:HE2	1:B:100:THR:HG23	1.78	0.48
1:C:174:THR:HG22	1:C:230:LYS:HE3	1.95	0.48
1:C:107:PHE:O	1:C:111:VAL:HG11	2.14	0.48
1:B:38:TYR:OH	1:B:66:VAL:O	2.29	0.47
1:C:105:VAL:HG13	1:C:111:VAL:HG23	1.96	0.47
1:B:192:TYR:CD2	1:B:217:ILE:HD11	2.49	0.47
1:B:95:SER:HA	1:B:100:THR:O	2.14	0.47
1:B:236:ASP:OD2	1:B:237:VAL:N	2.48	0.47
1:C:71:ALA:HB1	1:C:199:THR:HG23	1.96	0.47
1:C:131:MET:O	1:C:304:THR:C	2.57	0.47
1:C:81:ASP:OD2	1:C:83:SER:OG	2.15	0.46
1:A:54:ASN:O	1:A:58:GLU:HG3	2.15	0.46
1:C:263:PRO:HG3	1:C:278:TYR:CE1	2.51	0.46
2:H:330:PRO:HB3	2:H:350:THR:HG22	1.96	0.46
1:B:28:ASN:HD22	1:B:31:ASP:CB	2.21	0.46
1:C:82:ALA:N	1:C:105:VAL:O	2.46	0.46
1:C:244:SER:HB3	1:C:246:LEU:HD11	1.97	0.46
1:A:199:ARG:C	1:A:200:ARG:HD3	2.41	0.46
1:A:200:ARG:HH11	1:C:184:ASN:ND2	2.13	0.46
2:F:330:PRO:HA	2:F:349:ILE:O	2.16	0.46
1:C:265:ARG:HB2	1:C:265:ARG:CZ	2.45	0.46
2:G:358:ASP:O	2:G:359:THR:HB	2.16	0.45
1:B:126:ARG:C	1:B:128:LEU:H	2.24	0.45
1:A:186:LEU:O	1:A:188:VAL:N	2.50	0.45
1:C:142:SER:HB3	1:C:293:ARG:CD	2.45	0.45
1:A:20:LEU:HD22	1:A:221:LEU:HD22	1.98	0.45
1:C:197:ARG:O	1:C:198:GLN:CD	2.60	0.45
1:B:28:ASN:HD21	1:B:31:ASP:HB2	1.76	0.44
1:A:128:SER:HA	1:A:284:LYS:HE2	1.99	0.44
1:A:40:LEU:HD22	1:A:251:LEU:HD13	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:124:SER:HB2	1:C:283:LEU:HD12	2.00	0.44
1:B:84:VAL:HG11	1:B:282:LEU:HD11	1.99	0.44
1:B:185:GLU:HG3	1:B:186:THR:O	2.17	0.44
1:C:50:ALA:HB1	1:C:271:ALA:HB2	2.00	0.44
1:A:31:ASP:O	2:G:350:THR:HG22	2.18	0.44
1:C:81:ASP:HA	1:C:105:VAL:O	2.17	0.44
1:C:272:SER:CB	1:C:277:LEU:HD11	2.48	0.44
1:B:31:ASP:CB	3:D:1:NAG:H62	2.43	0.44
1:B:289:GLN:HG3	1:B:306:THR:OG1	2.18	0.44
1:A:227:TRP:O	1:A:231:MET:HG3	2.18	0.44
1:C:138:ARG:HH22	1:C:300:ASP:CG	2.26	0.44
1:A:291:ILE:HD11	2:G:345:PHE:HB3	1.98	0.43
1:A:200:ARG:NH1	1:C:184:ASN:HD21	2.15	0.43
1:C:156:PRO:HB2	1:C:158:TYR:CZ	2.53	0.43
1:B:81:ASP:OD2	1:B:83:SER:OG	2.30	0.43
1:B:181:THR:HG23	1:B:236:ASP:O	2.19	0.43
1:C:270:GLN:H	1:C:270:GLN:HG2	1.57	0.43
1:B:189:LEU:H	1:B:190:VAL:N	2.17	0.43
1:C:23:PHE:HA	1:C:26:VAL:HG22	2.00	0.43
1:B:146:MET:HE1	2:F:335:ILE:HD13	1.99	0.43
1:A:120:VAL:HG21	1:A:124:PRO:CG	2.48	0.43
1:C:28:ASN:CG	3:E:1:NAG:C1	2.84	0.42
1:B:223:LEU:HD23	1:B:223:LEU:HA	1.90	0.42
1:C:107:PHE:C	1:C:127:ARG:NE	2.77	0.42
3:D:2:NAG:H3	3:D:2:NAG:C8	2.50	0.42
1:B:28:ASN:CG	3:D:1:NAG:O5	2.40	0.42
1:B:172:LYS:HB2	2:F:356:ILE:HG13	2.00	0.42
1:A:121:LYS:HA	1:A:121:LYS:HD2	1.56	0.42
1:B:136:THR:C	1:B:137:LYS:HE2	2.44	0.42
1:A:148:ARG:HH21	1:A:199:ARG:NH2	2.18	0.42
1:B:250:VAL:HG22	1:B:290:THR:HG22	2.00	0.42
1:C:245:VAL:HG22	1:C:285:THR:HG22	2.01	0.42
1:C:218:LEU:HD23	1:C:218:LEU:HA	1.87	0.42
1:C:105:VAL:CG1	1:C:111:VAL:HG23	2.50	0.42
1:C:247:ASP:HA	1:C:283:LEU:HD23	2.01	0.42
1:B:185:GLU:HG3	1:B:186:THR:N	2.36	0.41
1:B:230:ILE:HA	1:B:233:MET:HE2	2.02	0.41
1:A:112:VAL:HA	1:A:115:VAL:HG12	2.03	0.41
1:A:131:ARG:C	1:A:131:ARG:CD	2.91	0.41
1:C:233:CYS:SG	2:H:323:THR:HG21	2.59	0.41
1:C:246:LEU:HD13	6:C:502:HOH:O	2.19	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:18:ILE:HD13	1:B:18:ILE:HG21	1.85	0.41
1:B:135:ASP:C	1:B:137:LYS:HE2	2.45	0.41
1:B:146:MET:HB3	1:B:146:MET:HE2	1.63	0.41
1:A:212:LEU:HD21	2:G:348:THR:HB	2.03	0.41
1:C:124:SER:CB	1:C:283:LEU:HD12	2.51	0.41
1:C:224:TRP:O	1:C:228:MET:HG3	2.19	0.41
1:A:81:ASP:HA	1:A:105:VAL:O	2.20	0.41
1:A:278:ASN:O	1:A:278:ASN:CG	2.64	0.41
1:C:127:ARG:HD2	1:C:127:ARG:HA	1.94	0.41
1:C:182:LEU:HD23	1:C:183:ARG:HG3	2.03	0.41
1:B:32:ASN:HD21	2:F:351:LYS:H	1.69	0.41
1:B:81:ASP:HA	1:B:105:VAL:O	2.20	0.41
1:C:276:ASP:O	1:C:276:ASP:OD1	2.38	0.41
1:A:186:LEU:O	1:A:187:ARG:HB3	2.22	0.40
1:C:225:ILE:HA	1:C:228:MET:HE2	2.02	0.40
1:B:54:ASN:O	1:B:58:GLU:HG3	2.21	0.40
1:B:161:ASP:OD1	1:B:174:ARG:NH1	2.51	0.40
1:A:87:ARG:HG2	1:A:276:PRO:HD3	2.03	0.40
1:A:148:ARG:HH21	1:A:199:ARG:HH21	1.70	0.40
1:A:200:ARG:HH11	1:C:184:ASN:HD21	1.69	0.40
1:B:93:GLU:O	1:B:97:ARG:HG3	2.21	0.40
1:B:154:PHE:CZ	1:B:178:LYS:HE2	2.56	0.40
1:A:188:VAL:HG21	1:A:214:GLU:HG2	2.02	0.40
1:C:121:ASP:OD1	1:C:121:ASP:C	2.64	0.40
2:F:356:ILE:HA	2:F:356:ILE:HD13	1.83	0.40
2:G:330:PRO:HA	2:G:349:ILE:O	2.22	0.40
1:C:90:PHE:CG	1:C:271:ALA:HA	2.56	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	284/315 (90%)	275 (97%)	8 (3%)	1 (0%)	30	49
1	B	287/315 (91%)	278 (97%)	6 (2%)	3 (1%)	12	24
1	C	282/315 (90%)	271 (96%)	10 (4%)	1 (0%)	30	49
2	F	36/54 (67%)	36 (100%)	0	0	100	100
2	G	37/54 (68%)	37 (100%)	0	0	100	100
2	H	38/54 (70%)	38 (100%)	0	0	100	100
All	All	964/1107 (87%)	935 (97%)	24 (2%)	5 (0%)	24	43

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	127	VAL
1	B	125	PRO
1	B	211	PRO
1	A	209	PRO
1	C	206	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	264/285 (93%)	260 (98%)	4 (2%)	57	80
1	B	266/285 (93%)	262 (98%)	4 (2%)	57	80
1	C	261/285 (92%)	255 (98%)	6 (2%)	44	72
2	F	34/48 (71%)	34 (100%)	0	100	100
2	G	33/48 (69%)	33 (100%)	0	100	100
2	H	35/48 (73%)	35 (100%)	0	100	100
All	All	893/999 (89%)	879 (98%)	14 (2%)	55	79

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	86	LEU
1	B	184	THR
1	B	189	LEU
1	B	221	LEU
1	A	34	VAL
1	A	128	SER
1	A	135	MET
1	A	221	LEU
1	C	61	VAL
1	C	64	SER
1	C	83	SER
1	C	86	LEU
1	C	114	VAL
1	C	142	SER

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

Mol	Chain	Res	Type
1	B	32	ASN
1	B	103	GLN
1	B	117	ASN
1	A	103	GLN
1	A	201	GLN
1	A	230	ASN
1	C	116	ASN
1	C	170	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 ⓘ

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	D	1	3	14,14,15	0.26	0	17,19,21	0.49	0
3	NAG	D	2	3	14,14,15	0.51	0	17,19,21	1.75	3 (17%)
3	NAG	E	1	3	14,14,15	0.28	0	17,19,21	0.55	0
3	NAG	E	2	3	14,14,15	0.29	0	17,19,21	0.49	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	D	1	3	-	4/6/23/26	0/1/1/1
3	NAG	D	2	3	-	4/6/23/26	0/1/1/1
3	NAG	E	1	3	-	0/6/23/26	0/1/1/1
3	NAG	E	2	3	1/1/5/7	3/6/23/26	0/1/1/1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	2	NAG	C1-O5-C5	5.38	119.40	112.19
3	D	2	NAG	O5-C1-C2	3.71	117.03	111.29
3	D	2	NAG	C4-C3-C2	2.13	114.14	111.02

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	E	2	NAG	C1

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	D	1	NAG	C3-C2-N2-C7
3	D	2	NAG	C1-C2-N2-C7
3	D	2	NAG	C8-C7-N2-C2

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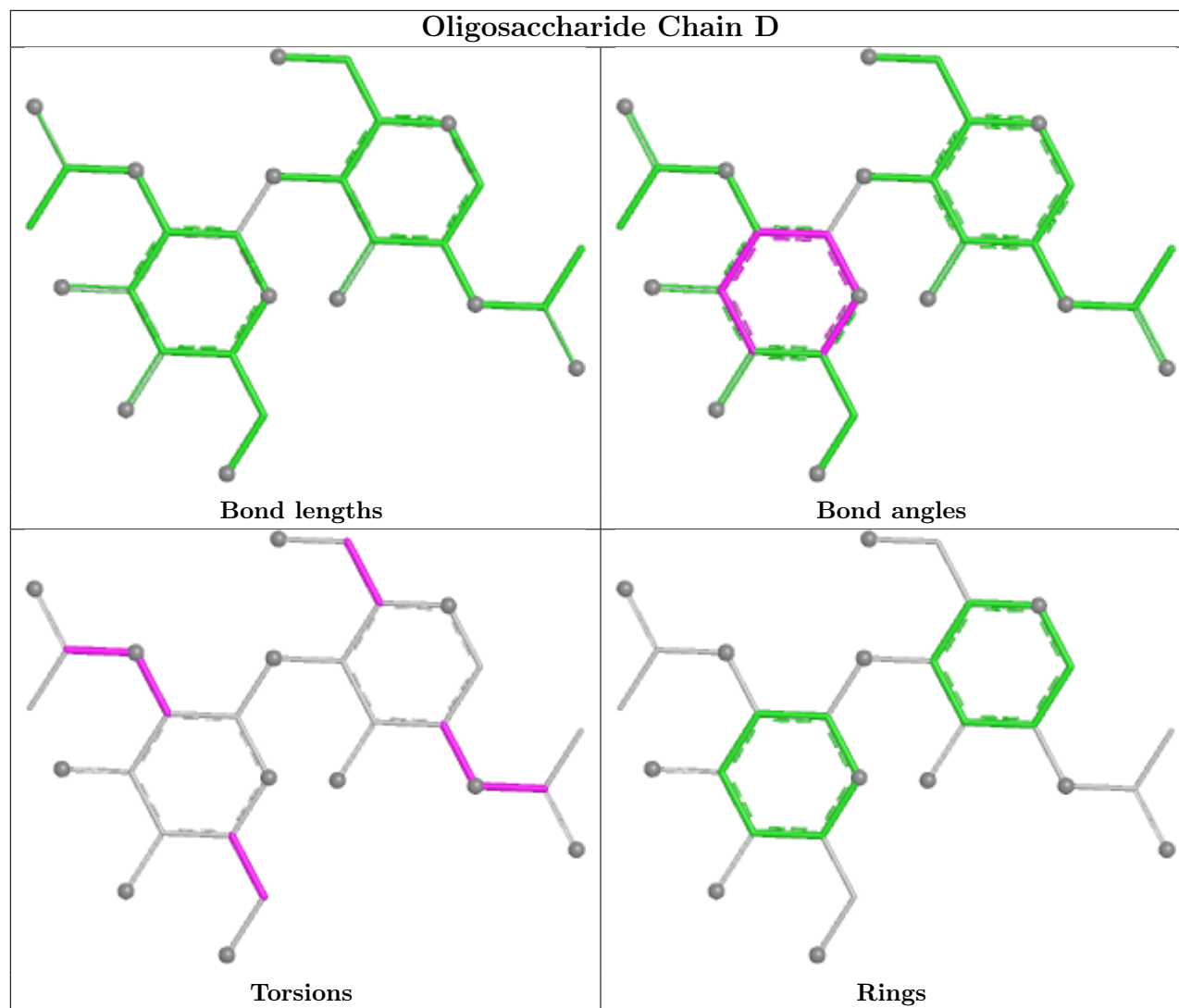
Mol	Chain	Res	Type	Atoms
3	D	2	NAG	O7-C7-N2-C2
3	E	2	NAG	C3-C2-N2-C7
3	E	2	NAG	C8-C7-N2-C2
3	E	2	NAG	O7-C7-N2-C2
3	D	1	NAG	C8-C7-N2-C2
3	D	1	NAG	O7-C7-N2-C2
3	D	1	NAG	O5-C5-C6-O6
3	D	2	NAG	O5-C5-C6-O6

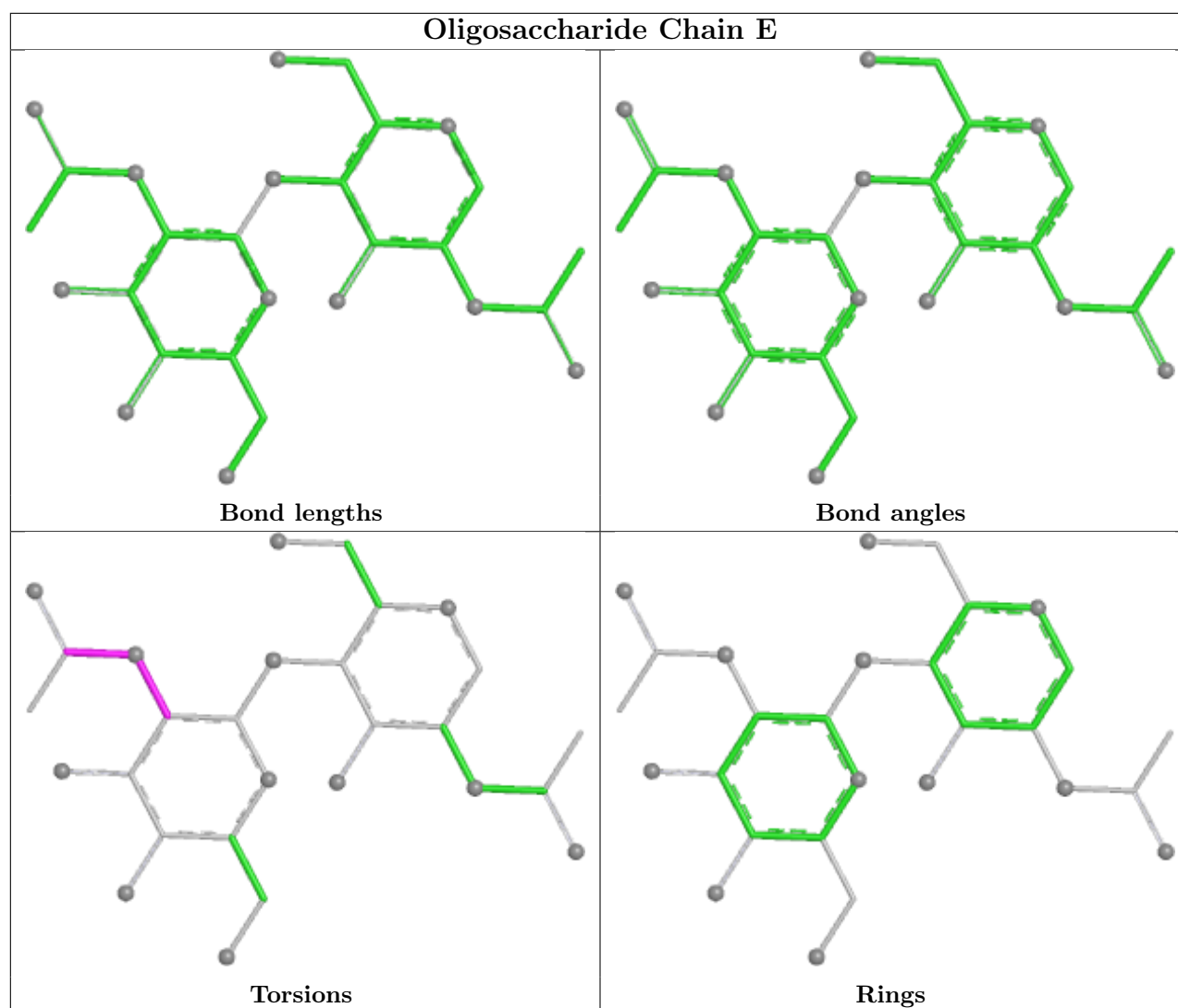
There are no ring outliers.

4 monomers are involved in 18 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	2	NAG	3	0
3	D	1	NAG	10	0
3	E	1	NAG	4	0
3	E	2	NAG	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.





5.6 Ligand geometry [i](#)

2 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	A	401	1	14,14,15	0.77	1 (7%)	17,19,21	0.73	1 (5%)
5	P15	C	401	-	19,19,19	0.73	1 (5%)	18,18,18	0.64	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	A	401	1	-	2/6/23/26	0/1/1/1
5	P15	C	401	-	-	12/17/17/17	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	401	NAG	C1-C2	2.45	1.55	1.52
5	C	401	P15	O6-C12	2.31	1.53	1.40

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	401	NAG	C1-O5-C5	2.47	115.49	112.19

There are no chirality outliers.

All (14) torsion outliers are listed below:

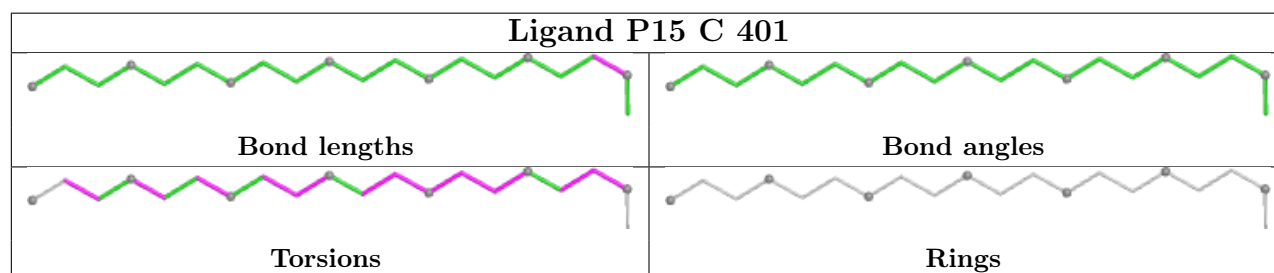
Mol	Chain	Res	Type	Atoms
5	C	401	P15	O5-C11-C12-O6
4	A	401	NAG	O5-C5-C6-O6
5	C	401	P15	O2-C5-C6-O3
5	C	401	P15	O3-C7-C8-O4
5	C	401	P15	C5-C6-O3-C7
4	A	401	NAG	C4-C5-C6-O6
5	C	401	P15	C9-C10-O5-C11
5	C	401	P15	C11-C12-O6-C13
5	C	401	P15	C7-C8-O4-C9
5	C	401	P15	C10-C9-O4-C8
5	C	401	P15	C3-C4-O2-C5
5	C	401	P15	C4-C3-O1-C2
5	C	401	P15	OXT-C1-C2-O1
5	C	401	P15	O5-C10-C9-O4

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	C	401	P15	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 all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

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	292/315 (92%)	0.18	14 (4%) 35 31	15, 42, 92, 128	0
1	B	293/315 (93%)	0.11	16 (5%) 30 27	17, 40, 84, 129	0
1	C	288/315 (91%)	0.24	24 (8%) 17 15	16, 43, 88, 124	0
2	F	38/54 (70%)	-0.15	1 (2%) 57 52	22, 37, 59, 85	0
2	G	39/54 (72%)	-0.24	2 (5%) 33 29	19, 32, 67, 79	0
2	H	40/54 (74%)	-0.29	0 100 100	19, 33, 59, 93	0
All	All	990/1107 (89%)	0.13	57 (5%) 29 25	15, 40, 87, 129	0

All (57) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	108	ASN	5.5
1	C	233	CYS	5.5
1	C	276	ASP	5.4
1	A	200	ARG	5.1
1	B	175	MET	5.0
1	B	201	ARG	5.0
1	A	111	ASN	4.5
1	C	110	ASN	4.3
1	B	282	LEU	3.9
1	C	115	ILE	3.9
1	A	268	ARG	3.8
1	A	123	VAL	3.8
1	B	108	ASN	3.7
1	C	106	THR	3.5
1	A	124	PRO	3.5
1	C	108	SER	3.4
1	C	126	ASP	3.3
1	B	189	LEU	3.3
1	C	114	VAL	3.2

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Mol	Chain	Res	Type	RSRZ
1	B	135	ASP	3.2
1	C	113	ASP	3.1
1	B	28	ASN	3.1
1	A	64	SER	3.1
1	A	279	ASP	3.0
1	A	131	ARG	3.0
1	B	111	ASN	2.9
1	C	127	ARG	2.9
1	C	197	ARG	2.8
1	B	127	VAL	2.8
1	C	91	THR	2.8
1	C	198	GLN	2.8
1	C	107	PHE	2.8
1	A	112	VAL	2.8
1	C	185	VAL	2.8
1	C	116	ASN	2.7
1	C	265	ARG	2.6
2	G	358	ASP	2.6
1	B	133	ASP	2.5
1	B	281	ASP	2.5
1	A	165	ASN	2.5
1	B	110	GLU	2.4
1	C	117	SER	2.4
1	A	125	LEU	2.4
1	B	70	ASP	2.4
2	F	322	LEU	2.4
1	C	93	GLU	2.4
1	C	111	VAL	2.3
1	B	190	VAL	2.3
1	C	263	PRO	2.3
1	A	110	GLU	2.2
1	C	121	ASP	2.2
1	A	107	PHE	2.2
1	C	281	LYS	2.2
1	C	304	THR	2.1
1	B	126	ARG	2.1
1	B	202	ARG	2.0
2	G	353	GLU	2.0

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

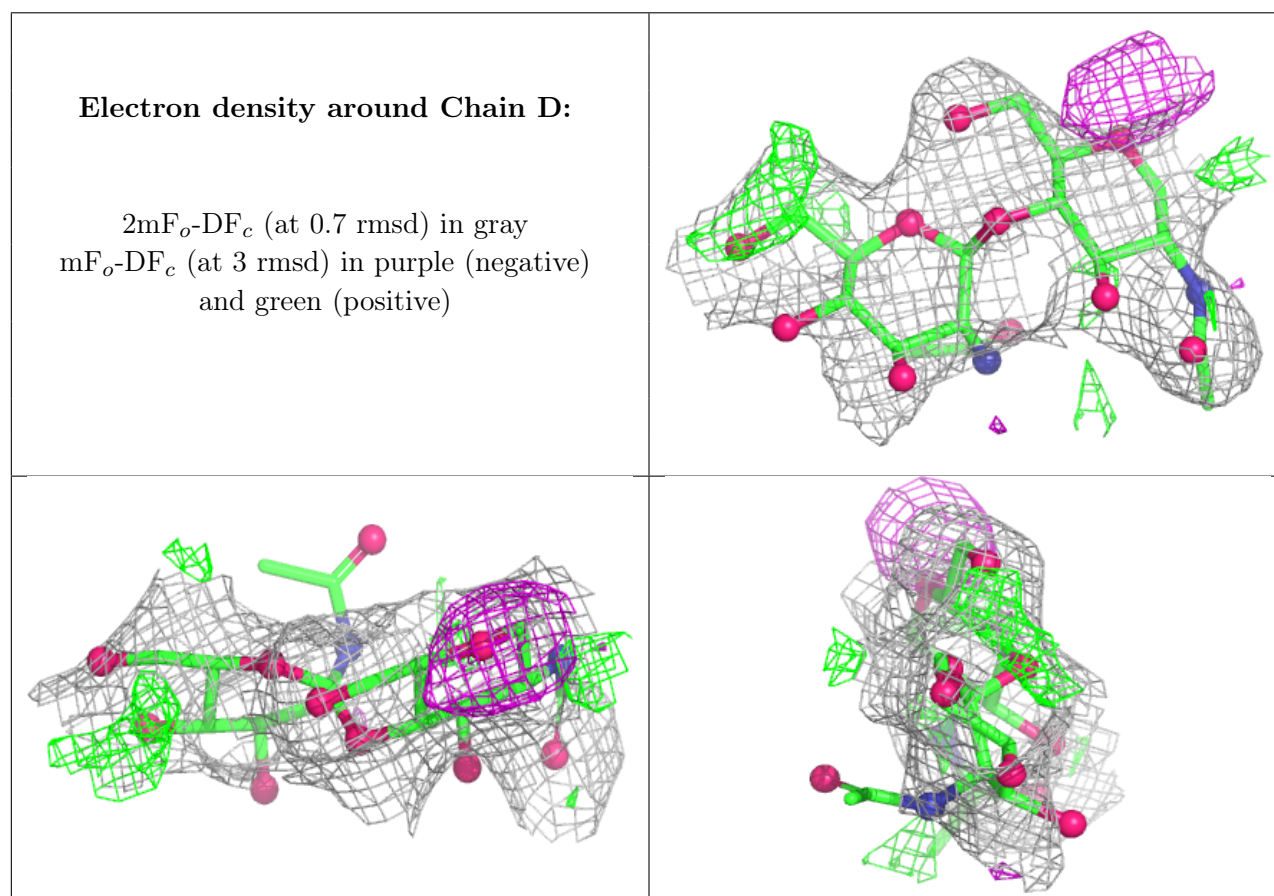
There are no non-standard protein/DNA/RNA residues in this entry.

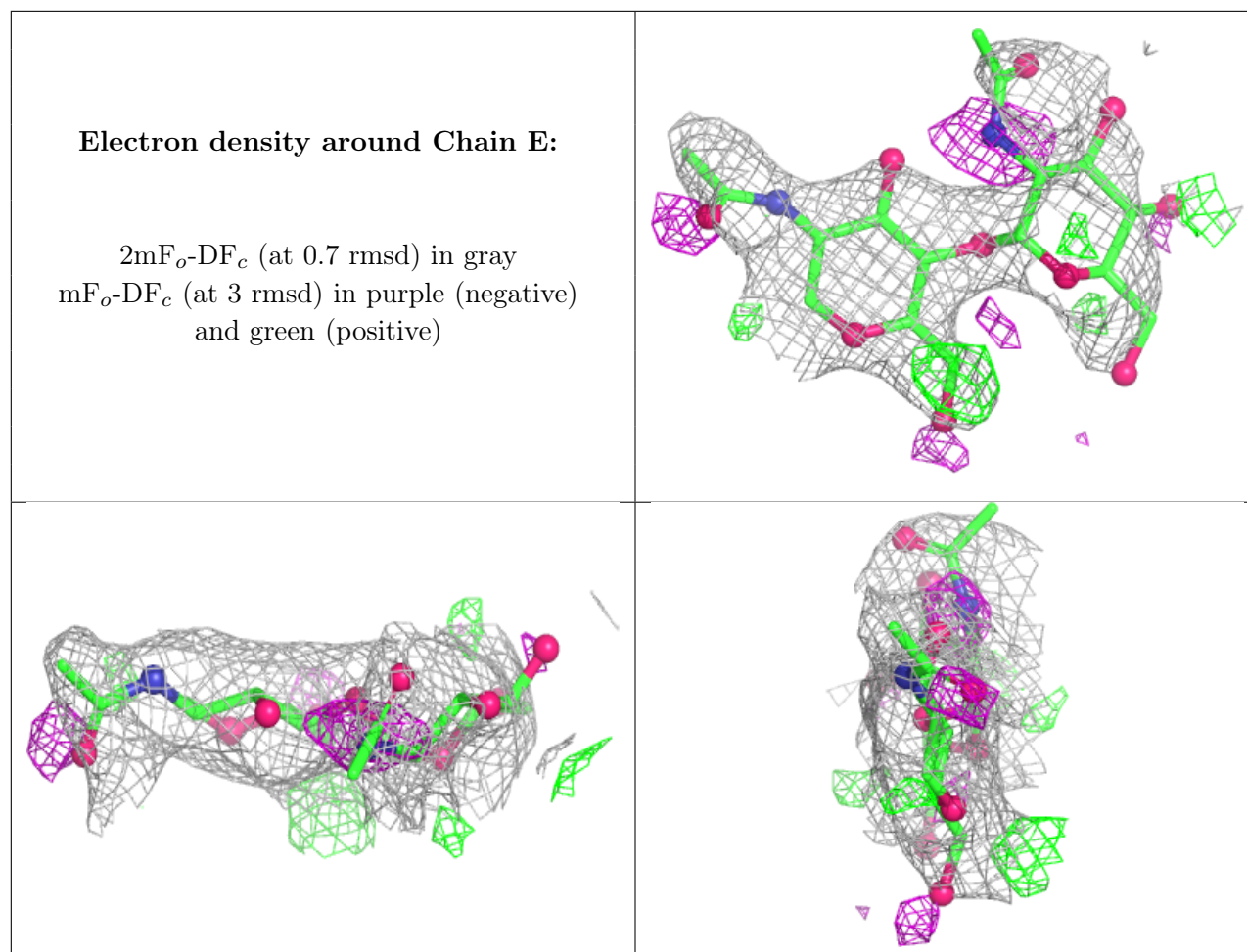
6.3 Carbohydrates ⓘ

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	NAG	D	1	14/15	-	-	44,71,87,91	0
3	NAG	D	2	14/15	-	-	56,94,99,101	0
3	NAG	E	1	14/15	-	-	38,73,83,83	0
3	NAG	E	2	14/15	-	-	65,87,99,103	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 ⓘ

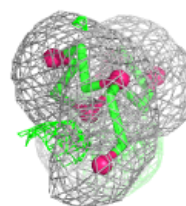
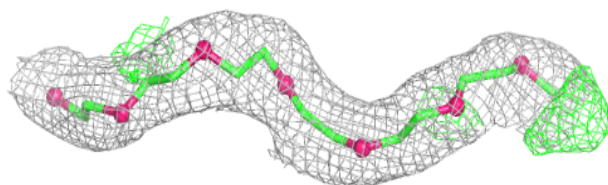
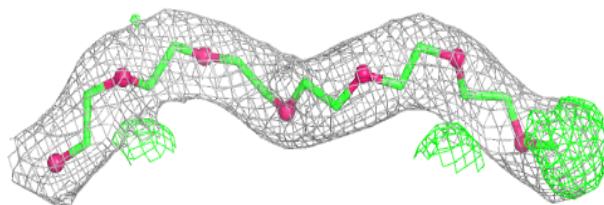
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists 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
4	NAG	A	401	14/15	0.76	0.13	66,73,77,78	0
5	P15	C	401	20/20	0.90	0.10	13,31,46,50	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around P15 C 401:

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



6.5 Other polymers [i](#)

There are no such residues in this entry.