



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 03:57 AM UTC

PDB ID : 2AOS / pdb_00002aos
Title : Protein-protein Interactions of protective signalling factor: Crystal structure of ternary complex involving signalling protein from goat (SPG-40), tetrasaccharide and a tripeptide Trp-pro-Trp at 2.9 Å resolution
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Deposited on : 2005-08-14
Resolution : 2.90 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)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

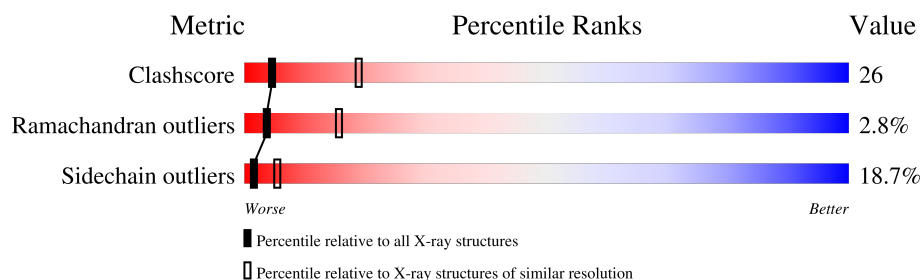
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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))
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	361	
2	D	3	
3	B	4	
4	C	3	

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 3107 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 Signaling protein from goat, SPG-40.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	361	Total	C	N	O	S	0	0	0
			2877	1836	508	524	9			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	33	VAL	ILE	SEE REMARK 999	GB 19526603
A	131	ALA	GLY	SEE REMARK 999	GB 19526603
A	205	ASN	GLN	SEE REMARK 999	GB 19526603
A	206	SER	GLU	SEE REMARK 999	GB 19526603
A	?	-	ASP	SEE REMARK 999	GB 19526603
A	361	ARG	GLU	SEE REMARK 999	GB 19526603

- Molecule 2 is a protein called Trp-Pro-Trp tripeptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	D	3	Total	C	N	O	0	0	0
			36	27	5	4			

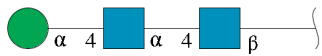
- Molecule 3 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	B	4	Total	C	N	O	0	0	0
			56	32	4	20			

- Molecule 4 is an oligosaccharide called alpha-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-

alpha-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



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

- Molecule 5 is water.

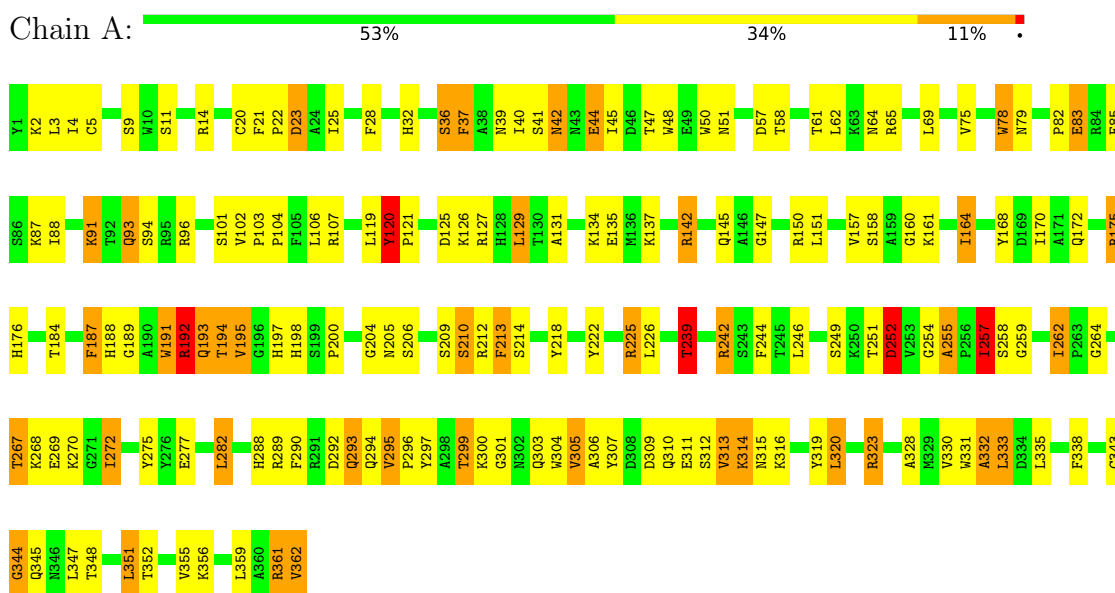
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	97	Total	O	0	0
			97	97		
5	D	2	Total	O	0	0
			2	2		

3 Residue-property plots

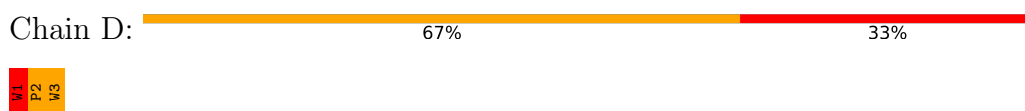
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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. 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.

Note EDS was not executed.

- Molecule 1: Signaling protein from goat, SPG-40



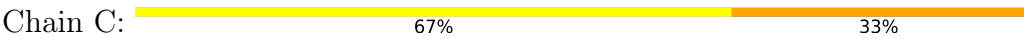
- Molecule 2: Trp-Pro-Trp tripeptide



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



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



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	62.87Å 66.51Å 107.62Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	56.00 – 2.90	Depositor
% Data completeness (in resolution range)	98.1 (56.00-2.90)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.11	Depositor
Refinement program	REFMAC 5.0	Depositor
R, R_{free}	0.192 , 0.239	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3107	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MAN, NDG, 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	0.95	0/2953	1.18	24/4001 (0.6%)
2	D	0.92	0/40	2.49	3/55 (5.5%)
All	All	0.95	0/2993	1.20	27/4056 (0.7%)

There are no bond length outliers.

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	1	TRP	CA-C-N	9.05	131.16	119.84
2	D	1	TRP	C-N-CA	9.05	131.16	119.84
1	A	187	PHE	N-CA-C	7.71	119.77	111.36
1	A	36	SER	N-CA-C	7.66	121.14	108.96
1	A	78	TRP	N-CA-C	7.62	119.66	111.36
1	A	120	TYR	CA-C-N	7.09	127.65	119.99
1	A	120	TYR	C-N-CA	7.09	127.65	119.99
1	A	36	SER	CA-C-N	6.92	134.76	121.54
1	A	36	SER	C-N-CA	6.92	134.76	121.54
1	A	129	LEU	N-CA-C	-6.50	104.11	111.07
1	A	264	GLY	N-CA-C	-6.46	103.33	112.37
2	D	2	PRO	N-CA-C	6.22	125.29	112.47
1	A	102	VAL	N-CA-C	5.97	121.77	108.88
1	A	239	THR	N-CA-C	-5.88	105.41	112.59
1	A	37	PHE	CA-CB-CG	-5.75	108.05	113.80
1	A	257	ILE	N-CA-C	5.75	116.56	108.17
1	A	225	ARG	N-CA-C	-5.69	104.98	111.07
1	A	25	ILE	N-CA-C	5.55	116.73	108.23
1	A	268	LYS	N-CA-C	5.54	119.21	111.90
1	A	213	PHE	N-CA-C	-5.40	106.88	112.93
1	A	192	ARG	N-CA-C	5.38	122.26	110.80
1	A	333	LEU	N-CA-C	-5.37	105.53	111.71

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	314	LYS	N-CA-C	-5.36	105.52	111.36
1	A	170	ILE	N-CA-C	5.35	116.08	110.62
1	A	37	PHE	N-CA-C	5.29	122.08	110.80
1	A	290	PHE	N-CA-C	-5.18	101.52	109.76
1	A	305	VAL	N-CA-C	5.09	115.31	107.77

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2877	0	2818	151	0
2	D	36	0	29	12	0
3	B	56	0	49	7	0
4	C	39	0	33	4	0
5	A	97	0	0	9	0
5	D	2	0	0	0	0
All	All	3107	0	2929	155	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 26.

All (155) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:82:PRO:HG3	2:D:1:TRP:HB2	1.40	1.03
1:A:204:GLY:HA2	1:A:292:ASP:HB3	1.41	1.01
1:A:78:TRP:CD1	2:D:1:TRP:HZ2	1.79	1.01
1:A:82:PRO:HD3	2:D:1:TRP:HD1	1.27	0.98
1:A:78:TRP:HD1	2:D:1:TRP:HZ2	1.05	0.93
1:A:78:TRP:HD1	2:D:1:TRP:CZ2	1.88	0.91
1:A:252:ASP:HB3	5:A:408:HOH:O	1.72	0.90
1:A:39:ASN:ND2	4:C:1:NAG:C1	2.38	0.86
1:A:192:ARG:NE	1:A:192:ARG:HA	1.90	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:192:ARG:HA	1:A:192:ARG:HE	1.38	0.84
1:A:39:ASN:HD21	4:C:1:NAG:C1	1.92	0.83
1:A:82:PRO:HD3	2:D:1:TRP:CD1	2.14	0.82
1:A:78:TRP:CD1	2:D:1:TRP:CZ2	2.64	0.82
1:A:200:PRO:HB2	1:A:293:GLN:HG2	1.61	0.80
1:A:142:ARG:HB2	1:A:142:ARG:NH1	1.97	0.79
1:A:332:ALA:CB	1:A:335:LEU:HD12	2.15	0.77
1:A:323:ARG:HH11	1:A:323:ARG:HG2	1.50	0.76
1:A:91:LYS:HB2	1:A:94:SER:OG	1.87	0.74
1:A:44:GLU:HG2	1:A:101:SER:HB2	1.69	0.73
1:A:61:THR:O	1:A:64:ASN:HB2	1.90	0.72
1:A:323:ARG:HG2	1:A:323:ARG:NH1	2.05	0.69
5:A:437:HOH:O	2:D:1:TRP:HZ3	1.76	0.69
1:A:137:LYS:HE3	1:A:150:ARG:NH2	2.08	0.67
1:A:39:ASN:CG	4:C:1:NAG:C1	2.68	0.67
1:A:78:TRP:HZ3	1:A:331:TRP:CZ2	2.13	0.65
5:A:424:HOH:O	3:B:4:NAG:C8	2.45	0.65
1:A:120:TYR:CE1	1:A:158:SER:HB2	2.32	0.65
5:A:424:HOH:O	3:B:4:NAG:H81	1.97	0.65
1:A:192:ARG:HD3	1:A:193:GLN:H	1.62	0.64
1:A:191:TRP:HB2	2:D:3:TRP:CZ3	2.34	0.62
1:A:209:SER:O	1:A:210:SER:C	2.43	0.62
1:A:332:ALA:HB2	1:A:335:LEU:HD12	1.81	0.61
1:A:343:CYS:O	1:A:344:GLY:C	2.42	0.61
1:A:212:ARG:O	1:A:212:ARG:HD2	2.01	0.61
1:A:83:GLU:HG2	5:A:394:HOH:O	2.01	0.59
1:A:267:THR:HB	1:A:277:GLU:OE1	2.03	0.59
1:A:361:ARG:O	1:A:362:VAL:HG12	2.03	0.59
1:A:330:VAL:HG21	1:A:352:THR:HG23	1.84	0.58
1:A:323:ARG:HH11	1:A:323:ARG:CG	2.16	0.58
1:A:289:ARG:NH2	1:A:309:ASP:OD2	2.36	0.58
1:A:5:CYS:HB3	1:A:333:LEU:HG	1.85	0.57
1:A:239:THR:HG22	1:A:335:LEU:HB2	1.87	0.57
1:A:272:ILE:HG21	3:B:1:NAG:H83	1.86	0.57
1:A:39:ASN:CG	1:A:40:ILE:N	2.63	0.56
1:A:39:ASN:OD1	4:C:1:NAG:C1	2.53	0.56
1:A:331:TRP:O	1:A:332:ALA:HB3	2.05	0.56
1:A:188:HIS:H	1:A:198:HIS:HA	1.71	0.56
1:A:57:ASP:O	1:A:61:THR:HG23	2.07	0.55
1:A:161:LYS:HE2	1:A:218:TYR:OH	2.07	0.55
1:A:239:THR:HG23	1:A:239:THR:O	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:83:GLU:H	1:A:83:GLU:CD	2.13	0.55
1:A:160:GLY:O	1:A:164:ILE:HG13	2.08	0.54
1:A:50:TRP:CD1	1:A:50:TRP:H	2.24	0.54
1:A:296:PRO:HD2	1:A:307:TYR:O	2.07	0.54
1:A:87:LYS:O	1:A:91:LYS:HD2	2.08	0.54
1:A:239:THR:HG21	1:A:332:ALA:O	2.08	0.54
1:A:142:ARG:HB2	1:A:142:ARG:HH11	1.70	0.53
1:A:191:TRP:CZ3	1:A:192:ARG:HG2	2.44	0.53
1:A:304:TRP:CZ3	1:A:306:ALA:HB2	2.43	0.53
1:A:313:VAL:HG13	1:A:355:VAL:HG22	1.90	0.53
1:A:262:ILE:HG13	1:A:303:GLN:NE2	2.24	0.52
1:A:137:LYS:HE3	1:A:150:ARG:CZ	2.39	0.52
1:A:239:THR:O	1:A:239:THR:CG2	2.56	0.52
1:A:142:ARG:HB2	1:A:142:ARG:CZ	2.39	0.52
1:A:164:ILE:HA	1:A:168:TYR:HD2	1.75	0.52
1:A:120:TYR:CD1	1:A:120:TYR:N	2.74	0.52
1:A:2:LYS:HB3	1:A:4:ILE:CD1	2.40	0.51
1:A:96:ARG:HB3	1:A:96:ARG:NH1	2.25	0.51
1:A:142:ARG:O	1:A:145:GLN:HB2	2.10	0.51
1:A:39:ASN:HB3	1:A:47:THR:O	2.10	0.51
1:A:2:LYS:HB3	1:A:4:ILE:HD11	1.91	0.51
1:A:213:PHE:HA	1:A:218:TYR:CD2	2.45	0.51
5:A:437:HOH:O	2:D:1:TRP:CZ3	2.54	0.51
1:A:172:GLN:O	1:A:175:ARG:HG3	2.10	0.51
1:A:191:TRP:HB3	2:D:3:TRP:CE3	2.46	0.50
1:A:131:ALA:O	1:A:135:GLU:HG3	2.11	0.50
1:A:299:THR:HA	1:A:303:GLN:O	2.11	0.49
1:A:188:HIS:CE1	1:A:194:THR:O	2.65	0.49
1:A:269:GLU:OE1	3:B:1:NAG:O3	2.24	0.49
1:A:62:LEU:C	1:A:64:ASN:H	2.19	0.48
1:A:22:PRO:HB2	1:A:58:THR:HG22	1.94	0.48
1:A:93:GLN:HG2	1:A:94:SER:N	2.25	0.48
1:A:195:VAL:CG1	1:A:304:TRP:CE2	2.97	0.47
1:A:11:SER:HB2	1:A:14:ARG:NH1	2.29	0.47
1:A:320:LEU:HD11	1:A:328:ALA:HB2	1.96	0.47
1:A:39:ASN:HB2	1:A:48:TRP:CE3	2.49	0.47
1:A:204:GLY:O	1:A:206:SER:N	2.47	0.47
1:A:39:ASN:CG	1:A:40:ILE:H	2.23	0.47
1:A:191:TRP:CG	1:A:192:ARG:H	2.33	0.47
1:A:44:GLU:HG2	1:A:101:SER:CB	2.43	0.47
1:A:313:VAL:HG13	1:A:355:VAL:CG2	2.45	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:242:ARG:HH21	1:A:272:ILE:HD11	1.80	0.47
1:A:41:SER:O	1:A:42:ASN:HB2	2.15	0.46
1:A:332:ALA:HB1	1:A:335:LEU:CG	2.45	0.46
1:A:200:PRO:CB	1:A:293:GLN:HG2	2.39	0.46
1:A:288:HIS:HB2	1:A:297:TYR:CZ	2.50	0.46
1:A:320:LEU:C	1:A:320:LEU:HD23	2.41	0.45
1:A:191:TRP:CE3	1:A:192:ARG:HG2	2.51	0.45
1:A:93:GLN:HB2	5:A:398:HOH:O	2.16	0.45
1:A:78:TRP:CH2	5:A:395:HOH:O	2.56	0.45
1:A:189:GLY:HA3	1:A:191:TRP:CD1	2.51	0.45
1:A:192:ARG:CD	5:A:444:HOH:O	2.65	0.45
1:A:157:VAL:HG12	1:A:158:SER:O	2.17	0.45
1:A:11:SER:CB	1:A:14:ARG:NH1	2.80	0.45
1:A:195:VAL:HG13	1:A:304:TRP:CZ2	2.52	0.45
1:A:195:VAL:HG11	1:A:304:TRP:CE2	2.51	0.45
1:A:310:GLN:O	1:A:314:LYS:CG	2.65	0.45
1:A:78:TRP:HB2	3:B:1:NAG:O5	2.17	0.45
1:A:332:ALA:HB1	1:A:335:LEU:HB2	1.98	0.44
1:A:37:PHE:CD1	1:A:37:PHE:N	2.78	0.44
1:A:78:TRP:HB2	3:B:1:NAG:O6	2.17	0.44
1:A:119:LEU:HA	1:A:119:LEU:HD23	1.45	0.44
1:A:62:LEU:C	1:A:64:ASN:N	2.75	0.44
1:A:120:TYR:HA	1:A:121:PRO:HD3	1.64	0.44
1:A:134:LYS:HG3	1:A:176:HIS:NE2	2.33	0.44
1:A:269:GLU:OE2	1:A:272:ILE:HD13	2.17	0.44
1:A:191:TRP:CB	2:D:3:TRP:CE3	3.00	0.44
1:A:257:ILE:HG23	1:A:259:GLY:H	1.83	0.44
1:A:187:PHE:HB2	1:A:198:HIS:O	2.17	0.44
1:A:79:ASN:ND2	3:B:2:NAG:C7	2.81	0.43
1:A:226:LEU:HD13	1:A:226:LEU:HA	1.80	0.43
1:A:145:GLN:C	1:A:147:GLY:H	2.26	0.43
1:A:28:PHE:O	1:A:356:LYS:NZ	2.52	0.43
1:A:78:TRP:NE1	1:A:119:LEU:HD13	2.33	0.43
1:A:96:ARG:HB3	1:A:96:ARG:CZ	2.49	0.43
1:A:103:PRO:HB2	1:A:104:PRO:CD	2.49	0.43
1:A:195:VAL:CG1	1:A:304:TRP:CZ2	3.02	0.43
1:A:246:LEU:HA	1:A:257:ILE:HD13	2.00	0.43
1:A:282:LEU:HA	1:A:282:LEU:HD12	1.76	0.43
1:A:244:PHE:HB3	1:A:257:ILE:HD12	2.00	0.43
1:A:320:LEU:CD1	1:A:328:ALA:HB2	2.49	0.43
1:A:204:GLY:C	1:A:206:SER:N	2.76	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:210:SER:O	1:A:210:SER:OG	2.32	0.42
1:A:310:GLN:O	1:A:314:LYS:HG3	2.19	0.42
1:A:319:TYR:CZ	1:A:323:ARG:HD2	2.54	0.42
1:A:195:VAL:HG13	1:A:304:TRP:CH2	2.55	0.42
1:A:254:GLY:O	1:A:255:ALA:C	2.62	0.42
1:A:20:CYS:HB2	1:A:338:PHE:CZ	2.55	0.41
1:A:106:LEU:HD23	1:A:151:LEU:HD22	2.01	0.41
1:A:212:ARG:C	1:A:214:SER:H	2.28	0.41
1:A:275:TYR:CG	1:A:351:LEU:HD22	2.56	0.41
1:A:21:PHE:C	1:A:23:ASP:N	2.76	0.41
1:A:312:SER:O	1:A:316:LYS:HG3	2.20	0.41
1:A:125:ASP:O	1:A:126:LYS:C	2.64	0.41
1:A:222:TYR:O	1:A:226:LEU:HD23	2.21	0.41
1:A:197:HIS:O	1:A:200:PRO:HD3	2.21	0.41
1:A:300:LYS:O	1:A:301:GLY:C	2.62	0.41
1:A:78:TRP:CD1	1:A:119:LEU:HD13	2.56	0.40
1:A:297:TYR:HA	1:A:305:VAL:O	2.21	0.40
1:A:332:ALA:CB	1:A:335:LEU:HB2	2.50	0.40
1:A:212:ARG:O	1:A:212:ARG:CD	2.68	0.40
1:A:51:ASN:OD1	1:A:51:ASN:N	2.54	0.40
1:A:257:ILE:HD13	1:A:257:ILE:HA	1.82	0.40
1:A:4:ILE:HG13	1:A:32:HIS:HB2	2.04	0.40
1:A:295:VAL:HA	1:A:296:PRO:HD3	1.79	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	359/361 (99%)	320 (89%)	29 (8%)	10 (3%)	4 15
2	D	1/3 (33%)	1 (100%)	0	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	360/364 (99%)	321 (89%)	29 (8%)	10 (3%)	4	15

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	191	TRP
1	A	192	ARG
1	A	205	ASN
1	A	332	ALA
1	A	344	GLY
1	A	252	ASP
1	A	345	GLN
1	A	120	TYR
1	A	255	ALA
1	A	249	SER

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	302/302 (100%)	248 (82%)	54 (18%)	2	6
2	D	3/3 (100%)	0	3 (100%)	0	0
All	All	305/305 (100%)	248 (81%)	57 (19%)	1	5

All (57) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	LEU
1	A	9	SER
1	A	23	ASP
1	A	36	SER
1	A	42	ASN
1	A	44	GLU
1	A	45	ILE
1	A	65	ARG

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Mol	Chain	Res	Type
1	A	69	LEU
1	A	75	VAL
1	A	83	GLU
1	A	85	PHE
1	A	88	ILE
1	A	91	LYS
1	A	93	GLN
1	A	107	ARG
1	A	127	ARG
1	A	129	LEU
1	A	142	ARG
1	A	164	ILE
1	A	175	ARG
1	A	184	THR
1	A	192	ARG
1	A	193	GLN
1	A	194	THR
1	A	195	VAL
1	A	210	SER
1	A	225	ARG
1	A	239	THR
1	A	242	ARG
1	A	251	THR
1	A	252	ASP
1	A	257	ILE
1	A	258	SER
1	A	262	ILE
1	A	267	THR
1	A	270	LYS
1	A	272	ILE
1	A	282	LEU
1	A	293	GLN
1	A	294	GLN
1	A	295	VAL
1	A	299	THR
1	A	311	GLU
1	A	313	VAL
1	A	315	ASN
1	A	320	LEU
1	A	323	ARG
1	A	347	LEU
1	A	348	THR

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Mol	Chain	Res	Type
1	A	351	LEU
1	A	359	LEU
1	A	361	ARG
1	A	362	VAL
2	D	1	TRP
2	D	2	PRO
2	D	3	TRP

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

Mol	Chain	Res	Type
1	A	42	ASN
1	A	60	ASN
1	A	79	ASN
1	A	188	HIS
1	A	205	ASN
1	A	288	HIS
1	A	303	GLN
1	A	310	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

7 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	B	1	3	14,14,15	0.74	0	17,19,21	2.41	7 (41%)
3	NAG	B	2	3	14,14,15	1.43	3 (21%)	17,19,21	2.15	6 (35%)
3	NAG	B	3	3	14,14,15	1.41	1 (7%)	17,19,21	2.45	5 (29%)
3	NAG	B	4	3	14,14,15	0.92	1 (7%)	17,19,21	1.83	4 (23%)
4	NAG	C	1	4	14,14,15	0.93	1 (7%)	17,19,21	3.97	4 (23%)
4	NDG	C	2	4	14,14,15	0.76	0	17,19,21	3.88	11 (64%)
4	MAN	C	3	4	11,11,12	1.07	0	15,15,17	2.32	5 (33%)

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	B	1	3	-	3/6/23/26	0/1/1/1
3	NAG	B	2	3	-	5/6/23/26	0/1/1/1
3	NAG	B	3	3	-	3/6/23/26	0/1/1/1
3	NAG	B	4	3	-	2/6/23/26	0/1/1/1
4	NAG	C	1	4	-	0/6/23/26	0/1/1/1
4	NDG	C	2	4	-	2/6/23/26	0/1/1/1
4	MAN	C	3	4	-	2/2/19/22	0/1/1/1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	3	NAG	C1-C2	4.10	1.57	1.52
3	B	2	NAG	C1-C2	2.91	1.56	1.52
3	B	2	NAG	O5-C5	-2.57	1.38	1.43
4	C	1	NAG	O5-C1	-2.34	1.39	1.43
3	B	2	NAG	C4-C3	2.19	1.58	1.52
3	B	4	NAG	O5-C5	-2.18	1.39	1.43

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	1	NAG	C2-N2-C7	-14.94	102.89	122.90
4	C	2	NDG	C2-N2-C7	-10.00	109.50	122.90
3	B	3	NAG	O5-C1-C2	7.23	122.48	111.29
4	C	2	NDG	C1-C2-N2	6.70	120.99	110.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	2	NDG	C1-O5-C5	6.61	121.04	112.19
4	C	3	MAN	C1-O5-C5	5.46	119.51	112.19
3	B	1	NAG	C2-N2-C7	4.89	129.46	122.90
3	B	1	NAG	C1-O5-C5	4.26	117.90	112.19
3	B	2	NAG	C2-N2-C7	4.16	128.47	122.90
4	C	3	MAN	O5-C1-C2	4.15	120.68	110.79
3	B	2	NAG	O5-C1-C2	4.12	117.67	111.29
3	B	4	NAG	O5-C5-C4	-4.04	100.99	110.83
3	B	3	NAG	C2-N2-C7	4.01	128.27	122.90
4	C	2	NDG	C4-C3-C2	3.93	116.78	111.02
4	C	3	MAN	O5-C5-C4	3.25	118.73	110.83
3	B	1	NAG	C3-C4-C5	3.22	116.08	110.23
4	C	1	NAG	O5-C1-C2	-3.22	106.32	111.29
3	B	2	NAG	C8-C7-N2	-3.18	110.85	116.12
3	B	4	NAG	C1-O5-C5	3.16	116.42	112.19
3	B	1	NAG	O5-C5-C4	3.14	118.47	110.83
3	B	2	NAG	C1-C2-N2	-3.10	105.55	110.43
3	B	1	NAG	C4-C3-C2	-3.09	106.49	111.02
4	C	2	NDG	O7-C7-N2	2.99	127.27	121.98
3	B	4	NAG	C4-C3-C2	2.99	115.39	111.02
3	B	3	NAG	C3-C4-C5	-2.95	104.88	110.23
4	C	3	MAN	O4-C4-C5	2.94	116.56	109.32
4	C	2	NDG	O5-C5-C4	2.86	117.79	110.83
3	B	3	NAG	O5-C5-C4	2.71	117.41	110.83
4	C	1	NAG	C3-C4-C5	-2.63	105.46	110.23
4	C	2	NDG	C6-C5-C4	-2.51	106.86	113.02
4	C	1	NAG	C8-C7-N2	2.47	120.21	116.12
3	B	4	NAG	C1-C2-N2	-2.46	106.55	110.43
4	C	2	NDG	O3-C3-C2	-2.42	104.38	109.40
3	B	1	NAG	O5-C1-C2	-2.31	107.72	111.29
3	B	1	NAG	C1-C2-N2	2.31	114.07	110.43
4	C	2	NDG	O3-C3-C4	-2.23	105.12	110.38
4	C	3	MAN	O5-C5-C6	-2.16	103.47	107.66
4	C	2	NDG	C3-C4-C5	2.15	114.13	110.23
3	B	2	NAG	O7-C7-N2	2.10	125.69	121.98
3	B	3	NAG	O7-C7-N2	2.09	125.67	121.98
4	C	2	NDG	O7-C7-C8	-2.09	118.34	122.05
3	B	2	NAG	C4-C3-C2	2.06	114.03	111.02

There are no chirality outliers.

All (17) torsion outliers are listed below:

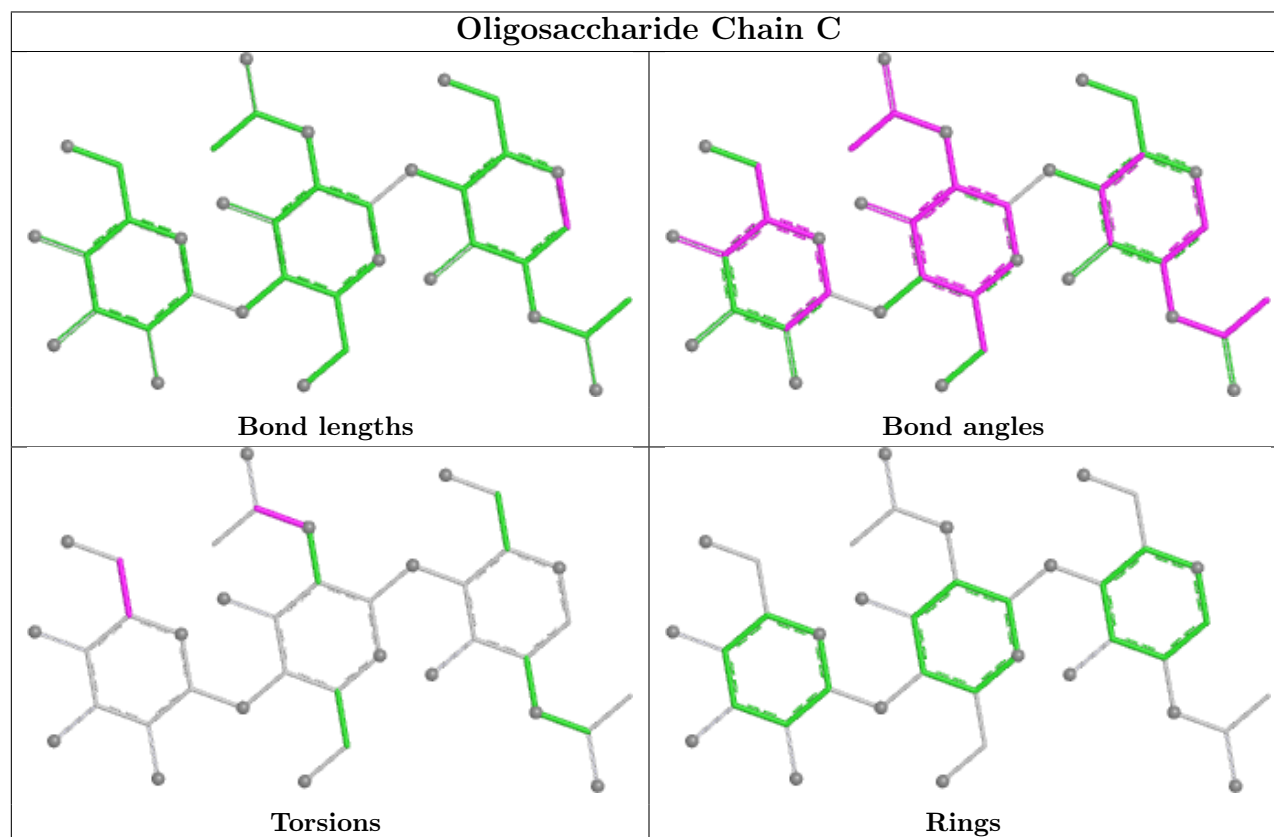
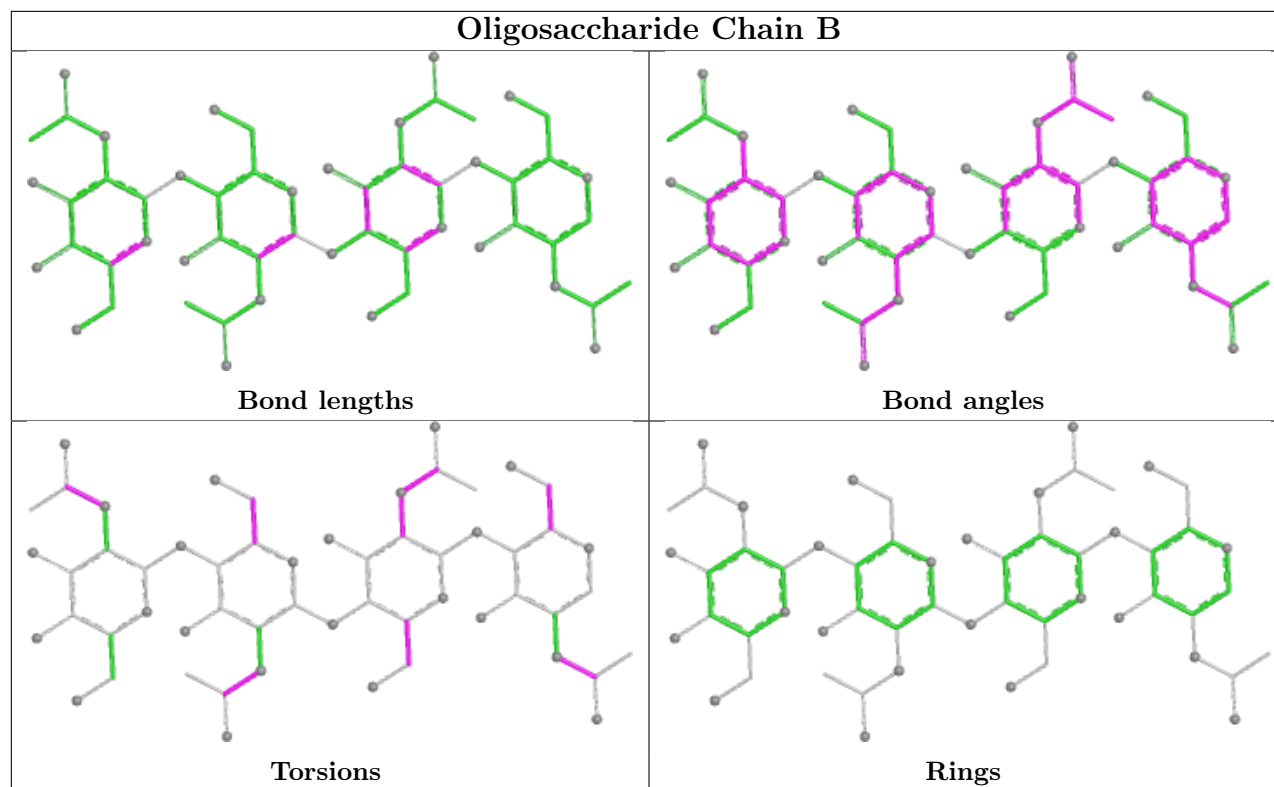
Mol	Chain	Res	Type	Atoms
3	B	4	NAG	O7-C7-N2-C2
3	B	4	NAG	C8-C7-N2-C2
3	B	2	NAG	O5-C5-C6-O6
4	C	3	MAN	C4-C5-C6-O6
3	B	2	NAG	C4-C5-C6-O6
4	C	3	MAN	O5-C5-C6-O6
3	B	3	NAG	C8-C7-N2-C2
4	C	2	NDG	C8-C7-N2-C2
4	C	2	NDG	O7-C7-N2-C2
3	B	1	NAG	C4-C5-C6-O6
3	B	2	NAG	C8-C7-N2-C2
3	B	2	NAG	O7-C7-N2-C2
3	B	3	NAG	O7-C7-N2-C2
3	B	3	NAG	C4-C5-C6-O6
3	B	2	NAG	C3-C2-N2-C7
3	B	1	NAG	O7-C7-N2-C2
3	B	1	NAG	O5-C5-C6-O6

There are no ring outliers.

4 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	2	NAG	1	0
3	B	1	NAG	4	0
3	B	4	NAG	2	0
4	C	1	NAG	4	0

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



5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.