



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 03:51 PM UTC

PDB ID : 1ZAG / pdb_00001zag
Title : HUMAN ZINC-ALPHA-2-GLYCOPROTEIN
Authors : Chirino, A.J.; Sanchez, L.M.; Bjorkman, P.J.
Deposited on : 1999-02-02
Resolution : 2.80 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

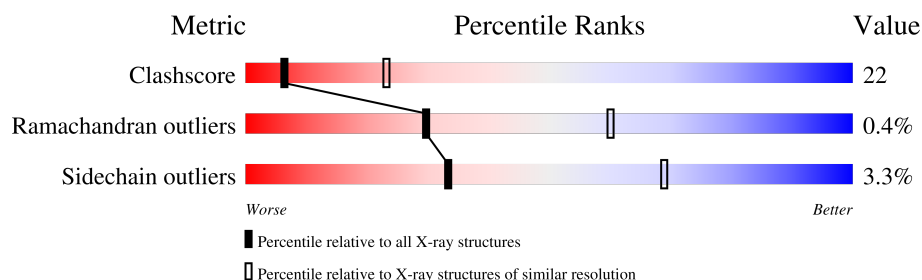
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)

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	274	59% 35% 5%
1	B	274	59% 39% .
1	C	274	57% 38% .
1	D	274	55% 39% . .
2	E	7	29% 43% 29%
3	F	9	22% 67% 11%
4	G	2	50% 50%
5	H	3	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
6	NAG	C	310	X	-	-	-

2 Entry composition [i](#)

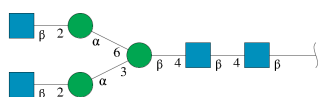
There are 7 unique types of molecules in this entry. The entry contains 9256 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 PROTEIN (ZINC-ALPHA-2-GLYCOPROTEIN).

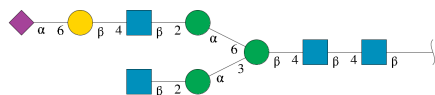
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	273	Total	C	N	O	S	0	0	0
			2232	1425	379	421	7			
1	B	274	Total	C	N	O	S	0	0	0
			2239	1428	380	424	7			
1	C	273	Total	C	N	O	S	0	0	0
			2231	1424	379	421	7			
1	D	263	Total	C	N	O	S	0	0	0
			2162	1382	368	405	7			

- Molecule 2 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



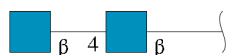
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	E	7	Total	C	N	O	0	0	0
			89	50	4	35			

- Molecule 3 is an oligosaccharide called N-acetyl-alpha-neuraminic acid-(2-6)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-6)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-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.



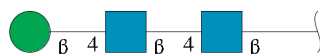
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	F	9	Total	C	N	O	0	0	0
			120	67	5	48			

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



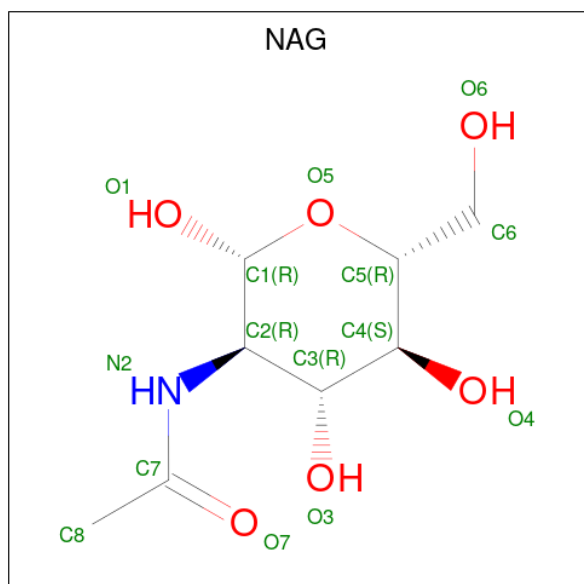
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	G	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 5 is an oligosaccharide called 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
5	H	3	Total	C	N	O	0	0	0
			39	22	2	15			

- Molecule 6 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



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

- Molecule 7 is water.

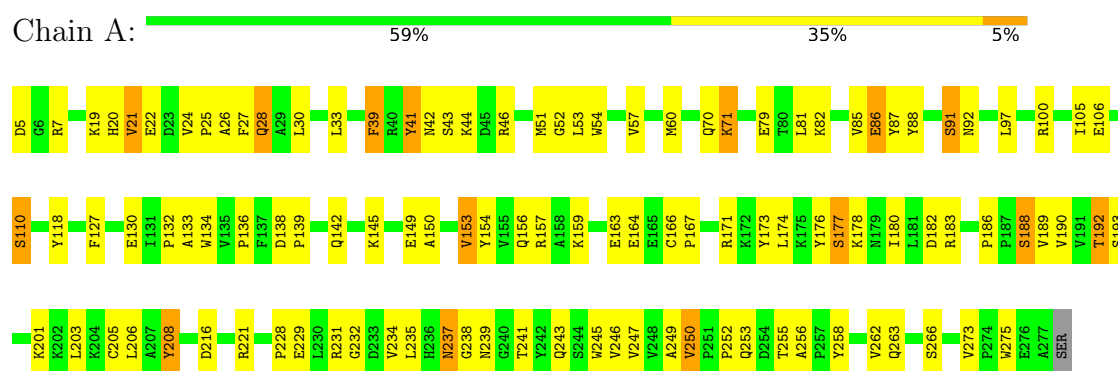
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	1	Total	O	0	0
			1	1		
7	B	1	Total	O	0	0
			1	1		
7	C	1	Total	O	0	0
			1	1		
7	D	1	Total	O	0	0
			1	1		

3 Residue-property plots

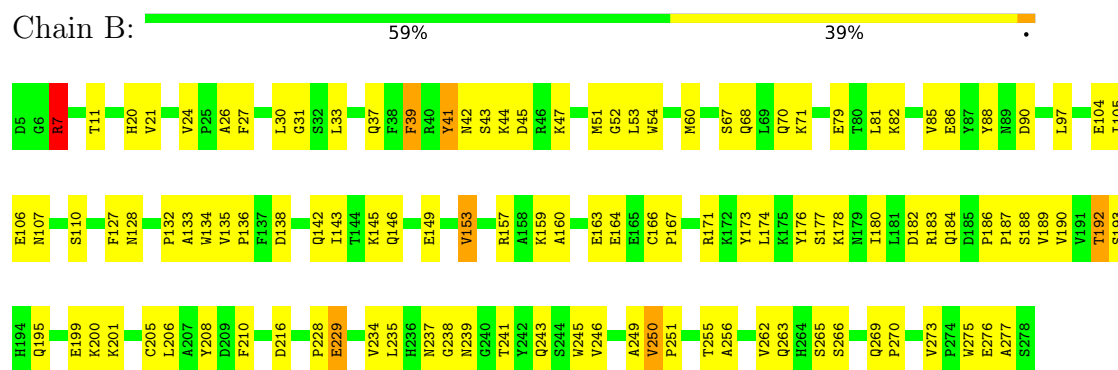
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded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

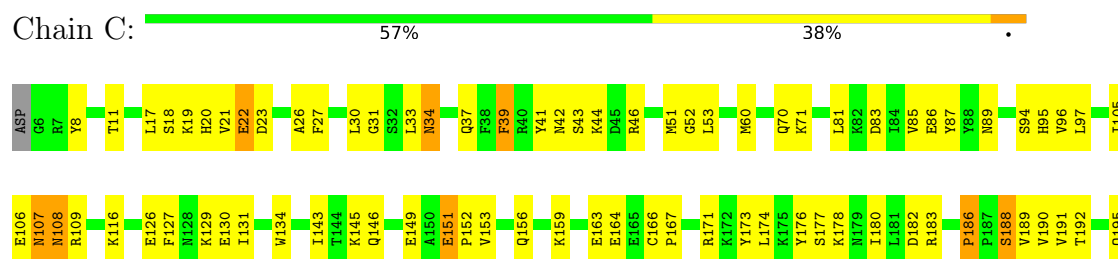
• Molecule 1: PROTEIN (ZINC-ALPHA-2-GLYCOPROTEIN)



• Molecule 1: PROTEIN (ZINC-ALPHA-2-GLYCOPROTEIN)

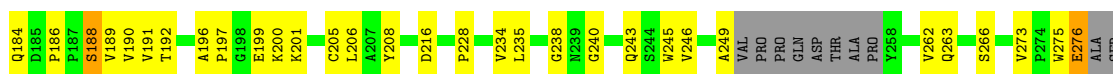
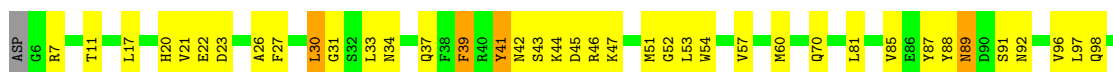


• Molecule 1: PROTEIN (ZINC-ALPHA-2-GLYCOPROTEIN)

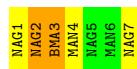




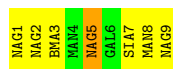
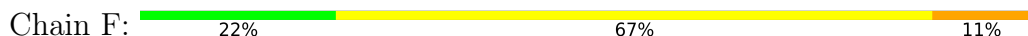
• Molecule 1: PROTEIN (ZINC-ALPHA-2-GLYCOPROTEIN)



• Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



• Molecule 3: N-acetyl-alpha-neuraminic acid-(2-6)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-6)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-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



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



• Molecule 5: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



MAG1
MAG2
BMA3

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 2	Depositor
Cell constants a, b, c, α , β , γ	105.60Å 131.70Å 118.50Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.80	Depositor
% Data completeness (in resolution range)	95.4 (20.00-2.80)	Depositor
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS 0.4	Depositor
R, R_{free}	0.229 , 0.288	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	9256	wwPDB-VP
Average B, all atoms (Å ²)	53.0	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: GAL, MAN, BMA, NAG, SIA

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	0.51	0/2298	0.98	7/3122 (0.2%)
1	B	0.48	0/2305	0.93	6/3130 (0.2%)
1	C	0.48	0/2297	0.98	12/3119 (0.4%)
1	D	0.46	0/2224	0.92	4/3015 (0.1%)
All	All	0.48	0/9124	0.95	29/12386 (0.2%)

There are no bond length outliers.

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	255	THR	N-CA-C	-9.67	101.48	113.28
1	C	151	GLU	CA-C-N	7.83	129.63	119.84
1	C	151	GLU	C-N-CA	7.83	129.63	119.84
1	A	177	SER	N-CA-C	6.99	122.92	113.97
1	A	188	SER	N-CA-C	-6.82	98.57	109.76
1	B	7	ARG	N-CA-C	6.79	119.06	110.24
1	C	186	PRO	CA-C-N	6.67	127.67	120.13
1	C	186	PRO	C-N-CA	6.67	127.67	120.13
1	B	237	ASN	N-CA-C	6.62	118.18	110.97
1	A	22	GLU	N-CA-C	6.53	118.95	111.11
1	C	188	SER	N-CA-C	-6.35	99.29	109.96
1	A	153	VAL	N-CA-C	6.26	116.93	110.36
1	B	188	SER	N-CA-C	-6.07	99.81	109.76
1	D	188	SER	N-CA-C	-5.93	100.83	110.20
1	C	254	ASP	N-CA-C	5.75	118.81	110.42
1	A	33	LEU	N-CA-C	-5.68	99.16	108.41
1	A	237	ASN	N-CA-C	5.63	117.10	110.97
1	B	33	LEU	N-CA-C	-5.39	99.62	108.41
1	D	130	GLU	N-CA-C	5.38	119.60	113.19
1	C	34	ASN	CB-CA-C	-5.37	110.40	116.63

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	107	ASN	N-CA-C	-5.36	104.64	110.91
1	C	208	TYR	N-CA-C	5.34	116.91	108.96
1	B	153	VAL	CB-CA-C	-5.33	105.16	111.97
1	A	208	TYR	N-CA-C	5.24	116.76	108.96
1	D	34	ASN	CB-CA-C	-5.22	110.58	116.63
1	D	33	LEU	N-CA-C	-5.21	99.93	108.41
1	C	229	GLU	N-CA-C	-5.18	106.11	112.90
1	C	33	LEU	N-CA-C	-5.16	100.01	108.41
1	B	229	GLU	N-CA-C	-5.11	106.20	112.90

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	2232	0	2124	96	0
1	B	2239	0	2129	97	0
1	C	2231	0	2124	116	0
1	D	2162	0	2060	98	0
2	E	89	0	76	4	0
3	F	120	0	101	3	0
4	G	28	0	25	0	0
5	H	39	0	34	0	0
6	A	28	0	26	1	0
6	B	28	0	26	1	0
6	C	28	0	26	0	0
6	D	28	0	26	1	0
7	A	1	0	0	0	0
7	B	1	0	0	0	0
7	C	1	0	0	0	0
7	D	1	0	0	0	0
All	All	9256	0	8777	397	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 22.

All (397) 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:C:37:GLN:HE21	1:C:51:MET:HE1	1.03	1.15
1:C:52:GLY:HA2	1:C:238:GLY:HA2	1.20	1.15
1:D:52:GLY:HA2	1:D:238:GLY:HA2	1.36	1.01
1:D:37:GLN:NE2	1:D:51:MET:HE1	1.81	0.95
1:C:37:GLN:NE2	1:C:51:MET:HE1	1.79	0.95
1:B:82:LYS:O	1:B:86:GLU:HG2	1.69	0.91
1:A:60:MET:HE1	1:A:177:SER:OG	1.69	0.91
1:B:54:TRP:HE1	1:B:177:SER:HB3	1.40	0.86
1:D:92:ASN:ND2	6:D:300:NAG:H83	1.89	0.86
1:C:19:LYS:HB2	1:C:94:SER:HB3	1.58	0.85
1:D:37:GLN:HE21	1:D:51:MET:HE1	1.41	0.85
1:C:8:TYR:HD1	1:C:105:ILE:HD11	1.42	0.84
1:B:53:LEU:HD12	1:B:180:ILE:HG23	1.60	0.83
1:B:7:ARG:HH11	1:B:7:ARG:HB2	1.45	0.81
1:B:153:VAL:HG22	1:D:235:LEU:HB3	1.62	0.80
1:B:157:ARG:NH2	1:D:234:VAL:HG22	1.98	0.79
1:B:269:GLN:HE21	1:B:269:GLN:HA	1.49	0.78
1:C:200:LYS:HB3	1:C:251:PRO:HB3	1.65	0.77
1:C:37:GLN:HG2	1:C:51:MET:CE	2.15	0.77
1:D:200:LYS:HD2	1:D:249:ALA:HB1	1.66	0.76
1:B:45:ASP:O	1:B:47:LYS:HG2	1.85	0.76
1:C:60:MET:HE1	1:C:177:SER:OG	1.85	0.76
1:A:180:ILE:O	1:A:183:ARG:HD3	1.86	0.76
1:D:60:MET:HE3	1:D:176:TYR:HB2	1.67	0.75
1:C:39:PHE:C	1:C:51:MET:HE2	2.13	0.74
1:D:131:ILE:HD11	1:D:133:ALA:HB3	1.69	0.74
1:D:275:TRP:O	1:D:276:GLU:HB2	1.85	0.74
1:A:60:MET:HE1	1:A:177:SER:HG	1.53	0.73
1:B:180:ILE:O	1:B:183:ARG:HG2	1.88	0.73
1:A:60:MET:HE3	1:A:176:TYR:HB2	1.69	0.73
1:D:39:PHE:C	1:D:51:MET:HE2	2.14	0.73
1:A:189:VAL:HG23	1:A:273:VAL:HG21	1.71	0.73
1:B:37:GLN:NE2	1:B:51:MET:HE1	2.03	0.72
1:C:60:MET:HE3	1:C:176:TYR:HB2	1.72	0.71
1:B:60:MET:HE3	1:B:176:TYR:HB2	1.73	0.71
1:C:183:ARG:NH2	1:C:240:GLY:O	2.24	0.70
1:D:131:ILE:HG13	1:D:132:PRO:C	2.14	0.70
1:D:167:PRO:O	1:D:171:ARG:HG2	1.90	0.70
1:D:189:VAL:HG23	1:D:273:VAL:HG21	1.72	0.70
1:B:39:PHE:C	1:B:51:MET:HE2	2.17	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:46:ARG:NE	1:C:71:LYS:HZ3	1.89	0.69
1:C:268:ALA:HB2	1:D:175:LYS:O	1.92	0.69
1:C:189:VAL:HG23	1:C:273:VAL:HG21	1.74	0.69
1:C:195:GLN:HB2	1:C:201:LYS:HA	1.73	0.69
1:D:129:LYS:HD2	1:D:163:GLU:HG2	1.75	0.69
1:B:184:GLN:NE2	1:B:265:SER:OG	2.26	0.68
1:B:189:VAL:HG23	1:B:273:VAL:HG21	1.75	0.68
1:C:167:PRO:O	1:C:171:ARG:HG2	1.94	0.68
1:C:37:GLN:HG2	1:C:51:MET:HE3	1.74	0.68
1:A:52:GLY:HA2	1:A:238:GLY:HA2	1.76	0.67
1:A:167:PRO:O	1:A:171:ARG:HG2	1.94	0.67
1:C:200:LYS:HD3	1:C:251:PRO:HG3	1.76	0.67
2:E:2:NAG:O3	2:E:3:BMA:H2	1.94	0.67
1:B:67:SER:O	1:B:71:LYS:HG3	1.95	0.67
1:D:151:GLU:HB3	1:D:153:VAL:HG23	1.76	0.67
1:A:157:ARG:NH2	1:C:234:VAL:HG22	2.10	0.67
1:D:216:ASP:HB3	1:D:263:GLN:HB3	1.77	0.66
1:B:269:GLN:HA	1:B:269:GLN:NE2	2.10	0.66
1:A:39:PHE:C	1:A:51:MET:HE2	2.21	0.66
1:C:216:ASP:HB3	1:C:263:GLN:HB3	1.78	0.66
1:B:184:GLN:HE21	1:B:266:SER:N	1.94	0.65
1:C:52:GLY:CA	1:C:238:GLY:HA2	2.13	0.65
1:C:105:ILE:C	1:C:105:ILE:HD12	2.21	0.65
1:D:53:LEU:HD12	1:D:180:ILE:HG23	1.79	0.65
1:B:167:PRO:O	1:B:171:ARG:HG2	1.97	0.65
1:C:46:ARG:HE	1:C:71:LYS:HZ3	1.45	0.64
1:C:250:VAL:HG11	1:C:258:TYR:CE1	2.31	0.64
1:A:153:VAL:HG21	1:C:235:LEU:O	1.96	0.64
1:B:157:ARG:HH22	1:D:234:VAL:HG22	1.62	0.64
1:C:191:VAL:HG11	1:C:275:TRP:HB2	1.79	0.64
1:D:106:GLU:HB2	1:D:111:SER:OG	1.98	0.64
1:D:174:LEU:O	1:D:178:LYS:HG2	1.98	0.64
1:B:52:GLY:HA2	1:B:238:GLY:HA2	1.79	0.64
1:B:178:LYS:HD3	1:B:182:ASP:OD2	1.98	0.63
1:C:196:ALA:HB3	1:C:199:GLU:OE1	1.99	0.63
1:C:199:GLU:HG2	1:C:200:LYS:H	1.62	0.63
1:D:136:PRO:HG3	1:D:142:GLN:NE2	2.14	0.62
1:C:52:GLY:HA2	1:C:238:GLY:CA	2.14	0.62
1:D:54:TRP:HA	1:D:57:VAL:HG23	1.82	0.62
1:D:152:PRO:O	1:D:156:GLN:HG2	1.99	0.62
1:B:54:TRP:NE1	1:B:177:SER:HB3	2.11	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:269:GLN:HE21	1:B:270:PRO:HD2	1.64	0.62
1:B:68:GLN:HE22	1:B:71:LYS:NZ	1.98	0.61
1:D:7:ARG:HB3	1:D:106:GLU:HG3	1.81	0.61
1:C:199:GLU:CG	1:C:200:LYS:H	2.13	0.61
1:B:184:GLN:HG3	1:B:266:SER:HB3	1.81	0.61
1:D:21:VAL:HG12	1:D:22:GLU:N	2.16	0.61
1:B:30:LEU:HD12	1:B:30:LEU:O	2.01	0.61
1:A:216:ASP:HB3	1:A:263:GLN:HB2	1.83	0.60
1:D:164:GLU:O	1:D:167:PRO:HG2	2.01	0.60
1:A:60:MET:HE2	1:A:173:TYR:HA	1.83	0.60
1:D:21:VAL:HG12	1:D:22:GLU:H	1.67	0.60
1:B:184:GLN:HG3	1:B:266:SER:CB	2.31	0.60
1:B:30:LEU:HB2	1:B:37:GLN:HE21	1.67	0.59
1:A:157:ARG:HH22	1:C:234:VAL:HG22	1.67	0.59
1:B:189:VAL:HG11	1:B:262:VAL:HG21	1.83	0.59
1:A:60:MET:HE2	1:A:173:TYR:CA	2.33	0.59
1:A:88:TYR:OH	1:A:139:PRO:HG2	2.02	0.59
1:A:105:ILE:HG22	1:A:110:SER:HA	1.83	0.59
1:C:8:TYR:CD1	1:C:105:ILE:HD11	2.33	0.58
2:E:2:NAG:H4	2:E:3:BMA:O2	2.03	0.58
1:B:30:LEU:HB2	1:B:37:GLN:NE2	2.19	0.58
1:C:19:LYS:HB2	1:C:94:SER:CB	2.31	0.58
1:C:60:MET:HE2	1:C:173:TYR:CA	2.34	0.58
1:D:275:TRP:O	1:D:276:GLU:CB	2.52	0.58
1:D:105:ILE:HG22	1:D:110:SER:HA	1.85	0.58
1:C:183:ARG:HH11	1:C:183:ARG:HG2	1.69	0.58
1:C:180:ILE:O	1:C:183:ARG:NH1	2.37	0.57
1:B:60:MET:HE2	1:B:173:TYR:HA	1.86	0.57
1:C:20:HIS:CD2	1:C:26:ALA:HB2	2.39	0.57
1:C:199:GLU:HG2	1:C:200:LYS:N	2.19	0.57
1:C:60:MET:HE2	1:C:173:TYR:HA	1.86	0.57
1:D:228:PRO:HB3	1:D:246:VAL:HG13	1.86	0.57
1:A:201:LYS:HB3	1:A:250:VAL:HG13	1.87	0.57
1:C:200:LYS:CB	1:C:251:PRO:HB3	2.34	0.57
1:D:88:TYR:OH	1:D:138:ASP:OD1	2.23	0.57
1:A:189:VAL:HG11	1:A:262:VAL:HG21	1.86	0.56
1:A:30:LEU:O	1:A:30:LEU:HD12	2.05	0.56
1:C:83:ASP:O	1:C:86:GLU:HB3	2.05	0.56
1:A:60:MET:CE	1:A:177:SER:OG	2.48	0.56
1:B:60:MET:HE2	1:B:173:TYR:CA	2.34	0.56
1:C:37:GLN:HG2	1:C:51:MET:HE1	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:142:GLN:CD	1:C:17:LEU:HD13	2.30	0.56
1:C:251:PRO:N	1:C:252:PRO:CD	2.69	0.56
1:C:197:PRO:C	1:C:199:GLU:H	2.13	0.56
1:C:186:PRO:HA	1:C:266:SER:OG	2.06	0.56
1:C:228:PRO:HB3	1:C:246:VAL:HG13	1.87	0.56
1:B:7:ARG:HH11	1:B:7:ARG:CB	2.17	0.55
1:B:269:GLN:HE21	1:B:269:GLN:CA	2.14	0.55
1:C:46:ARG:CD	1:C:71:LYS:HZ3	2.18	0.55
1:D:30:LEU:HD12	1:D:30:LEU:O	2.05	0.55
1:A:54:TRP:HA	1:A:57:VAL:HG23	1.87	0.55
1:A:150:ALA:HB3	1:A:154:TYR:HE2	1.72	0.55
1:A:228:PRO:HB3	1:A:246:VAL:HG13	1.89	0.55
1:B:228:PRO:HB3	1:B:246:VAL:HG13	1.89	0.55
1:D:228:PRO:HB3	1:D:246:VAL:CG1	2.36	0.55
1:A:189:VAL:CG2	1:A:273:VAL:HG21	2.34	0.55
1:B:184:GLN:HE22	1:B:265:SER:CB	2.19	0.55
1:D:106:GLU:N	1:D:111:SER:OG	2.39	0.55
1:D:189:VAL:CG2	1:D:273:VAL:HG21	2.36	0.55
1:D:208:TYR:CB	1:D:243:GLN:HG2	2.37	0.55
1:C:189:VAL:CG2	1:C:273:VAL:HG21	2.37	0.54
1:A:100:ARG:HD2	1:A:118:TYR:CE1	2.43	0.54
1:B:208:TYR:CB	1:B:243:GLN:HG2	2.38	0.54
1:C:41:TYR:CD1	1:C:41:TYR:C	2.85	0.54
1:A:20:HIS:CD2	1:A:26:ALA:HB2	2.41	0.54
1:B:30:LEU:HD12	1:B:30:LEU:C	2.33	0.54
1:D:108:ASN:OD1	1:D:171:ARG:NH1	2.40	0.54
1:B:239:ASN:OD1	1:B:241:THR:HG23	2.08	0.54
1:C:164:GLU:O	1:C:167:PRO:HG2	2.07	0.54
1:C:189:VAL:HG11	1:C:262:VAL:HG21	1.90	0.54
1:A:41:TYR:CD1	1:A:41:TYR:C	2.86	0.54
1:A:192:THR:CG2	1:A:193:SER:N	2.71	0.54
1:A:208:TYR:CB	1:A:243:GLN:HG2	2.37	0.53
1:B:189:VAL:CG2	1:B:273:VAL:HG21	2.38	0.53
1:C:208:TYR:CB	1:C:243:GLN:HG2	2.37	0.53
1:D:60:MET:HE2	1:D:173:TYR:HA	1.88	0.53
1:A:139:PRO:O	1:A:142:GLN:HB2	2.09	0.53
1:A:228:PRO:HB3	1:A:246:VAL:CG1	2.39	0.53
1:B:216:ASP:HB3	1:B:263:GLN:HB2	1.91	0.53
1:B:88:TYR:OH	1:B:138:ASP:OD1	2.26	0.52
1:D:60:MET:HE2	1:D:173:TYR:CA	2.38	0.52
1:B:7:ARG:HB2	1:B:7:ARG:NH1	2.19	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:41:TYR:CD1	1:D:41:TYR:C	2.87	0.52
1:B:201:LYS:HB3	1:B:250:VAL:HG22	1.90	0.52
1:B:228:PRO:HB3	1:B:246:VAL:CG1	2.39	0.52
1:C:229:GLU:OE1	1:C:249:ALA:HB2	2.10	0.52
1:B:192:THR:CG2	1:B:193:SER:N	2.73	0.52
1:D:27:PHE:HB3	1:D:43:SER:HB3	1.91	0.52
1:A:208:TYR:HB3	1:A:243:GLN:HG2	1.92	0.52
1:A:221:ARG:HG3	1:A:258:TYR:CZ	2.45	0.52
1:B:41:TYR:CD1	1:B:41:TYR:C	2.87	0.52
1:C:199:GLU:CG	1:C:200:LYS:N	2.73	0.52
1:D:191:VAL:HG11	1:D:275:TRP:HB2	1.92	0.52
1:B:20:HIS:CD2	1:B:26:ALA:HB2	2.45	0.51
1:D:189:VAL:HG11	1:D:262:VAL:HG21	1.92	0.51
1:A:234:VAL:HG12	1:A:235:LEU:N	2.25	0.51
1:C:252:PRO:O	1:C:254:ASP:N	2.41	0.51
1:A:132:PRO:O	1:A:133:ALA:HB2	2.10	0.51
1:A:255:THR:O	1:A:255:THR:HG22	2.10	0.51
1:D:200:LYS:HG2	1:D:201:LYS:N	2.25	0.51
1:A:85:VAL:CG1	1:A:91:SER:HA	2.41	0.51
1:C:251:PRO:N	1:C:252:PRO:HD2	2.26	0.51
1:A:27:PHE:HB3	1:A:43:SER:HB3	1.93	0.51
1:A:30:LEU:HD12	1:A:30:LEU:C	2.36	0.51
1:C:228:PRO:HB3	1:C:246:VAL:CG1	2.41	0.51
1:B:234:VAL:HG12	1:B:235:LEU:N	2.25	0.50
1:D:20:HIS:CD2	1:D:26:ALA:HB2	2.47	0.50
1:D:131:ILE:HG13	1:D:132:PRO:CA	2.40	0.50
1:A:53:LEU:HD12	1:A:180:ILE:HG23	1.93	0.50
1:B:27:PHE:HB3	1:B:43:SER:HB3	1.92	0.50
1:D:196:ALA:HB1	1:D:197:PRO:HD2	1.93	0.50
1:C:30:LEU:HD12	1:C:30:LEU:O	2.11	0.50
1:A:164:GLU:O	1:A:167:PRO:HG2	2.11	0.50
1:C:234:VAL:HG12	1:C:235:LEU:N	2.25	0.50
1:A:100:ARG:HD2	1:A:118:TYR:HE1	1.76	0.49
1:B:229:GLU:OE1	1:B:249:ALA:HB2	2.12	0.49
1:D:30:LEU:HD12	1:D:30:LEU:C	2.37	0.49
1:B:208:TYR:HB3	1:B:243:GLN:HG2	1.93	0.49
1:D:131:ILE:H	1:D:132:PRO:HA	1.76	0.49
1:A:21:VAL:HG22	1:A:24:VAL:HG23	1.94	0.49
1:A:221:ARG:HE	1:A:256:ALA:CB	2.26	0.49
1:D:186:PRO:HA	1:D:266:SER:OG	2.11	0.49
1:C:18:SER:OG	1:C:95:HIS:N	2.45	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:27:PHE:HB3	1:C:43:SER:HB3	1.93	0.49
1:D:208:TYR:HB3	1:D:243:GLN:HG2	1.94	0.49
1:C:197:PRO:C	1:C:199:GLU:N	2.70	0.49
1:D:60:MET:HE1	1:D:177:SER:OG	2.13	0.49
1:A:87:TYR:O	6:A:300:NAG:H82	2.12	0.49
1:A:239:ASN:OD1	1:A:241:THR:HG23	2.12	0.49
1:A:130:GLU:HA	1:A:130:GLU:OE1	2.12	0.48
1:C:208:TYR:HB3	1:C:243:GLN:HG2	1.94	0.48
1:C:127:PHE:CE2	1:C:129:LYS:HA	2.48	0.48
1:D:131:ILE:CD1	1:D:133:ALA:HB3	2.40	0.48
3:F:2:NAG:H83	3:F:5:NAG:O3	2.13	0.48
1:D:85:VAL:CG1	1:D:91:SER:HA	2.43	0.48
1:B:164:GLU:O	1:B:167:PRO:HG2	2.13	0.48
1:A:43:SER:O	1:A:46:ARG:HD2	2.14	0.48
1:A:192:THR:HG23	1:A:193:SER:N	2.29	0.48
1:D:234:VAL:HG12	1:D:235:LEU:N	2.27	0.48
1:A:153:VAL:O	1:A:156:GLN:HB2	2.14	0.48
1:A:208:TYR:CE2	2:E:2:NAG:H61	2.48	0.48
1:C:30:LEU:HD12	1:C:30:LEU:C	2.38	0.48
1:D:45:ASP:O	1:D:46:ARG:CG	2.62	0.48
1:C:46:ARG:HG2	1:C:71:LYS:HZ3	1.79	0.48
1:D:89:ASN:CG	1:D:89:ASN:O	2.56	0.48
3:F:1:NAG:O3	3:F:2:NAG:O5	2.32	0.48
1:A:136:PRO:HG2	1:C:96:VAL:HG21	1.96	0.47
1:C:250:VAL:C	1:C:252:PRO:HD2	2.39	0.47
1:B:143:ILE:O	1:B:146:GLN:HB2	2.13	0.47
1:D:190:VAL:HG23	1:D:208:TYR:HE1	1.79	0.47
1:A:201:LYS:HZ3	1:A:275:TRP:HH2	1.63	0.47
1:B:205:CYS:O	1:B:245:TRP:HA	2.14	0.47
1:C:221:ARG:HD2	1:C:258:TYR:OH	2.14	0.47
1:A:157:ARG:HH21	1:C:234:VAL:HA	1.78	0.47
1:B:177:SER:O	1:B:178:LYS:C	2.56	0.47
1:D:45:ASP:O	1:D:47:LYS:HG2	2.15	0.47
1:C:174:LEU:HD23	1:C:174:LEU:HA	1.72	0.47
1:D:139:PRO:O	1:D:142:GLN:HB2	2.15	0.47
1:A:19:LYS:NZ	1:A:92:ASN:ND2	2.62	0.47
1:A:54:TRP:HA	1:A:57:VAL:CG2	2.44	0.47
1:B:184:GLN:HE21	1:B:265:SER:C	2.23	0.47
1:B:200:LYS:HB3	1:B:250:VAL:O	2.14	0.47
1:A:205:CYS:O	1:A:245:TRP:HA	2.15	0.47
1:B:189:VAL:CG1	1:B:262:VAL:HG21	2.44	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:53:LEU:HD21	1:C:239:ASN:O	2.15	0.46
1:A:85:VAL:HG12	1:A:91:SER:HA	1.97	0.46
1:C:116:LYS:HG3	1:C:126:GLU:HB2	1.98	0.46
1:D:127:PHE:HB2	1:D:134:TRP:CZ3	2.50	0.46
2:E:1:NAG:O3	2:E:2:NAG:O5	2.32	0.46
1:C:153:VAL:O	1:C:156:GLN:N	2.49	0.46
1:C:190:VAL:HG23	1:C:208:TYR:HE1	1.81	0.46
1:C:180:ILE:O	1:C:183:ARG:HG2	2.16	0.46
1:D:174:LEU:HD23	1:D:174:LEU:HA	1.70	0.46
1:A:25:PRO:HG2	1:A:28:GLN:CD	2.40	0.46
1:A:127:PHE:HB2	1:A:134:TRP:CZ3	2.51	0.46
1:A:145:LYS:O	1:A:149:GLU:HG3	2.16	0.46
1:C:46:ARG:CG	1:C:71:LYS:HZ3	2.28	0.46
1:A:41:TYR:CE2	1:A:70:GLN:HB2	2.51	0.46
1:A:79:GLU:O	1:A:82:LYS:HB3	2.15	0.46
1:B:128:ASN:O	1:B:132:PRO:HA	2.16	0.46
1:A:7:ARG:HG2	1:A:106:GLU:HG3	1.98	0.46
1:B:81:LEU:HD13	1:B:97:LEU:HB2	1.97	0.46
1:B:184:GLN:NE2	1:B:265:SER:CB	2.78	0.46
1:B:190:VAL:HG23	1:B:208:TYR:HE1	1.81	0.46
1:C:106:GLU:O	1:C:107:ASN:HB3	2.15	0.46
1:B:106:GLU:O	1:B:107:ASN:HB2	2.16	0.46
1:B:157:ARG:O	1:B:160:ALA:HB3	2.16	0.46
1:C:41:TYR:CE2	1:C:70:GLN:HB2	2.51	0.45
1:C:46:ARG:HG2	1:C:71:LYS:NZ	2.30	0.45
1:C:145:LYS:O	1:C:149:GLU:HG3	2.15	0.45
1:D:153:VAL:O	1:D:156:GLN:HB2	2.16	0.45
1:D:177:SER:HB3	1:D:181:LEU:HG	1.98	0.45
1:A:186:PRO:HA	1:A:266:SER:OG	2.17	0.45
1:B:142:GLN:OE1	1:D:17:LEU:HD13	2.16	0.45
1:C:159:LYS:O	1:C:163:GLU:HG3	2.17	0.45
1:A:153:VAL:HG22	1:C:235:LEU:HB3	1.97	0.45
1:A:221:ARG:HE	1:A:256:ALA:HB3	1.81	0.45
1:B:41:TYR:CE2	1:B:70:GLN:HB2	2.51	0.45
1:B:88:TYR:HB3	1:B:90:ASP:OD2	2.17	0.45
1:D:159:LYS:O	1:D:163:GLU:HG3	2.17	0.45
1:B:88:TYR:HA	6:B:300:NAG:H82	1.99	0.45
1:A:86:GLU:HA	1:A:86:GLU:OE1	2.16	0.45
1:A:159:LYS:O	1:A:163:GLU:HG3	2.17	0.45
1:B:174:LEU:HD23	1:B:174:LEU:HA	1.72	0.45
1:C:89:ASN:CG	1:C:89:ASN:O	2.60	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:205:CYS:O	1:C:245:TRP:HA	2.17	0.44
1:D:41:TYR:CE2	1:D:70:GLN:HB2	2.51	0.44
1:D:166:CYS:HB3	1:D:167:PRO:CD	2.47	0.44
1:A:178:LYS:HD3	1:A:182:ASP:OD2	2.17	0.44
1:A:188:SER:HB2	1:A:208:TYR:CZ	2.52	0.44
1:B:275:TRP:CZ3	1:B:277:ALA:HB2	2.52	0.44
1:B:68:GLN:HE22	1:B:71:LYS:HZ1	1.64	0.44
1:B:166:CYS:HB3	1:B:167:PRO:CD	2.47	0.44
1:C:42:ASN:OD1	1:C:44:LYS:HB3	2.17	0.44
1:D:11:THR:O	1:D:31:GLY:HA2	2.18	0.44
1:C:252:PRO:HG2	1:C:254:ASP:OD2	2.17	0.44
1:A:189:VAL:CG1	1:A:262:VAL:HG21	2.47	0.44
1:B:136:PRO:HG2	1:D:96:VAL:HG21	1.98	0.44
1:B:189:VAL:HA	1:B:206:LEU:O	2.18	0.44
1:D:42:ASN:OD1	1:D:44:LYS:HB3	2.18	0.44
1:D:166:CYS:N	1:D:167:PRO:HD2	2.33	0.44
1:C:105:ILE:HD12	1:C:105:ILE:O	2.18	0.43
1:C:143:ILE:O	1:C:146:GLN:HB3	2.18	0.43
1:D:54:TRP:HA	1:D:57:VAL:CG2	2.47	0.43
1:D:116:LYS:HG3	1:D:126:GLU:HB2	1.99	0.43
1:A:201:LYS:NZ	1:A:275:TRP:HH2	2.16	0.43
1:B:192:THR:HG23	1:B:193:SER:N	2.33	0.43
1:D:205:CYS:O	1:D:245:TRP:HA	2.18	0.43
1:B:127:PHE:HB2	1:B:134:TRP:CZ3	2.53	0.43
1:A:190:VAL:HG23	1:A:208:TYR:HE1	1.82	0.43
1:B:81:LEU:O	1:B:85:VAL:HG23	2.18	0.43
1:B:186:PRO:HA	1:B:266:SER:OG	2.18	0.43
1:D:178:LYS:O	1:D:182:ASP:HB2	2.18	0.43
1:B:142:GLN:CD	1:D:17:LEU:HD13	2.44	0.43
1:C:19:LYS:CB	1:C:94:SER:HB3	2.41	0.43
1:C:151:GLU:HA	1:C:151:GLU:OE1	2.19	0.43
1:B:79:GLU:O	1:B:82:LYS:HB3	2.19	0.43
1:C:107:ASN:O	1:C:108:ASN:C	2.61	0.43
1:C:188:SER:HB2	1:C:208:TYR:CZ	2.53	0.43
1:C:222:ALA:HB2	1:C:256:ALA:HB1	2.00	0.43
1:A:166:CYS:HB3	1:A:167:PRO:CD	2.48	0.43
1:A:237:ASN:C	1:A:239:ASN:N	2.75	0.43
1:B:187:PRO:HB3	1:B:210:PHE:HB3	1.99	0.43
1:C:81:LEU:HD13	1:C:97:LEU:HB2	2.00	0.43
1:B:255:THR:O	1:B:256:ALA:C	2.61	0.43
1:C:60:MET:HE1	1:C:177:SER:HG	1.82	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:86:GLU:OE1	1:A:86:GLU:CA	2.66	0.43
1:A:229:GLU:OE1	1:A:249:ALA:HB2	2.19	0.43
1:A:232:GLY:HA3	1:A:245:TRP:CE2	2.54	0.43
1:A:234:VAL:CG1	1:A:235:LEU:N	2.82	0.43
1:B:21:VAL:CG2	1:B:24:VAL:HG23	2.49	0.43
1:B:276:GLU:HA	1:B:276:GLU:OE1	2.19	0.43
1:D:81:LEU:HD13	1:D:97:LEU:HB2	2.00	0.42
1:A:19:LYS:NZ	1:A:92:ASN:CG	2.78	0.42
1:A:71:LYS:HE2	1:A:71:LYS:O	2.19	0.42
1:A:174:LEU:HD23	1:A:174:LEU:HA	1.74	0.42
1:C:127:PHE:HB2	1:C:134:TRP:CZ3	2.54	0.42
1:D:196:ALA:HB3	1:D:199:GLU:OE1	2.19	0.42
1:B:135:VAL:HG22	1:D:98:GLN:HE22	1.84	0.42
1:C:11:THR:O	1:C:31:GLY:HA2	2.18	0.42
1:C:130:GLU:O	1:C:131:ILE:C	2.62	0.42
1:C:189:VAL:CG1	1:C:262:VAL:HG21	2.49	0.42
1:A:19:LYS:HZ2	1:A:92:ASN:ND2	2.16	0.42
1:A:177:SER:O	1:A:178:LYS:C	2.61	0.42
1:D:125:ILE:HD11	1:D:134:TRP:HB3	2.01	0.42
1:B:234:VAL:CG1	1:B:235:LEU:N	2.82	0.42
1:C:53:LEU:HD11	1:C:240:GLY:HA3	2.00	0.42
1:C:234:VAL:CG1	1:C:235:LEU:N	2.82	0.42
1:D:45:ASP:O	1:D:46:ARG:HG2	2.20	0.42
1:D:106:GLU:O	1:D:107:ASN:CB	2.68	0.42
1:C:151:GLU:HA	1:C:152:PRO:HD3	1.73	0.42
1:D:145:LYS:O	1:D:149:GLU:HG3	2.20	0.42
1:A:88:TYR:OH	1:A:138:ASP:OD1	2.36	0.42
1:C:81:LEU:O	1:C:85:VAL:HG23	2.19	0.42
1:D:22:GLU:O	1:D:23:ASP:HB2	2.19	0.42
1:B:145:LYS:O	1:B:149:GLU:HG3	2.20	0.41
1:D:54:TRP:NE1	1:D:177:SER:OG	2.48	0.41
1:A:81:LEU:O	1:A:85:VAL:HG23	2.20	0.41
1:B:160:ALA:O	1:B:164:GLU:HB2	2.20	0.41
1:D:87:TYR:CD1	1:D:87:TYR:C	2.98	0.41
1:A:189:VAL:HA	1:A:206:LEU:O	2.19	0.41
1:B:208:TYR:CE2	3:F:2:NAG:H61	2.55	0.41
1:C:195:GLN:HG3	1:C:200:LYS:C	2.45	0.41
1:C:227:GLU:CD	1:C:228:PRO:HD2	2.45	0.41
1:C:189:VAL:HA	1:C:206:LEU:O	2.20	0.41
1:C:227:GLU:OE2	1:C:228:PRO:HD2	2.20	0.41
1:D:189:VAL:CG1	1:D:262:VAL:HG21	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:234:VAL:CG1	1:D:235:LEU:N	2.84	0.41
1:A:81:LEU:HD13	1:A:97:LEU:HB2	2.03	0.41
1:B:42:ASN:OD1	1:B:44:LYS:HB3	2.21	0.41
1:B:105:ILE:HG22	1:B:110:SER:HA	2.02	0.41
1:C:8:TYR:HA	1:C:34:ASN:OD1	2.21	0.41
1:B:195:GLN:NE2	1:B:199:GLU:O	2.48	0.41
1:D:189:VAL:HA	1:D:206:LEU:O	2.20	0.41
1:C:166:CYS:HB3	1:C:167:PRO:CD	2.51	0.41
1:D:131:ILE:HG13	1:D:133:ALA:N	2.34	0.41
1:A:252:PRO:HG2	1:A:253:GLN:H	1.86	0.41
1:B:11:THR:O	1:B:31:GLY:HA2	2.21	0.41
1:B:250:VAL:HA	1:B:251:PRO:HD3	1.91	0.41
1:C:21:VAL:HG12	1:C:22:GLU:N	2.36	0.41
1:A:42:ASN:OD1	1:A:44:LYS:HB3	2.21	0.40
1:C:178:LYS:HA	1:C:182:ASP:OD2	2.21	0.40
1:C:195:GLN:CB	1:C:201:LYS:HA	2.45	0.40
1:A:166:CYS:N	1:A:167:PRO:HD2	2.37	0.40
1:B:132:PRO:O	1:B:133:ALA:HB2	2.20	0.40
1:D:54:TRP:HE1	1:D:60:MET:HE1	1.86	0.40
1:D:106:GLU:O	1:D:107:ASN:HB2	2.20	0.40
1:A:183:ARG:HD3	1:A:183:ARG:HH11	1.76	0.40
1:C:60:MET:HE2	1:C:173:TYR:CB	2.51	0.40
1:A:203:LEU:O	1:A:247:VAL:HA	2.22	0.40
1:A:231:ARG:HG3	1:A:232:GLY:H	1.86	0.40
1:B:159:LYS:O	1:B:163:GLU:HG3	2.22	0.40
1:C:166:CYS:N	1:C:167:PRO:HD2	2.35	0.40
1:D:162:LEU:O	1:D:167:PRO:HD3	2.21	0.40
1:D:183:ARG:NH2	1:D:240:GLY:O	2.45	0.40
1:D:188:SER:HB2	1:D:208:TYR:CZ	2.57	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	271/274 (99%)	252 (93%)	18 (7%)	1 (0%)	30	60
1	B	272/274 (99%)	249 (92%)	23 (8%)	0	100	100
1	C	271/274 (99%)	249 (92%)	19 (7%)	3 (1%)	11	36
1	D	259/274 (94%)	243 (94%)	16 (6%)	0	100	100
All	All	1073/1096 (98%)	993 (92%)	76 (7%)	4 (0%)	30	60

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	91	SER
1	C	108	ASN
1	C	87	TYR
1	C	109	ARG

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	239/240 (100%)	229 (96%)	10 (4%)	26	61
1	B	240/240 (100%)	234 (98%)	6 (2%)	42	76
1	C	239/240 (100%)	232 (97%)	7 (3%)	37	73
1	D	231/240 (96%)	223 (96%)	8 (4%)	32	67
All	All	949/960 (99%)	918 (97%)	31 (3%)	33	69

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	ASP
1	A	21	VAL
1	A	28	GLN
1	A	39	PHE
1	A	41	TYR
1	A	71	LYS
1	A	86	GLU

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Mol	Chain	Res	Type
1	A	110	SER
1	A	192	THR
1	A	250	VAL
1	B	7	ARG
1	B	39	PHE
1	B	41	TYR
1	B	104	GLU
1	B	192	THR
1	B	250	VAL
1	C	22	GLU
1	C	23	ASP
1	C	39	PHE
1	C	192	THR
1	C	202	LYS
1	C	254	ASP
1	C	255	THR
1	D	30	LEU
1	D	39	PHE
1	D	41	TYR
1	D	89	ASN
1	D	153	VAL
1	D	184	GLN
1	D	192	THR
1	D	276	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	92	ASN
1	A	107	ASN
1	A	184	GLN
1	A	218	HIS
1	A	236	HIS
1	A	261	HIS
1	B	37	GLN
1	B	68	GLN
1	B	92	ASN
1	B	184	GLN
1	B	218	HIS
1	B	226	GLN
1	B	253	GLN
1	B	261	HIS

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Mol	Chain	Res	Type
1	B	269	GLN
1	C	37	GLN
1	C	49	GLN
1	C	218	HIS
1	C	261	HIS
1	C	269	GLN
1	D	37	GLN
1	D	92	ASN
1	D	98	GLN
1	D	142	GLN
1	D	156	GLN
1	D	184	GLN
1	D	218	HIS
1	D	261	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 ⓘ

21 monosaccharides are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	NAG	E	1	1,2	14,14,15	0.67	0	17,19,21	0.68	0
2	NAG	E	2	2	14,14,15	0.57	0	17,19,21	0.85	1 (5%)
2	BMA	E	3	2	11,11,12	0.57	0	15,15,17	1.33	2 (13%)
2	MAN	E	4	2	11,11,12	0.54	0	15,15,17	0.89	1 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NAG	E	5	2	14,14,15	0.63	0	17,19,21	0.94	0
2	MAN	E	6	2	11,11,12	0.61	0	15,15,17	0.64	0
2	NAG	E	7	2	14,14,15	0.77	1 (7%)	17,19,21	0.77	1 (5%)
3	NAG	F	1	1,3	14,14,15	0.56	0	17,19,21	0.65	0
3	NAG	F	2	3	14,14,15	0.60	0	17,19,21	0.70	0
3	BMA	F	3	3	11,11,12	0.37	0	15,15,17	1.74	2 (13%)
3	MAN	F	4	3	11,11,12	0.73	0	15,15,17	0.70	0
3	NAG	F	5	3	14,14,15	0.66	0	17,19,21	1.16	3 (17%)
3	GAL	F	6	3	11,11,12	0.70	0	15,15,17	0.50	0
3	SIA	F	7	3	20,20,21	0.93	0	21,28,31	0.82	1 (4%)
3	MAN	F	8	3	11,11,12	1.04	1 (9%)	15,15,17	1.37	2 (13%)
3	NAG	F	9	3	14,14,15	0.67	0	17,19,21	1.05	2 (11%)
4	NAG	G	1	1,4	14,14,15	0.59	0	17,19,21	0.71	0
4	NAG	G	2	4	14,14,15	0.56	0	17,19,21	0.89	1 (5%)
5	NAG	H	1	1,5	14,14,15	0.58	0	17,19,21	0.64	0
5	NAG	H	2	5	14,14,15	0.51	0	17,19,21	0.67	0
5	BMA	H	3	5	11,11,12	0.64	0	15,15,17	0.34	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
2	NAG	E	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	E	2	2	-	0/6/23/26	0/1/1/1
2	BMA	E	3	2	-	0/2/19/22	0/1/1/1
2	MAN	E	4	2	-	1/2/19/22	0/1/1/1
2	NAG	E	5	2	-	0/6/23/26	0/1/1/1
2	MAN	E	6	2	-	2/2/19/22	0/1/1/1
2	NAG	E	7	2	-	2/6/23/26	0/1/1/1
3	NAG	F	1	1,3	-	2/6/23/26	0/1/1/1
3	NAG	F	2	3	-	0/6/23/26	0/1/1/1
3	BMA	F	3	3	-	2/2/19/22	0/1/1/1
3	MAN	F	4	3	-	2/2/19/22	0/1/1/1
3	NAG	F	5	3	-	0/6/23/26	0/1/1/1
3	GAL	F	6	3	-	0/2/19/22	0/1/1/1
3	SIA	F	7	3	-	1/18/34/38	0/1/1/1
3	MAN	F	8	3	-	1/2/19/22	0/1/1/1
3	NAG	F	9	3	-	1/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	G	1	1,4	-	2/6/23/26	0/1/1/1
4	NAG	G	2	4	-	2/6/23/26	0/1/1/1
5	NAG	H	1	1,5	-	0/6/23/26	0/1/1/1
5	NAG	H	2	5	-	2/6/23/26	0/1/1/1
5	BMA	H	3	5	-	1/2/19/22	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	F	8	MAN	C2-C3	2.64	1.56	1.52
2	E	7	NAG	C1-C2	2.08	1.55	1.52

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	3	BMA	C6-C5-C4	-4.51	101.94	113.02
3	F	8	MAN	C1-C2-C3	3.60	114.88	109.64
2	E	3	BMA	C2-C3-C4	-3.49	104.72	110.86
3	F	3	BMA	C3-C4-C5	3.26	116.15	110.23
3	F	8	MAN	C2-C3-C4	3.08	116.28	110.86
2	E	3	BMA	C3-C4-C5	-2.88	105.00	110.23
3	F	5	NAG	C4-C3-C2	-2.72	107.03	111.02
2	E	4	MAN	C1-O5-C5	2.66	115.75	112.19
3	F	9	NAG	C1-O5-C5	2.65	115.74	112.19
4	G	2	NAG	C2-N2-C7	-2.54	119.50	122.90
3	F	9	NAG	C2-N2-C7	-2.49	119.57	122.90
3	F	7	SIA	O1B-C1-C2	2.28	118.64	112.71
3	F	5	NAG	C2-N2-C7	-2.25	119.88	122.90
3	F	5	NAG	C1-O5-C5	2.25	115.20	112.19
2	E	2	NAG	C2-N2-C7	-2.24	119.90	122.90
2	E	7	NAG	C2-N2-C7	-2.16	120.00	122.90

There are no chirality outliers.

All (23) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	E	1	NAG	O5-C5-C6-O6
2	E	7	NAG	O5-C5-C6-O6
3	F	1	NAG	O5-C5-C6-O6
3	F	4	MAN	O5-C5-C6-O6

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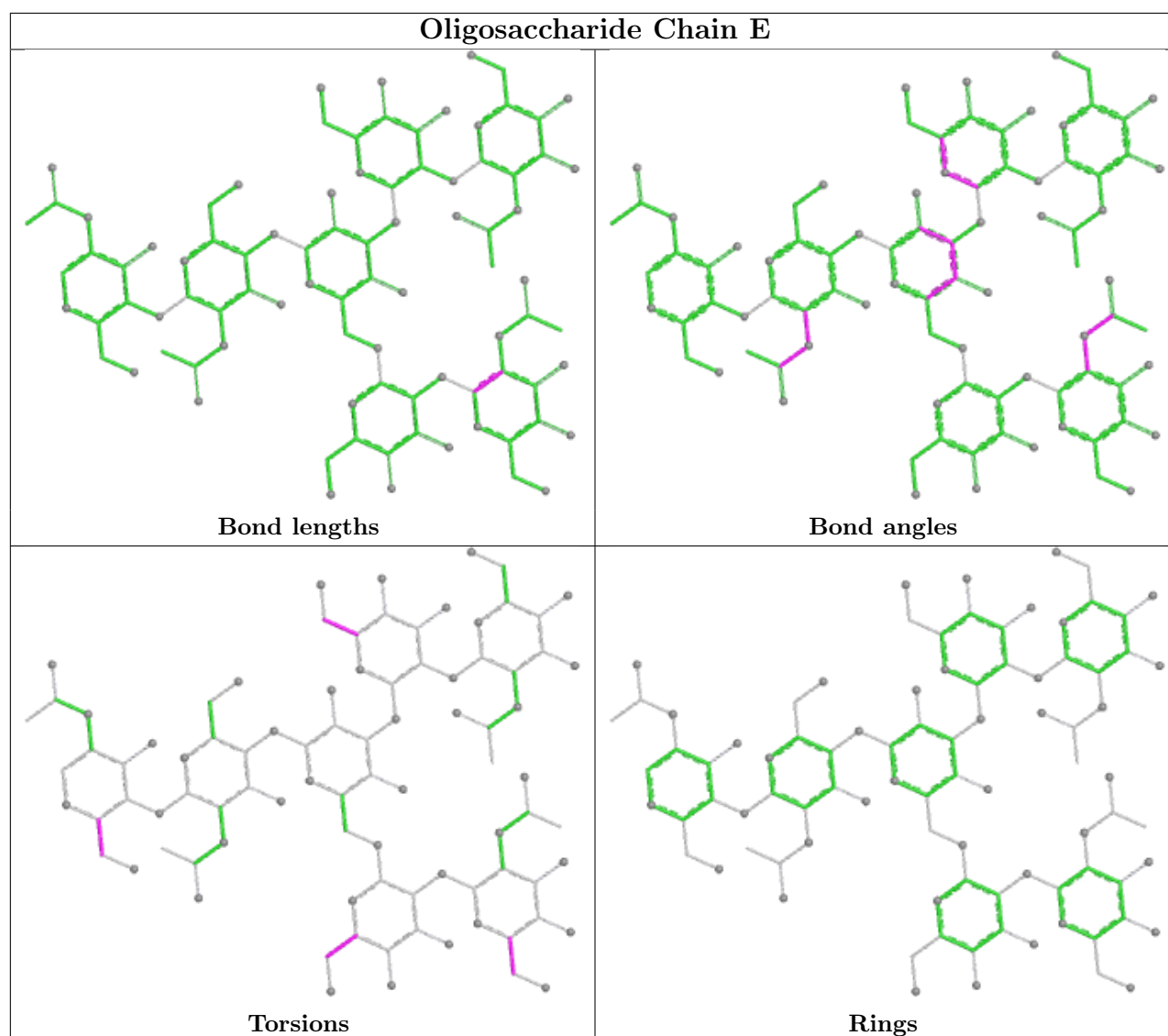
Mol	Chain	Res	Type	Atoms
5	H	2	NAG	O5-C5-C6-O6
3	F	1	NAG	C4-C5-C6-O6
2	E	7	NAG	C4-C5-C6-O6
5	H	2	NAG	C4-C5-C6-O6
3	F	4	MAN	C4-C5-C6-O6
2	E	6	MAN	O5-C5-C6-O6
2	E	1	NAG	C4-C5-C6-O6
4	G	1	NAG	O5-C5-C6-O6
5	H	3	BMA	O5-C5-C6-O6
4	G	1	NAG	C4-C5-C6-O6
2	E	6	MAN	C4-C5-C6-O6
4	G	2	NAG	C4-C5-C6-O6
4	G	2	NAG	O5-C5-C6-O6
3	F	9	NAG	O5-C5-C6-O6
3	F	8	MAN	O5-C5-C6-O6
3	F	3	BMA	C4-C5-C6-O6
3	F	3	BMA	O5-C5-C6-O6
3	F	7	SIA	C4-C5-N5-C10
2	E	4	MAN	O5-C5-C6-O6

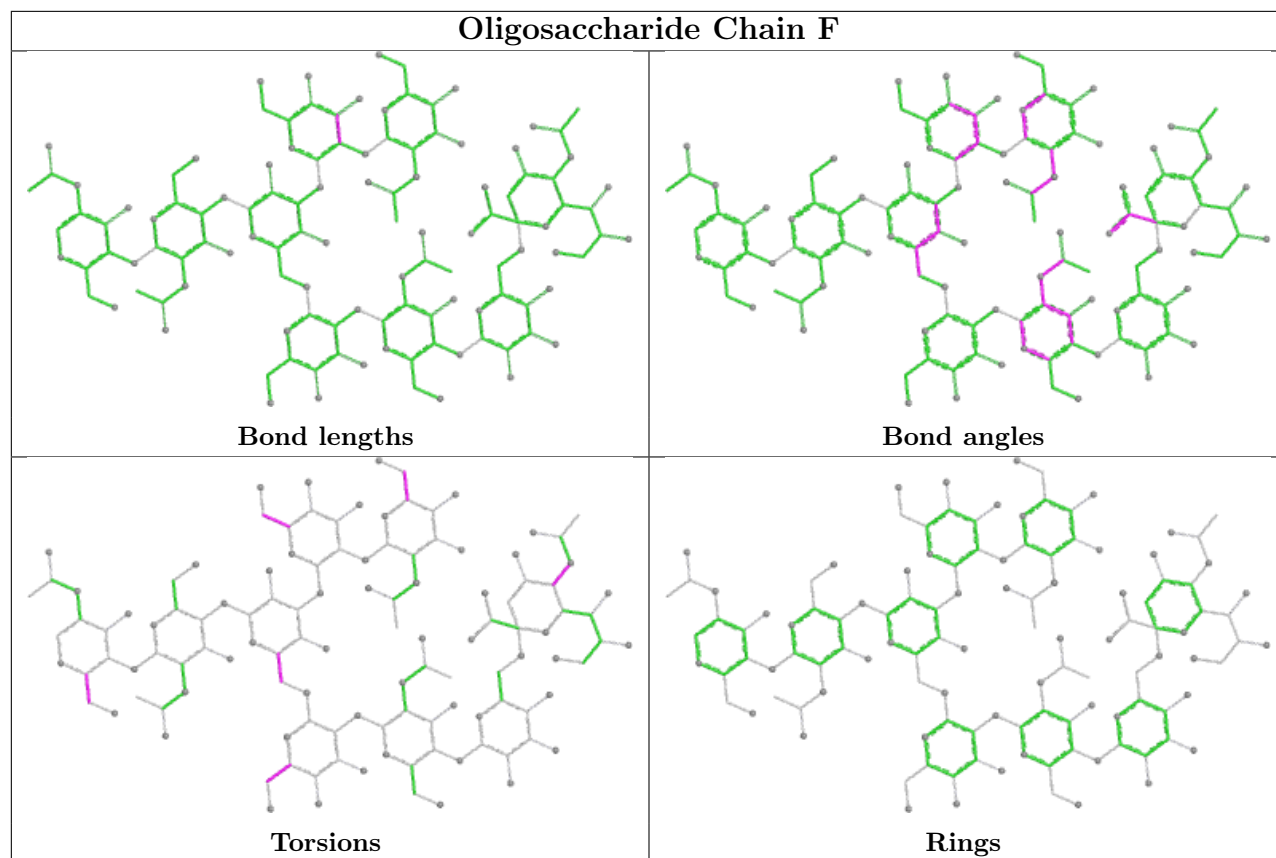
There are no ring outliers.

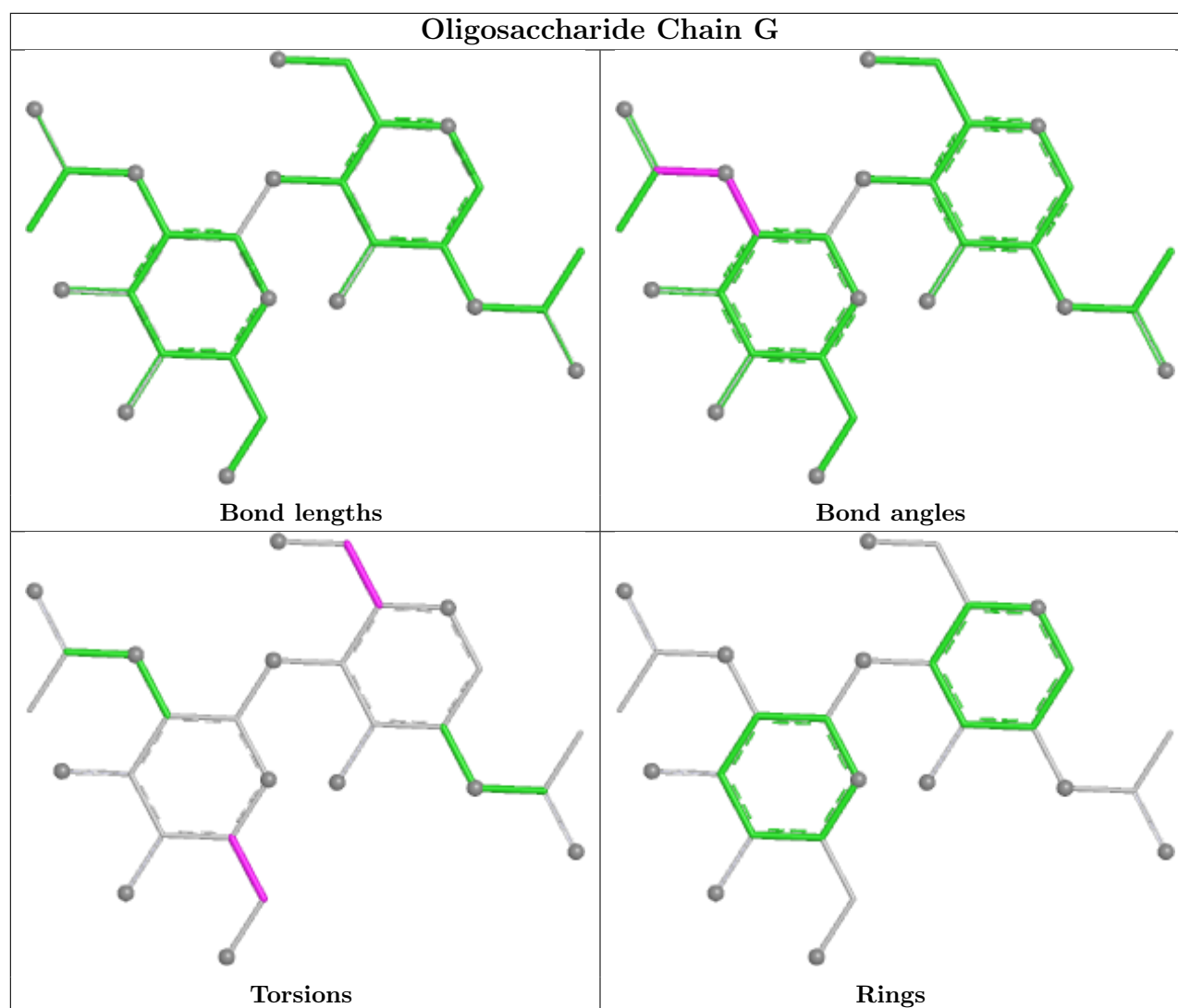
6 monomers are involved in 7 short contacts:

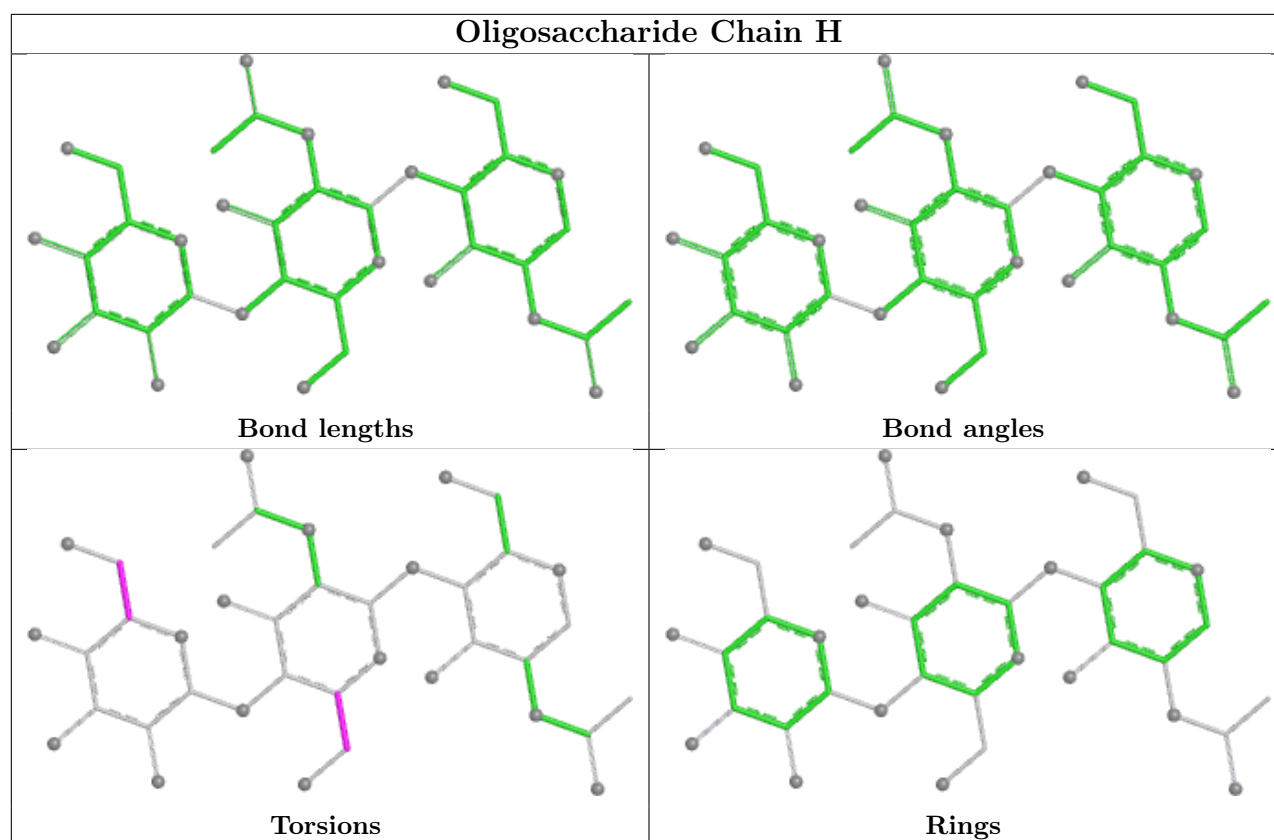
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	F	5	NAG	1	0
3	F	1	NAG	1	0
3	F	2	NAG	3	0
2	E	1	NAG	1	0
2	E	3	BMA	2	0
2	E	2	NAG	4	0

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









5.6 Ligand geometry [i](#)

8 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	$\# Z > 2$
6	NAG	D	310	1	14,14,15	0.67	0	17,19,21	0.75	0
6	NAG	C	300	1	14,14,15	0.72	1 (7%)	17,19,21	0.72	0
6	NAG	A	300	1	14,14,15	0.69	0	17,19,21	0.72	0
6	NAG	C	310	1	14,14,15	0.57	0	17,19,21	0.74	1 (5%)
6	NAG	D	300	1	14,14,15	0.55	0	17,19,21	0.76	1 (5%)
6	NAG	B	300	1	14,14,15	0.76	1 (7%)	17,19,21	0.62	0
6	NAG	B	310	1	14,14,15	0.58	0	17,19,21	0.64	1 (5%)
6	NAG	A	310	1	14,14,15	0.49	0	17,19,21	0.91	1 (5%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	NAG	D	310	1	-	2/6/23/26	0/1/1/1
6	NAG	C	300	1	-	2/6/23/26	0/1/1/1
6	NAG	A	300	1	-	0/6/23/26	0/1/1/1
6	NAG	C	310	1	1/1/5/7	4/6/23/26	0/1/1/1
6	NAG	D	300	1	-	2/6/23/26	0/1/1/1
6	NAG	B	300	1	-	1/6/23/26	0/1/1/1
6	NAG	B	310	1	-	1/6/23/26	0/1/1/1
6	NAG	A	310	1	-	2/6/23/26	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	C	300	NAG	C1-C2	2.14	1.55	1.52
6	B	300	NAG	C1-C2	2.10	1.55	1.52

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	310	NAG	C2-N2-C7	-2.40	119.69	122.90
6	D	300	NAG	C2-N2-C7	-2.37	119.72	122.90
6	B	310	NAG	C2-N2-C7	-2.13	120.05	122.90
6	C	310	NAG	C2-N2-C7	-2.02	120.19	122.90

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
6	C	310	NAG	C1

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	D	300	NAG	O5-C5-C6-O6
6	C	310	NAG	C4-C5-C6-O6
6	D	300	NAG	C4-C5-C6-O6
6	C	300	NAG	O5-C5-C6-O6
6	C	300	NAG	C4-C5-C6-O6
6	D	310	NAG	O5-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
6	C	310	NAG	O5-C5-C6-O6
6	D	310	NAG	C4-C5-C6-O6
6	B	300	NAG	O5-C5-C6-O6
6	B	310	NAG	O5-C5-C6-O6
6	C	310	NAG	C3-C2-N2-C7
6	C	310	NAG	C1-C2-N2-C7
6	A	310	NAG	C4-C5-C6-O6
6	A	310	NAG	O5-C5-C6-O6

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	300	NAG	1	0
6	D	300	NAG	1	0
6	B	300	NAG	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

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.