



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 11:19 AM UTC

PDB ID : 7Y4Q / pdb_00007y4q
Title : Semaphorin 6D in complex with Plexin A1
Authors : Tanaka, T.; Neyazaki, M.; Nogi, T.
Deposited on : 2022-06-16
Resolution : 4.70 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

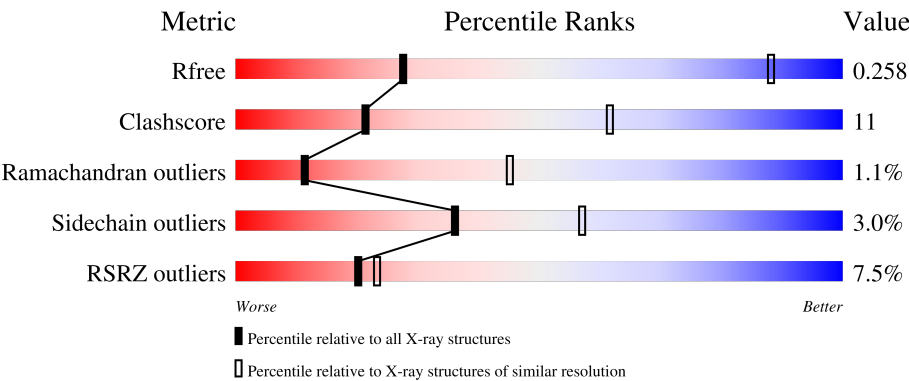
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 4.70 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	1020 (5.48-3.92)
Clashscore	190562	1005 (5.40-3.96)
Ramachandran outliers	187476	1095 (5.50-3.90)
Sidechain outliers	187428	1077 (5.50-3.90)
RSRZ outliers	180081	1015 (5.48-3.92)

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	694	9% 70% 25% • •
1	B	694	7% 70% 25% • •
2	C	552	6% 68% 26% • 5%
2	D	552	7% 69% 25% • 5%
3	E	2	50% 50%

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Mol	Chain	Length	Quality of chain
3	F	2	 50%50%
3	G	2	 50%50%
3	H	2	 100%
3	I	2	 100%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 19075 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 Plexin-A1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	670	Total	C	N	O	S	0	0	0
			5235	3304	906	992	33			
1	A	670	Total	C	N	O	S	0	0	0
			5235	3304	906	992	33			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	25	GLY	-	expression tag	UNP Q9UIW2
B	26	ARG	-	expression tag	UNP Q9UIW2
B	709	SER	-	expression tag	UNP Q9UIW2
B	710	ARG	-	expression tag	UNP Q9UIW2
B	711	LEU	-	expression tag	UNP Q9UIW2
B	712	GLU	-	expression tag	UNP Q9UIW2
B	713	ASN	-	expression tag	UNP Q9UIW2
B	714	LEU	-	expression tag	UNP Q9UIW2
B	715	TYR	-	expression tag	UNP Q9UIW2
B	716	PHE	-	expression tag	UNP Q9UIW2
B	717	GLN	-	expression tag	UNP Q9UIW2
B	718	GLY	-	expression tag	UNP Q9UIW2
A	25	GLY	-	expression tag	UNP Q9UIW2
A	26	ARG	-	expression tag	UNP Q9UIW2
A	709	SER	-	expression tag	UNP Q9UIW2
A	710	ARG	-	expression tag	UNP Q9UIW2
A	711	LEU	-	expression tag	UNP Q9UIW2
A	712	GLU	-	expression tag	UNP Q9UIW2
A	713	ASN	-	expression tag	UNP Q9UIW2
A	714	LEU	-	expression tag	UNP Q9UIW2
A	715	TYR	-	expression tag	UNP Q9UIW2
A	716	PHE	-	expression tag	UNP Q9UIW2
A	717	GLN	-	expression tag	UNP Q9UIW2
A	718	GLY	-	expression tag	UNP Q9UIW2

- Molecule 2 is a protein called Semaphorin 6D.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	527	Total	C	N	O	S	0	0	0
			4213	2673	730	784	26			
2	C	525	Total	C	N	O	S	0	0	0
			4196	2664	725	781	26			

There are 8 discrepancies between the modelled and reference sequences:

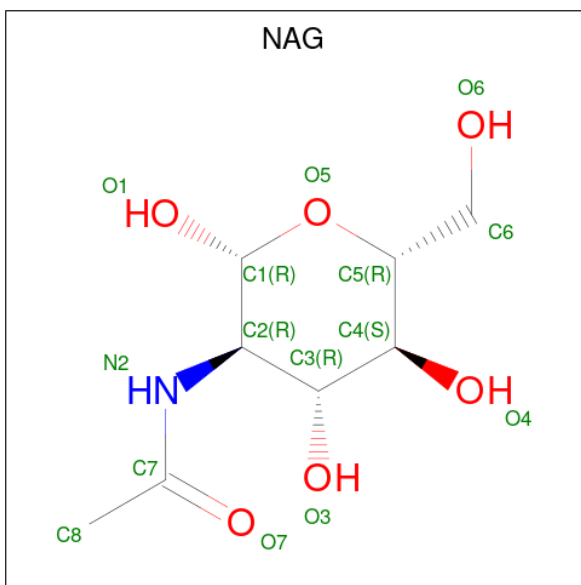
Chain	Residue	Modelled	Actual	Comment	Reference
D	19	SER	-	expression tag	UNP A0A0G2JZC4
D	20	SER	-	expression tag	UNP A0A0G2JZC4
D	21	ARG	-	expression tag	UNP A0A0G2JZC4
D	332	GLY	SER	engineered mutation	UNP A0A0G2JZC4
C	19	SER	-	expression tag	UNP A0A0G2JZC4
C	20	SER	-	expression tag	UNP A0A0G2JZC4
C	21	ARG	-	expression tag	UNP A0A0G2JZC4
C	332	GLY	SER	engineered mutation	UNP A0A0G2JZC4

- 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	F	2	Total	C	N	O		0	0	0
			28	16	2	10				
3	H	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				
3	G	2	Total	C	N	O		0	0	0
			28	16	2	10				
3	I	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) (labeled as "Ligand of Interest" by depositor).



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

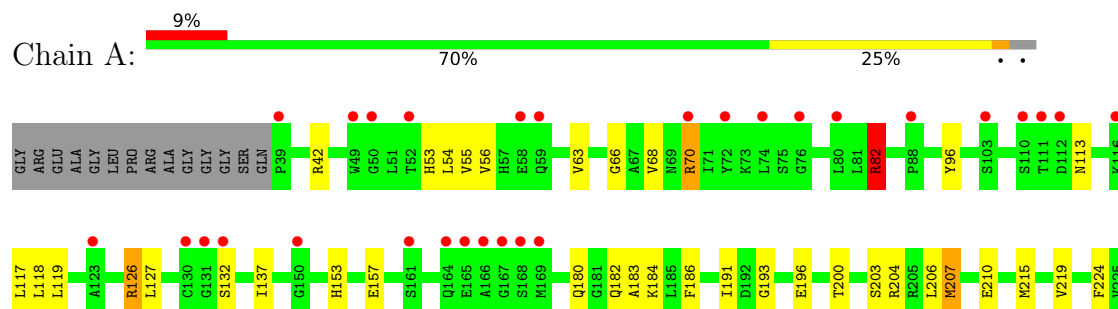
3 Residue-property plots [i](#)

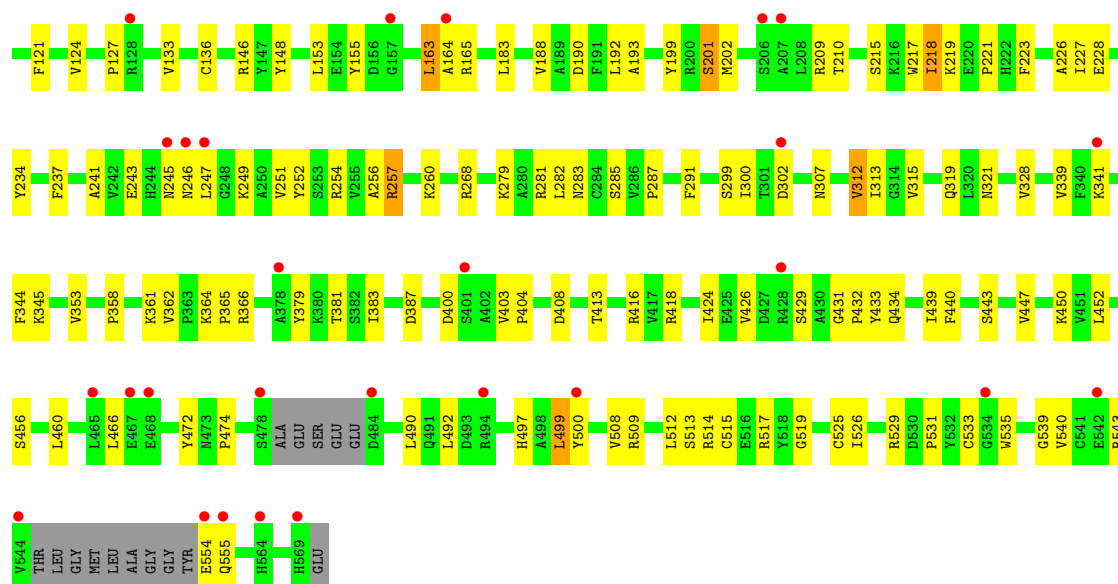
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Plexin-A1



• Molecule 1: Plexin-A1





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

Chain F: 50%



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

Chain H: 100%



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

Chain E: 50%



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

Chain G: 50%



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

Chain I:

100%

MAG1
MAG2

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	173.14Å 198.51Å 180.50Å 90.00° 117.83° 90.00°	Depositor
Resolution (Å)	49.43 – 4.70 49.43 – 4.70	Depositor EDS
% Data completeness (in resolution range)	95.1 (49.43-4.70) 93.2 (49.43-4.70)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.36 (at 4.64Å)	Xtriage
Refinement program	PHENIX 1.13_2998+SVN	Depositor
R, R_{free}	0.241 , 0.258 0.239 , 0.258	Depositor DCC
R_{free} test set	2523 reflections (4.48%)	wwPDB-VP
Wilson B-factor (Å ²)	157.7	Xtriage
Anisotropy	0.081	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.23 , 134.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.39$, $\langle L^2 \rangle = 0.22$	Xtriage
Estimated twinning fraction	0.059 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.87	EDS
Total number of atoms	19075	wwPDB-VP
Average B, all atoms (Å ²)	200.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.58% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: NAG

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	A	0.33	0/5357	0.77	4/7272 (0.1%)
1	B	0.30	0/5357	0.77	3/7272 (0.0%)
2	C	0.35	2/4298 (0.0%)	0.78	1/5821 (0.0%)
2	D	0.28	0/4315	0.81	6/5843 (0.1%)
All	All	0.32	2/19327 (0.0%)	0.78	14/26208 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	4
1	B	0	2
All	All	0	6

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	87	GLN	CA-C	12.34	1.60	1.53
2	C	404	PRO	CA-C	-5.45	1.48	1.52

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	572	ASN	N-CA-CB	-5.98	100.42	111.13
1	A	353	LYS	CD-CE-NZ	-5.75	93.50	111.90
1	A	70	ARG	CG-CD-NE	-5.72	99.42	112.00
2	D	66	ARG	NE-CZ-NH2	5.61	124.25	119.20
2	C	404	PRO	O-C-N	-5.41	118.61	121.15

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	484	ASP	CA-C-N	5.41	131.87	121.54
2	D	484	ASP	C-N-CA	5.41	131.87	121.54
1	A	82	ARG	NE-CZ-NH2	-5.35	114.39	119.20
1	B	80	LEU	CA-CB-CG	5.28	134.77	116.30
1	B	555	ARG	CB-CG-CD	-5.18	99.39	111.30
2	D	66	ARG	NE-CZ-NH1	-5.13	116.37	121.50
2	D	204	ASP	CA-C-N	5.12	131.44	121.41
2	D	204	ASP	C-N-CA	5.12	131.44	121.41
1	B	181	GLY	N-CA-C	5.12	125.31	113.18

There are no chirality outliers.

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	126	ARG	Mainchain
1	A	525	LEU	Mainchain
1	A	571	ARG	Mainchain
1	A	82	ARG	Mainchain
1	B	651	PHE	Mainchain
1	B	82	ARG	Mainchain

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	5235	0	5084	123	0
1	B	5235	0	5085	115	0
2	C	4196	0	4075	100	0
2	D	4213	0	4093	90	0
3	E	28	0	25	1	0
3	F	28	0	25	0	0
3	G	28	0	25	1	0
3	H	28	0	25	0	0
3	I	28	0	25	2	0
4	A	28	0	26	2	0
4	B	28	0	26	2	0
All	All	19075	0	18514	411	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 11.

All (411) 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:632:GLN:HG2	1:A:636:ARG:HB2	1.52	0.92
1:A:393:LEU:HD21	2:C:192:LEU:HD22	1.58	0.85
1:B:239:PHE:HZ	1:B:265:THR:HG22	1.43	0.83
2:D:123:LYS:HB3	2:D:174:ASN:HD21	1.43	0.83
2:D:58:ASP:HB2	2:D:74:ARG:HB2	1.60	0.83
2:D:43:VAL:HG21	2:D:476:LYS:HG2	1.62	0.81
2:C:165:ARG:HH21	2:C:201:SER:HB3	1.43	0.79
1:A:224:PHE:HE1	2:C:164:ALA:HB2	1.49	0.78
1:B:239:PHE:CZ	1:B:265:THR:HG22	2.21	0.75
2:D:218:ILE:HG12	2:D:221:PRO:HG3	1.66	0.75
2:D:426:VAL:HG22	2:D:439:ILE:HG12	1.68	0.74
2:D:490:LEU:HD13	2:D:499:LEU:HD21	1.69	0.74
1:A:324:LEU:HB2	1:A:446:THR:HG21	1.68	0.73
2:D:429:SER:HB2	2:D:434:GLN:HB3	1.70	0.73
2:D:321:ASN:ND2	2:C:287:PRO:O	2.21	0.73
2:D:529:ARG:NH2	2:D:558:GLU:O	2.23	0.72
1:A:426:LEU:HA	1:A:471:PRO:HG2	1.71	0.72
1:B:327:GLN:NE2	1:B:463:ASP:OD1	2.22	0.70
1:A:268:THR:HG22	1:A:277:PHE:H	1.58	0.69
2:C:345:LYS:HE2	2:C:400:ASP:HA	1.72	0.69
2:C:513:SER:OG	2:C:515:CYS:SG	2.44	0.69
2:D:103:GLN:O	2:D:107:ASN:ND2	2.26	0.69
1:A:56:VAL:HG22	1:A:63:VAL:HG22	1.73	0.69
1:A:700:VAL:HG13	1:A:705:ASP:HB2	1.76	0.67
2:C:500:TYR:CE2	2:C:509:ARG:HG3	2.30	0.67
1:B:324:LEU:HG	1:B:446:THR:HG21	1.76	0.67
1:A:389:LEU:HD21	1:A:414:PRO:HD2	1.75	0.67
1:B:370:LYS:O	1:B:374:GLN:HB2	1.95	0.67
1:B:459:LYS:NZ	1:B:512:VAL:O	2.29	0.66
1:A:706:CYS:HB3	1:A:707:PRO:HD2	1.78	0.65
2:D:33:VAL:HG13	2:D:37:TYR:HB3	1.79	0.65
1:A:649:LYS:HD2	1:A:650:LYS:H	1.61	0.65
2:D:124:VAL:HB	2:D:136:CYS:HB2	1.78	0.65
1:A:239:PHE:CZ	1:A:265:THR:HG22	2.32	0.65
1:B:155:ARG:HB2	1:B:158:HIS:CD2	2.32	0.64
2:C:513:SER:OG	2:C:533:CYS:SG	2.55	0.64
2:C:431:GLY:HA2	2:C:531:PRO:HB2	1.80	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:287:PRO:O	2:C:321:ASN:ND2	2.29	0.63
1:B:566:LEU:HB3	1:B:654:VAL:HB	1.81	0.63
1:A:446:THR:HG23	1:A:462:VAL:HG23	1.81	0.62
2:C:42:PRO:HG2	2:C:509:ARG:HB3	1.80	0.62
1:A:347:ASN:O	1:A:349:VAL:N	2.31	0.62
1:A:649:LYS:HD2	1:A:650:LYS:N	2.15	0.62
2:D:433:TYR:CD2	3:E:1:NAG:H62	2.35	0.61
1:A:426:LEU:HD13	1:A:462:VAL:HG11	1.83	0.61
1:A:203:SER:HB3	1:A:227:SER:HB2	1.80	0.61
1:A:118:LEU:HD11	1:A:127:LEU:HD21	1.80	0.61
1:B:373:ILE:HG22	1:B:415:LEU:HD23	1.83	0.61
2:C:65:ILE:HD12	2:C:70:TYR:HE2	1.66	0.61
1:A:323:ALA:O	1:A:327:GLN:HG3	2.01	0.61
1:B:196:GLU:OE2	2:D:394:LYS:NZ	2.32	0.60
1:B:466:ASN:O	1:B:468:GLY:N	2.30	0.60
2:D:35:TYR:CE1	2:D:474:PRO:HD2	2.35	0.60
2:D:454:LYS:HE3	2:D:460:LEU:O	2.01	0.60
1:B:411:PHE:HA	2:D:192:LEU:HD21	1.82	0.60
1:A:461:LEU:HD22	1:A:530:PRO:HG3	1.84	0.60
1:B:582:LEU:HD13	1:B:584:LEU:HD13	1.83	0.60
1:B:286:CYS:HB2	1:B:289:ASP:HB2	1.84	0.60
2:C:344:PHE:HE1	2:C:366:ARG:NH1	1.99	0.60
2:D:244:HIS:CE1	2:D:247:LEU:HD21	2.37	0.60
2:C:268:ARG:NE	2:C:387:ASP:OD2	2.34	0.60
1:B:700:VAL:HG13	1:B:705:ASP:HB2	1.84	0.59
2:C:315:VAL:HG22	2:C:328:VAL:HG22	1.83	0.59
2:D:342:GLY:O	2:D:366:ARG:NH2	2.34	0.59
1:A:373:ILE:HG22	1:A:415:LEU:HD23	1.84	0.58
2:D:62:MET:HE3	2:D:502:ALA:HB2	1.84	0.58
1:A:184:LYS:HD3	1:A:219:VAL:HG21	1.84	0.58
1:B:268:THR:HG22	1:B:277:PHE:H	1.68	0.58
2:C:228:GLU:HG3	2:C:260:LYS:HE2	1.86	0.58
1:B:451:GLY:HA2	1:B:457:ILE:HD13	1.86	0.58
2:D:62:MET:HB2	2:D:69:LEU:HD11	1.85	0.58
2:C:96:LEU:HD12	2:C:153:LEU:HB3	1.86	0.58
1:B:298:PHE:HZ	1:B:372:ARG:HB2	1.69	0.57
1:A:370:LYS:O	1:A:374:GLN:HB2	2.05	0.57
1:B:582:LEU:HD11	1:B:617:CYS:SG	2.45	0.57
1:A:537:LEU:HD21	1:A:646:GLU:HB3	1.87	0.57
2:D:339:VAL:HG13	2:D:401:SER:HB2	1.87	0.57
2:C:365:PRO:HG3	2:C:379:TYR:CE2	2.40	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:224:PHE:CD1	2:C:163:LEU:HD23	2.40	0.57
1:A:348:ARG:O	1:A:351:PRO:HD3	2.04	0.57
2:C:440:PHE:CE1	2:C:450:LYS:HG3	2.39	0.57
2:D:561:ASN:C	2:D:563:ALA:H	2.12	0.57
1:B:173:LEU:HG	1:B:185:LEU:HD13	1.86	0.56
1:A:566:LEU:HB3	1:A:654:VAL:HB	1.87	0.56
1:A:687:THR:HG21	1:A:692:ASP:HB2	1.86	0.56
1:A:96:TYR:O	2:C:268:ARG:HG2	2.05	0.56
2:C:146:ARG:HD3	2:C:155:TYR:CD1	2.41	0.56
1:A:641:TYR:CD2	1:A:650:LYS:HD2	2.40	0.56
1:B:281:LYS:HD3	1:B:299:PRO:HG3	1.86	0.56
1:A:281:LYS:HD3	1:A:299:PRO:HG3	1.88	0.56
1:A:536:VAL:HG21	1:A:590:PRO:HD3	1.88	0.56
2:C:433:TYR:CG	3:I:1:NAG:H62	2.40	0.56
1:B:411:PHE:CE1	2:D:111:LYS:HA	2.40	0.56
2:D:285:SER:O	2:D:410:PRO:HD3	2.06	0.56
1:A:429:ASP:OD1	1:A:434:LEU:HG	2.05	0.56
1:A:597:ASN:HB2	1:A:643:LYS:HB3	1.88	0.56
2:C:190:ASP:OD1	2:C:193:ALA:N	2.38	0.55
2:C:429:SER:HB2	2:C:434:GLN:HB3	1.87	0.55
2:C:65:ILE:HD12	2:C:70:TYR:CE2	2.41	0.55
2:C:241:ALA:HB2	2:C:252:TYR:HB2	1.87	0.55
2:D:497:HIS:CE1	2:D:512:LEU:HB2	2.42	0.55
1:A:226:SER:O	1:A:228:GLN:N	2.38	0.55
2:C:215:SER:HB2	2:C:219:LYS:HD2	1.89	0.55
2:C:339:VAL:HG21	2:C:403:VAL:HG22	1.88	0.55
1:B:42:ARG:HB2	1:B:509:ARG:HB3	1.89	0.55
2:C:127:PRO:HA	2:C:133:VAL:HA	1.89	0.55
1:B:140:PHE:HE1	1:B:216:PHE:CZ	2.25	0.55
1:A:462:VAL:HG12	1:A:472:ALA:HB2	1.88	0.55
2:D:25:GLU:HG2	2:D:26:ASP:H	1.72	0.55
1:B:470:ARG:HH21	1:B:470:ARG:HB2	1.73	0.54
2:D:531:PRO:HA	2:D:557:THR:HG21	1.90	0.54
1:B:239:PHE:HD1	1:B:242:PHE:HB2	1.73	0.54
1:B:383:LEU:HD23	1:B:383:LEU:H	1.72	0.54
1:A:186:PHE:O	1:A:258:TYR:OH	2.26	0.54
1:A:683:ARG:HG2	1:A:685:VAL:HG13	1.90	0.54
2:D:227:ILE:HG21	2:D:303:ILE:HG13	1.90	0.54
2:C:44:PHE:HZ	2:C:78:TYR:CG	2.25	0.54
2:C:513:SER:O	2:C:543:ARG:NH2	2.41	0.53
1:B:513:GLU:HG2	1:B:515:CYS:SG	2.49	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:405:ASP:OD1	1:A:405:ASP:N	2.42	0.53
2:C:246:ASN:O	2:C:247:LEU:HG	2.08	0.53
1:A:699:ARG:HD2	1:A:699:ARG:O	2.08	0.53
2:C:217:TRP:O	2:C:279:LYS:NZ	2.29	0.53
2:C:526:ILE:HD11	2:C:535:TRP:HB3	1.90	0.53
2:D:123:LYS:HB3	2:D:174:ASN:ND2	2.19	0.53
2:C:452:LEU:HB2	2:C:466:LEU:HD11	1.91	0.53
2:D:449:LEU:HD23	2:D:468:GLU:HG3	1.90	0.53
1:A:239:PHE:HZ	1:A:265:THR:HG22	1.73	0.53
2:C:440:PHE:HE1	2:C:450:LYS:HG3	1.73	0.53
2:C:472:TYR:O	2:C:474:PRO:HD3	2.09	0.53
2:C:490:LEU:HD13	2:C:499:LEU:HD21	1.91	0.53
1:B:585:GLN:HG3	1:B:614:ARG:HG3	1.91	0.52
1:B:640:LEU:HB3	1:B:654:VAL:HG13	1.91	0.52
2:D:283:ASN:O	2:D:405:PRO:HD3	2.09	0.52
1:B:671:VAL:HG11	1:B:703:SER:N	2.25	0.52
1:A:362:LEU:HA	1:A:365:ILE:HD12	1.92	0.52
1:A:568:VAL:HG22	1:A:584:LEU:HG	1.90	0.52
2:C:381:THR:HG23	2:C:383:ILE:H	1.74	0.52
2:C:433:TYR:CD2	3:I:1:NAG:H62	2.45	0.52
2:C:492:LEU:HA	2:C:499:LEU:HD12	1.92	0.52
1:B:234:ASP:O	1:B:237:SER:HB3	2.09	0.52
1:B:362:LEU:HA	1:B:365:ILE:HD12	1.91	0.52
1:A:560:LEU:C	1:A:562:GLN:H	2.18	0.52
2:C:74:ARG:C	2:C:76:GLN:H	2.16	0.52
1:B:139:GLN:HG2	1:B:151:GLU:HG3	1.92	0.52
1:B:466:ASN:C	1:B:468:GLY:H	2.16	0.52
2:D:345:LYS:HB2	2:D:400:ASP:HA	1.91	0.52
1:A:349:VAL:HG13	1:A:350:LYS:HG2	1.91	0.52
2:D:454:LYS:HG3	2:D:459:SER:HB3	1.92	0.52
2:D:147:TYR:CE2	2:D:202:MET:HE1	2.45	0.52
1:B:378:ARG:HD3	1:B:380:GLU:OE2	2.09	0.51
1:B:393:LEU:HD22	2:D:192:LEU:HB3	1.92	0.51
1:A:451:GLY:HA2	1:A:457:ILE:HD13	1.92	0.51
1:B:126:ARG:HH11	1:B:210:GLU:HG3	1.75	0.51
1:B:152:PRO:HB2	1:B:158:HIS:CE1	2.45	0.51
1:A:491:VAL:O	1:A:499:LEU:HD12	2.10	0.51
1:B:231:ILE:HD11	1:B:244:ILE:HD12	1.93	0.51
1:A:299:PRO:HB2	1:A:420:THR:HG22	1.92	0.51
1:B:298:PHE:HZ	1:B:372:ARG:CB	2.24	0.50
1:B:42:ARG:NH1	4:B:1001:NAG:H81	2.26	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:180:GLN:O	1:B:182:GLN:N	2.44	0.50
1:A:377:TYR:HE2	1:A:400:LEU:HD23	1.77	0.50
2:D:147:TYR:HE2	2:D:202:MET:HE1	1.76	0.50
2:C:251:VAL:H	2:C:319:GLN:NE2	2.08	0.50
1:A:137:ILE:HG21	1:A:153:HIS:HB3	1.93	0.50
2:C:94:LYS:HB3	2:C:153:LEU:HD11	1.93	0.50
1:B:391:LYS:H	1:B:391:LYS:HD2	1.77	0.50
1:A:196:GLU:OE1	1:A:196:GLU:N	2.42	0.50
2:C:533:CYS:O	2:C:555:GLN:NE2	2.45	0.50
2:D:344:PHE:HE1	2:D:366:ARG:NH1	2.10	0.50
2:C:30:LEU:HD12	2:C:466:LEU:HA	1.93	0.50
2:C:241:ALA:HB2	2:C:252:TYR:CG	2.46	0.50
2:C:514:ARG:HE	2:C:517:ARG:HB2	1.76	0.50
1:B:362:LEU:O	1:B:366:LYS:HG3	2.12	0.50
1:B:374:GLN:HG3	1:B:378:ARG:NH1	2.27	0.50
2:D:315:VAL:HG22	2:D:328:VAL:HG22	1.94	0.50
1:A:544:ARG:O	1:A:544:ARG:HG3	2.12	0.49
1:A:224:PHE:CG	2:C:163:LEU:HD23	2.47	0.49
2:C:35:TYR:CE1	2:C:474:PRO:HD2	2.47	0.49
1:A:117:LEU:HD21	1:A:119:LEU:HD12	1.93	0.49
2:C:256:ALA:HA	2:C:279:LYS:HA	1.94	0.49
1:B:117:LEU:HD21	1:B:119:LEU:HD12	1.94	0.49
1:A:534:TRP:HB3	1:A:557:ALA:HB3	1.95	0.49
1:B:456:ARG:HD3	1:B:477:SER:OG	2.13	0.49
1:A:322:ARG:H	1:A:444:GLY:HA2	1.77	0.49
1:A:674:SER:HB3	4:A:1002:NAG:H82	1.93	0.49
2:D:536:LEU:HD11	2:D:542:GLU:HB2	1.94	0.49
2:D:101:ARG:O	2:D:105:ARG:HG3	2.13	0.49
2:D:497:HIS:HE1	2:D:513:SER:H	1.61	0.49
2:C:199:TYR:CE2	2:C:201:SER:HB2	2.48	0.49
1:B:537:LEU:O	1:B:647:THR:HG21	2.13	0.49
2:C:365:PRO:HG3	2:C:379:TYR:HE2	1.76	0.48
1:B:251:ARG:HG2	1:B:256:VAL:HG12	1.94	0.48
2:C:124:VAL:HB	2:C:136:CYS:HB2	1.95	0.48
1:B:525:LEU:HD11	1:B:557:ALA:O	2.13	0.48
2:D:406:ILE:C	2:D:408:ASP:H	2.20	0.48
1:B:96:TYR:CE1	1:B:197:TYR:HB3	2.49	0.48
2:D:440:PHE:CE1	2:D:450:LYS:HG3	2.48	0.48
1:B:690:VAL:HG21	1:B:699:ARG:HD2	1.93	0.48
1:A:706:CYS:HB3	1:A:707:PRO:CD	2.44	0.48
2:D:174:ASN:OD1	2:D:174:ASN:N	2.46	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:234:TYR:CD1	2:C:257:ARG:HB2	2.49	0.48
1:B:495:ASN:OD1	1:B:495:ASN:N	2.44	0.48
1:A:273:ALA:HB3	1:A:275:GLU:OE1	2.14	0.48
2:D:472:TYR:HB2	2:D:503:PHE:CE2	2.49	0.48
2:C:364:LYS:HD3	2:C:364:LYS:HA	1.50	0.48
1:A:582:LEU:HD11	1:A:617:CYS:SG	2.53	0.48
2:D:288:GLY:HA3	2:C:321:ASN:ND2	2.29	0.48
2:C:43:VAL:HG22	2:C:508:VAL:HG22	1.96	0.48
1:B:96:TYR:OH	1:B:156:LYS:HG3	2.14	0.48
2:C:268:ARG:HE	2:C:387:ASP:CG	2.21	0.48
1:B:470:ARG:HB2	1:B:470:ARG:NH2	2.28	0.47
1:B:683:ARG:HG2	1:B:685:VAL:HG13	1.95	0.47
1:A:157:GLU:O	1:A:204:ARG:NH1	2.47	0.47
2:C:77:VAL:HB	2:C:96:LEU:HB3	1.95	0.47
1:A:234:ASP:O	1:A:237:SER:OG	2.27	0.47
1:A:391:LYS:O	1:A:393:LEU:N	2.47	0.47
1:B:448:VAL:HB	1:B:460:ILE:HB	1.95	0.47
2:D:513:SER:OG	2:D:533:CYS:SG	2.73	0.47
1:B:200:THR:HG22	1:B:229:LEU:O	2.15	0.47
1:B:370:LYS:O	1:B:374:GLN:CB	2.62	0.47
1:B:155:ARG:HB2	1:B:158:HIS:CG	2.50	0.47
1:B:487:LEU:HD23	1:B:487:LEU:HA	1.71	0.47
1:A:574:SER:HB2	1:A:659:TYR:CE2	2.50	0.47
2:D:316:PHE:O	2:D:327:ALA:N	2.37	0.47
2:D:454:LYS:H	2:D:463:SER:HG	1.58	0.47
2:C:245:ASN:C	2:C:247:LEU:H	2.23	0.47
1:A:683:ARG:O	1:A:684:HIS:ND1	2.46	0.47
2:C:101:ARG:O	2:C:105:ARG:HG3	2.14	0.47
1:B:126:ARG:NH1	1:B:210:GLU:HG3	2.30	0.47
2:C:226:ALA:O	2:C:227:ILE:HG13	2.15	0.47
1:B:348:ARG:O	1:B:351:PRO:HD3	2.15	0.47
1:A:42:ARG:HG3	1:A:509:ARG:NH2	2.30	0.47
1:A:126:ARG:NH2	1:A:207:MET:O	2.41	0.47
2:D:363:PRO:O	2:D:364:LYS:HD3	2.15	0.47
1:B:405:ASP:OD1	1:B:405:ASP:N	2.48	0.46
1:A:383:LEU:HD23	1:A:383:LEU:H	1.78	0.46
2:C:358:PRO:HD2	2:C:361:LYS:HD2	1.97	0.46
1:A:495:ASN:OD1	1:A:495:ASN:N	2.44	0.46
1:B:349:VAL:HG13	1:B:350:LYS:HG2	1.97	0.46
2:C:313:ILE:HG22	2:C:424:ILE:HD11	1.97	0.46
1:B:141:LEU:HB3	1:B:146:LEU:HA	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:167:PRO:HG3	2:D:174:ASN:HB3	1.97	0.46
1:A:612:ASP:OD1	1:A:612:ASP:N	2.47	0.46
2:C:227:ILE:HD11	2:C:302:ASP:HA	1.96	0.46
1:B:226:SER:O	1:B:228:GLN:N	2.40	0.46
1:B:157:GLU:O	1:B:204:ARG:NH1	2.48	0.46
1:A:569:GLN:HA	1:A:570:PRO:HA	1.62	0.46
1:B:568:VAL:HG22	1:B:584:LEU:HG	1.98	0.46
2:C:341:LYS:HA	2:C:366:ARG:HH21	1.80	0.46
1:A:183:ALA:O	1:A:206:LEU:HB2	2.14	0.46
1:B:382:LYS:HB2	1:B:394:GLY:HA2	1.98	0.45
1:A:463:ASP:O	1:A:471:PRO:HA	2.16	0.45
2:C:74:ARG:HB2	2:C:121:PHE:CE1	2.51	0.45
1:B:490:LEU:HD22	1:B:499:LEU:HD11	1.98	0.45
1:A:579:GLN:HA	1:A:619:SER:O	2.16	0.45
2:C:432:PRO:HG3	2:C:555:GLN:OE1	2.16	0.45
1:B:239:PHE:CD1	1:B:242:PHE:HB2	2.50	0.45
2:D:190:ASP:OD1	2:D:193:ALA:N	2.50	0.45
2:D:431:GLY:HA2	2:D:531:PRO:HB2	1.98	0.45
2:C:443:SER:HG	2:C:447:VAL:H	1.65	0.45
1:A:54:LEU:HD23	1:A:55:VAL:N	2.31	0.45
1:A:700:VAL:HG13	1:A:705:ASP:CB	2.46	0.45
2:C:188:VAL:HG11	2:C:193:ALA:HA	1.98	0.45
1:B:555:ARG:HD2	1:B:590:PRO:HG3	1.98	0.45
2:C:500:TYR:CZ	2:C:509:ARG:HG3	2.52	0.45
2:C:426:VAL:HG22	2:C:439:ILE:HG12	1.99	0.45
1:B:203:SER:HB3	1:B:227:SER:HB2	1.99	0.45
2:C:243:GLU:OE1	2:C:243:GLU:N	2.45	0.45
1:A:298:PHE:HZ	1:A:372:ARG:CB	2.30	0.44
1:A:298:PHE:HZ	1:A:372:ARG:HB3	1.82	0.44
2:D:195:ASP:CG	2:D:212:LYS:HG3	2.42	0.44
1:B:58:GLU:O	1:B:494:PRO:HB3	2.17	0.44
2:D:42:PRO:HG2	2:D:509:ARG:HB3	1.98	0.44
1:A:574:SER:OG	1:A:576:THR:HG23	2.17	0.44
2:C:165:ARG:NH2	2:C:201:SER:HB3	2.21	0.44
1:A:348:ARG:HE	1:A:348:ARG:HB3	1.63	0.44
1:A:113:ASN:OD1	1:A:132:SER:N	2.45	0.44
1:A:268:THR:CG2	1:A:277:PHE:H	2.30	0.44
1:A:625:VAL:HA	1:A:628:ILE:HD12	1.98	0.44
1:B:254:GLN:O	1:B:287:VAL:HG22	2.18	0.44
1:A:207:MET:SD	1:A:215:MET:HA	2.57	0.44
1:A:324:LEU:HA	1:A:327:GLN:HB2	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:155:TYR:OH	2:D:158:GLU:HA	2.16	0.44
1:A:481:GLN:OE1	1:A:505:LYS:HB2	2.17	0.44
1:A:632:GLN:O	1:A:632:GLN:HG3	2.18	0.44
1:B:675:PHE:HB3	1:B:676:PRO:HD2	2.00	0.44
1:A:261:LEU:HD11	1:A:278:PHE:HB3	2.00	0.44
1:A:402:ILE:HG22	1:A:403:ASP:O	2.18	0.44
1:A:537:LEU:HD13	1:A:555:ARG:NH1	2.33	0.44
1:B:54:LEU:HD23	1:B:55:VAL:N	2.33	0.43
1:B:191:ILE:HD12	1:B:193:GLY:N	2.33	0.43
2:C:223:PHE:HA	2:C:237:PHE:CB	2.47	0.43
1:A:699:ARG:NH1	1:A:701:ASN:ND2	2.66	0.43
2:D:65:ILE:HD12	2:D:70:TYR:HE2	1.83	0.43
1:A:286:CYS:HB2	1:A:289:ASP:HB2	1.99	0.43
2:D:319:GLN:O	2:D:322:SER:OG	2.28	0.43
1:B:116:LYS:HG2	1:B:166:ALA:HA	1.99	0.43
1:A:200:THR:HG22	1:A:229:LEU:O	2.18	0.43
1:A:369:ILE:O	1:A:373:ILE:HG23	2.18	0.43
2:C:281:ARG:HG2	2:C:282:LEU:N	2.33	0.43
2:C:283:ASN:OD1	2:C:285:SER:HB3	2.18	0.43
1:B:687:THR:HG21	1:B:692:ASP:HB2	2.00	0.43
2:C:243:GLU:CD	2:C:254:ARG:HH22	2.23	0.43
1:B:426:LEU:O	1:B:472:ALA:HB3	2.19	0.43
2:D:220:GLU:HB2	2:D:240:ILE:HD12	2.01	0.43
2:D:349:THR:OG1	2:D:352:SER:HB3	2.19	0.43
2:D:489:SER:O	2:D:490:LEU:HD23	2.18	0.43
2:D:490:LEU:HD21	2:D:501:VAL:HG13	2.00	0.43
1:A:53:HIS:O	1:A:66:GLY:N	2.46	0.43
2:D:221:PRO:HB3	2:D:239:GLU:HB3	2.01	0.43
2:C:300:ILE:HD11	2:C:312:VAL:HB	2.00	0.43
2:C:62:MET:HB2	2:C:69:LEU:HD11	2.01	0.43
1:B:415:LEU:HD12	1:B:415:LEU:HA	1.85	0.43
2:D:108:CYS:HB2	2:D:140:ALA:HB1	2.00	0.43
2:C:96:LEU:HD11	2:C:148:TYR:CG	2.54	0.43
1:B:569:GLN:HA	1:B:570:PRO:HA	1.77	0.43
1:A:383:LEU:HD12	1:A:389:LEU:HD11	2.01	0.43
2:D:25:GLU:HG2	2:D:26:ASP:N	2.34	0.43
1:B:393:LEU:HD21	2:D:192:LEU:HD22	2.01	0.42
2:D:436:TYR:HB3	2:D:452:LEU:HD11	1.99	0.42
1:A:226:SER:C	1:A:228:GLN:N	2.77	0.42
1:A:311:LEU:HB2	1:A:342:ALA:HB3	2.00	0.42
2:D:228:GLU:HG3	2:D:260:LYS:HE2	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:196:GLU:H	1:A:196:GLU:CD	2.26	0.42
2:D:235:PHE:HB2	2:D:256:ALA:HB3	2.01	0.42
1:B:560:LEU:C	1:B:562:GLN:H	2.28	0.42
1:A:470:ARG:CZ	1:A:470:ARG:HB2	2.50	0.42
2:D:163:LEU:HD12	2:D:163:LEU:HA	1.72	0.42
1:B:96:TYR:O	2:D:268:ARG:HG2	2.20	0.42
1:B:174:ILE:HG12	1:B:251:ARG:HD2	2.01	0.42
1:A:126:ARG:HH11	1:A:210:GLU:HG3	1.85	0.42
1:A:620:PRO:HB2	1:A:625:VAL:HG23	2.00	0.42
2:C:362:VAL:HG11	2:C:366:ARG:HH11	1.84	0.42
1:B:92:ASN:OD1	1:B:93:GLU:N	2.53	0.42
1:B:178:PRO:C	1:B:180:GLN:H	2.28	0.42
1:B:324:LEU:HD23	1:B:324:LEU:HA	1.79	0.42
1:B:383:LEU:HB3	1:B:415:LEU:HD12	2.00	0.42
1:B:492:LEU:HD12	1:B:499:LEU:HD13	2.02	0.42
2:D:214:ASP:C	2:D:216:LYS:H	2.27	0.42
2:D:297:LEU:HD13	2:D:316:PHE:CZ	2.55	0.42
2:D:353:VAL:HG22	2:C:353:VAL:HG22	2.00	0.42
1:B:42:ARG:HG3	1:B:509:ARG:NE	2.35	0.42
1:B:125:ASN:C	1:B:143:LEU:HG	2.45	0.42
1:A:68:VAL:C	1:A:70:ARG:H	2.28	0.42
1:A:555:ARG:HE	1:A:555:ARG:HB3	1.50	0.42
1:A:569:GLN:HB2	1:A:583:VAL:HG13	2.01	0.42
1:A:674:SER:HB3	4:A:1002:NAG:C8	2.50	0.42
2:D:65:ILE:HG22	2:D:66:ARG:HG2	2.01	0.42
2:D:74:ARG:C	2:D:76:GLN:H	2.28	0.42
2:D:230:GLY:O	2:D:260:LYS:HE3	2.19	0.42
2:C:209:ARG:NE	2:C:210:THR:O	2.48	0.42
1:B:592:LEU:C	1:B:594:ALA:N	2.77	0.41
2:D:497:HIS:CE1	2:D:513:SER:H	2.37	0.41
1:B:68:VAL:C	1:B:70:ARG:H	2.27	0.41
1:B:78:LEU:HD21	1:B:500:TYR:HD1	1.85	0.41
2:D:288:GLY:HA3	2:C:321:ASN:HD22	1.85	0.41
2:C:107:ASN:O	2:C:111:LYS:HG3	2.20	0.41
2:C:460:LEU:HA	2:C:460:LEU:HD23	1.85	0.41
1:A:191:ILE:HD12	1:A:193:GLY:N	2.35	0.41
1:A:466:ASN:HA	1:A:467:PRO:HD3	1.94	0.41
2:D:275:THR:O	2:D:393:ILE:HD13	2.20	0.41
2:D:490:LEU:HD22	2:D:501:VAL:HA	2.02	0.41
1:B:223:GLU:HB3	1:B:224:PHE:CE2	2.56	0.41
1:A:492:LEU:HD23	1:A:496:HIS:HA	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:678:HIS:ND1	1:A:701:ASN:OD1	2.53	0.41
2:D:347:GLN:HB2	2:D:354:TRP:CD2	2.55	0.41
2:C:497:HIS:CD2	2:C:512:LEU:HB2	2.56	0.41
1:B:161:SER:OG	1:B:163:VAL:HG22	2.21	0.41
1:B:490:LEU:HD23	1:B:490:LEU:HA	1.90	0.41
1:A:382:LYS:HB2	1:A:393:LEU:O	2.20	0.41
1:A:426:LEU:O	1:A:472:ALA:HB3	2.21	0.41
2:D:41:TYR:CD2	2:D:510:ILE:HD13	2.55	0.41
2:C:241:ALA:HB2	2:C:252:TYR:CB	2.50	0.41
1:B:650:LYS:HE3	1:B:650:LYS:HB2	1.90	0.41
1:A:372:ARG:NH2	1:A:382:LYS:O	2.54	0.41
1:A:622:ALA:HA	1:A:625:VAL:HB	2.02	0.41
2:D:315:VAL:HG23	2:D:424:ILE:HD12	2.03	0.41
2:C:26:ASP:OD2	2:C:416:ARG:HB2	2.21	0.41
2:C:37:TYR:HD1	2:C:514:ARG:NH2	2.19	0.41
2:D:533:CYS:O	2:D:555:GLN:NE2	2.53	0.41
2:C:307:ASN:OD1	2:C:456:SER:HB2	2.20	0.41
1:B:226:SER:C	1:B:228:GLN:N	2.79	0.41
1:B:311:LEU:HB2	1:B:342:ALA:HB3	2.01	0.41
1:A:82:ARG:HH21	1:A:82:ARG:HD2	1.51	0.41
1:A:180:GLN:HG2	1:A:182:GLN:H	1.86	0.41
1:A:426:LEU:HB3	1:A:472:ALA:HB2	2.02	0.41
2:C:535:TRP:CZ2	2:C:539:GLY:HA2	2.56	0.41
1:B:283:VAL:HA	1:B:296:VAL:O	2.21	0.41
1:A:383:LEU:HD21	1:A:393:LEU:HG	2.03	0.41
1:A:644:SER:HB2	1:A:651:PHE:CE2	2.56	0.41
2:D:183:LEU:O	2:D:200:ARG:HA	2.21	0.41
1:B:73:LYS:O	1:B:81:LEU:HB2	2.21	0.40
1:A:448:VAL:HB	1:A:460:ILE:HB	2.02	0.40
1:A:579:GLN:HE21	1:A:621:SER:HB3	1.86	0.40
1:A:621:SER:O	1:A:624:GLU:N	2.53	0.40
2:D:289:ASP:O	2:C:249:LYS:HB3	2.22	0.40
2:C:299:SER:OG	2:C:424:ILE:HG22	2.21	0.40
1:B:42:ARG:HH12	4:B:1001:NAG:H81	1.87	0.40
1:A:378:ARG:HD3	1:A:380:GLU:OE2	2.22	0.40
2:D:55:HIS:O	2:D:55:HIS:CG	2.74	0.40
2:C:291:PHE:CE2	3:G:1:NAG:H5	2.56	0.40
2:C:514:ARG:NE	2:C:517:ARG:HB2	2.36	0.40
1:B:92:ASN:HB2	1:B:108:LEU:HA	2.04	0.40
1:B:647:THR:HG23	1:B:649:LYS:H	1.86	0.40
1:B:699:ARG:HG2	1:B:701:ASN:HD21	1.86	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:78:LEU:HD21	1:B:500:TYR:CD1	2.57	0.40
1:B:152:PRO:HB2	1:B:158:HIS:ND1	2.37	0.40
1:B:183:ALA:HB3	1:B:206:LEU:HB2	2.03	0.40
1:B:207:MET:O	1:B:215:MET:HE1	2.21	0.40
2:C:218:ILE:HG12	2:C:221:PRO:HG3	2.04	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	668/694 (96%)	596 (89%)	67 (10%)	5 (1%)	18	55
1	B	668/694 (96%)	593 (89%)	64 (10%)	11 (2%)	7	37
2	C	517/552 (94%)	456 (88%)	58 (11%)	3 (1%)	21	59
2	D	519/552 (94%)	454 (88%)	59 (11%)	6 (1%)	10	43
All	All	2372/2492 (95%)	2099 (88%)	248 (10%)	25 (1%)	11	45

All (25) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	47	SER
1	B	106	HIS
1	B	181	GLY
1	B	273	ALA
1	A	684	HIS
2	D	466	LEU
2	D	485	ARG
2	D	543	ARG
2	C	525	CYS
1	B	108	LEU

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Mol	Chain	Res	Type
1	A	348	ARG
1	A	630	ARG
1	B	630	ARG
2	D	512	LEU
1	B	240	PRO
1	A	240	PRO
2	D	206	SER
2	C	519	GLY
1	B	148	LYS
1	B	347	ASN
1	A	561	LEU
1	B	105	PRO
2	C	408	ASP
1	B	467	PRO
2	D	374	GLY

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	580/596 (97%)	561 (97%)	19 (3%)	33	55
1	B	580/596 (97%)	562 (97%)	18 (3%)	35	56
2	C	461/482 (96%)	446 (97%)	15 (3%)	33	55
2	D	463/482 (96%)	453 (98%)	10 (2%)	45	64
All	All	2084/2156 (97%)	2022 (97%)	62 (3%)	36	57

All (62) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	82	ARG
1	B	116	LYS
1	B	148	LYS
1	B	196	GLU
1	B	207	MET
1	B	229	LEU

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Mol	Chain	Res	Type
1	B	235	THR
1	B	239	PHE
1	B	271	ASP
1	B	275	GLU
1	B	383	LEU
1	B	429	ASP
1	B	446	THR
1	B	528	ARG
1	B	544	ARG
1	B	554	GLN
1	B	583	VAL
1	B	695	PHE
1	A	207	MET
1	A	233	SER
1	A	237	SER
1	A	253	GLU
1	A	271	ASP
1	A	275	GLU
1	A	276	HIS
1	A	378	ARG
1	A	383	LEU
1	A	429	ASP
1	A	446	THR
1	A	528	ARG
1	A	543	ARG
1	A	582	LEU
1	A	610	LEU
1	A	612	ASP
1	A	649	LYS
1	A	655	ASP
1	A	695	PHE
2	D	87	GLN
2	D	93	SER
2	D	163	LEU
2	D	174	ASN
2	D	204	ASP
2	D	247	LEU
2	D	255	VAL
2	D	311	THR
2	D	456	SER
2	D	517	ARG
2	C	87	GLN

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Mol	Chain	Res	Type
2	C	95	LYS
2	C	163	LEU
2	C	183	LEU
2	C	201	SER
2	C	202	MET
2	C	218	ILE
2	C	257	ARG
2	C	312	VAL
2	C	413	THR
2	C	418	ARG
2	C	499	LEU
2	C	529	ARG
2	C	540	VAL
2	C	554	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	115	ASN
1	B	158	HIS
1	B	313	GLN
1	B	506	GLN
1	B	579	GLN
1	B	616	HIS
1	B	664	HIS
1	B	672	ASN
1	A	59	GLN
1	A	374	GLN
1	A	390	ASN
1	A	565	GLN
1	A	579	GLN
1	A	664	HIS
2	D	31	ASN
2	D	139	ASN
2	D	172	GLN
2	D	497	HIS
2	C	31	ASN
2	C	267	GLN
2	C	307	ASN
2	C	321	ASN
2	C	347	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 ⓘ

10 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	E	1	2,3	14,14,15	0.64	1 (7%)	17,19,21	0.51	0
3	NAG	E	2	3	14,14,15	0.50	0	17,19,21	0.45	0
3	NAG	F	1	3,1	14,14,15	0.72	1 (7%)	17,19,21	0.65	1 (5%)
3	NAG	F	2	3	14,14,15	0.59	0	17,19,21	0.50	0
3	NAG	G	1	2,3	14,14,15	0.65	0	17,19,21	0.61	1 (5%)
3	NAG	G	2	3	14,14,15	0.74	1 (7%)	17,19,21	0.48	0
3	NAG	H	1	2,3	14,14,15	0.59	0	17,19,21	0.55	0
3	NAG	H	2	3	14,14,15	0.51	0	17,19,21	0.46	0
3	NAG	I	1	2,3	14,14,15	0.46	0	17,19,21	0.56	0
3	NAG	I	2	3	14,14,15	0.73	1 (7%)	17,19,21	0.47	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	E	1	2,3	-	0/6/23/26	0/1/1/1
3	NAG	E	2	3	-	0/6/23/26	0/1/1/1
3	NAG	F	1	3,1	-	2/6/23/26	0/1/1/1
3	NAG	F	2	3	-	1/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	G	1	2,3	-	2/6/23/26	0/1/1/1
3	NAG	G	2	3	-	2/6/23/26	0/1/1/1
3	NAG	H	1	2,3	-	2/6/23/26	0/1/1/1
3	NAG	H	2	3	-	2/6/23/26	0/1/1/1
3	NAG	I	1	2,3	-	0/6/23/26	0/1/1/1
3	NAG	I	2	3	-	2/6/23/26	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	1	NAG	O5-C1	-2.30	1.39	1.43
3	I	2	NAG	C1-C2	2.27	1.55	1.52
3	F	1	NAG	C1-C2	2.16	1.55	1.52
3	G	2	NAG	C1-C2	2.07	1.55	1.52

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	1	NAG	C1-O5-C5	2.09	114.99	112.19
3	G	1	NAG	C1-O5-C5	2.04	114.92	112.19

There are no chirality outliers.

All (13) torsion outliers are listed below:

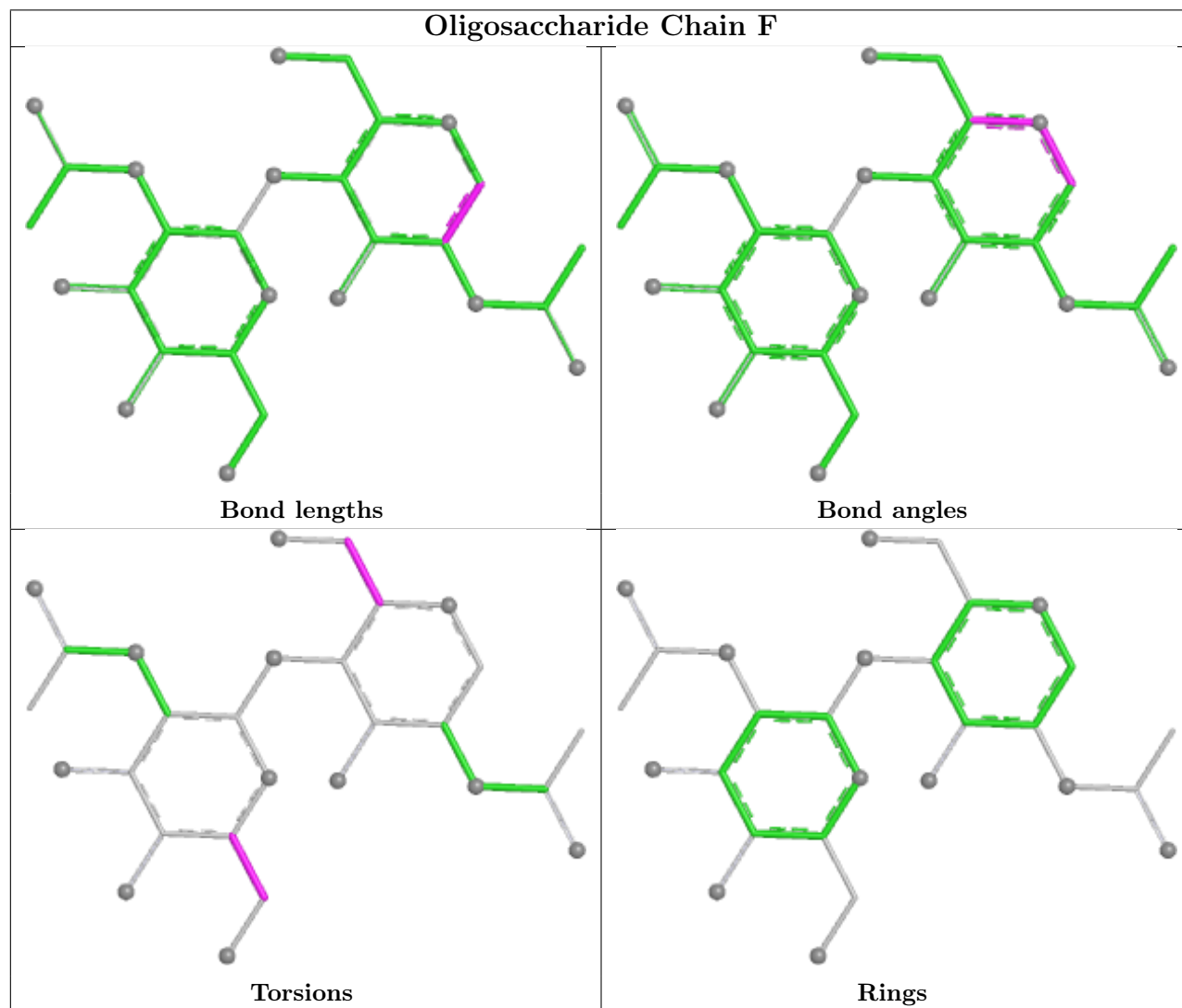
Mol	Chain	Res	Type	Atoms
3	H	2	NAG	O5-C5-C6-O6
3	I	2	NAG	O5-C5-C6-O6
3	H	2	NAG	C4-C5-C6-O6
3	G	2	NAG	O5-C5-C6-O6
3	I	2	NAG	C4-C5-C6-O6
3	G	1	NAG	C4-C5-C6-O6
3	G	1	NAG	O5-C5-C6-O6
3	H	1	NAG	C4-C5-C6-O6
3	H	1	NAG	O5-C5-C6-O6
3	G	2	NAG	C4-C5-C6-O6
3	F	2	NAG	O5-C5-C6-O6
3	F	1	NAG	C4-C5-C6-O6
3	F	1	NAG	O5-C5-C6-O6

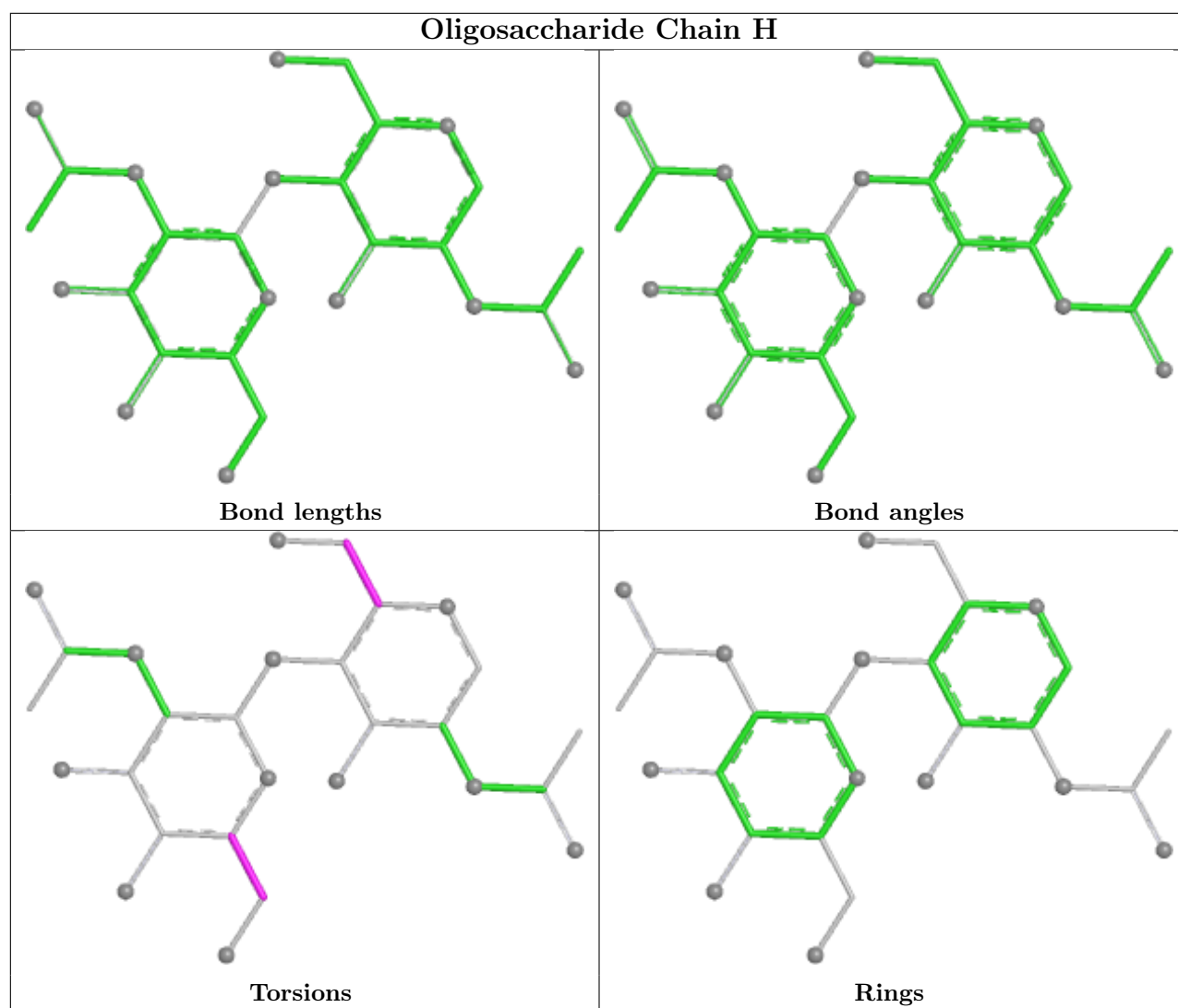
There are no ring outliers.

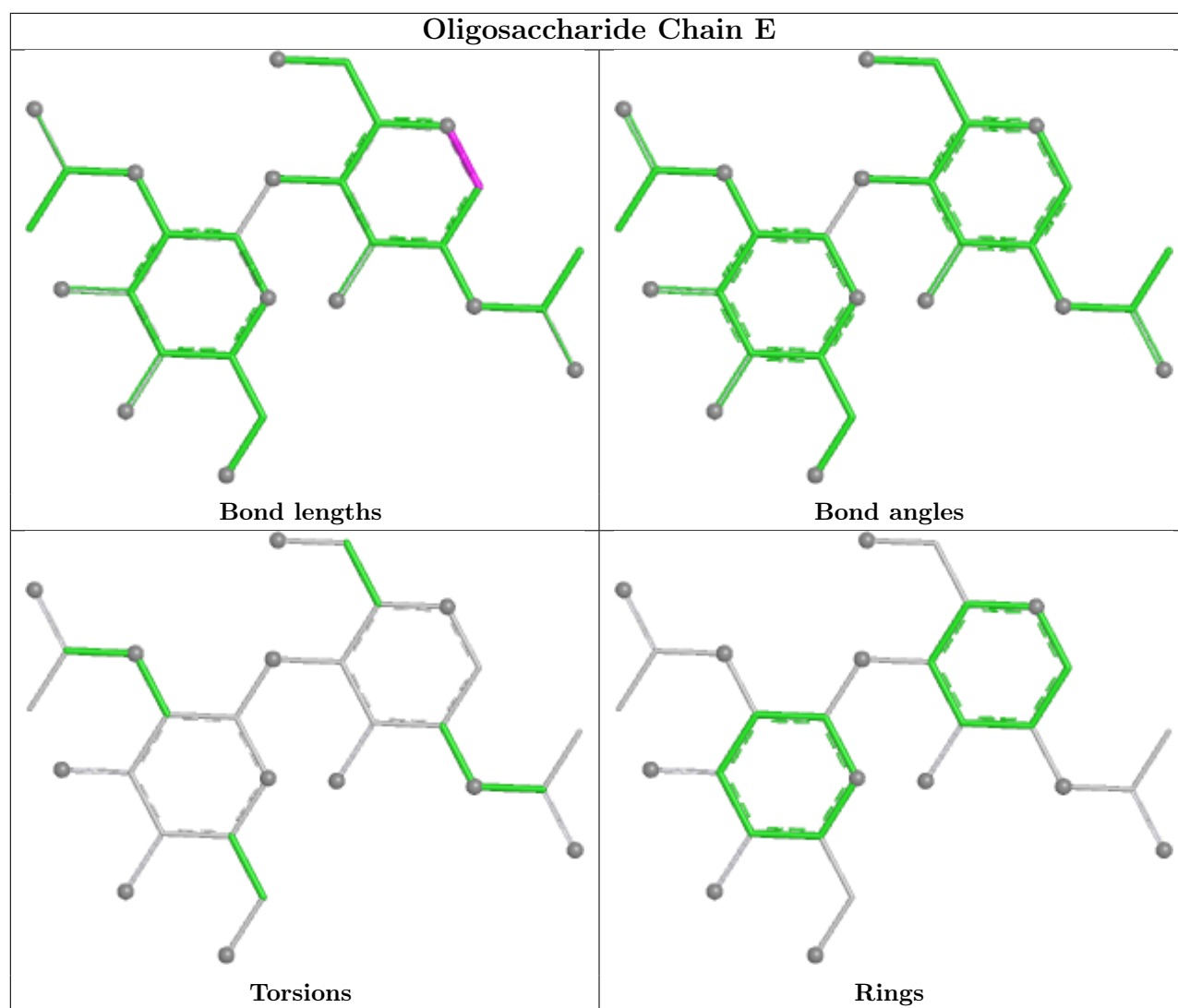
3 monomers are involved in 4 short contacts:

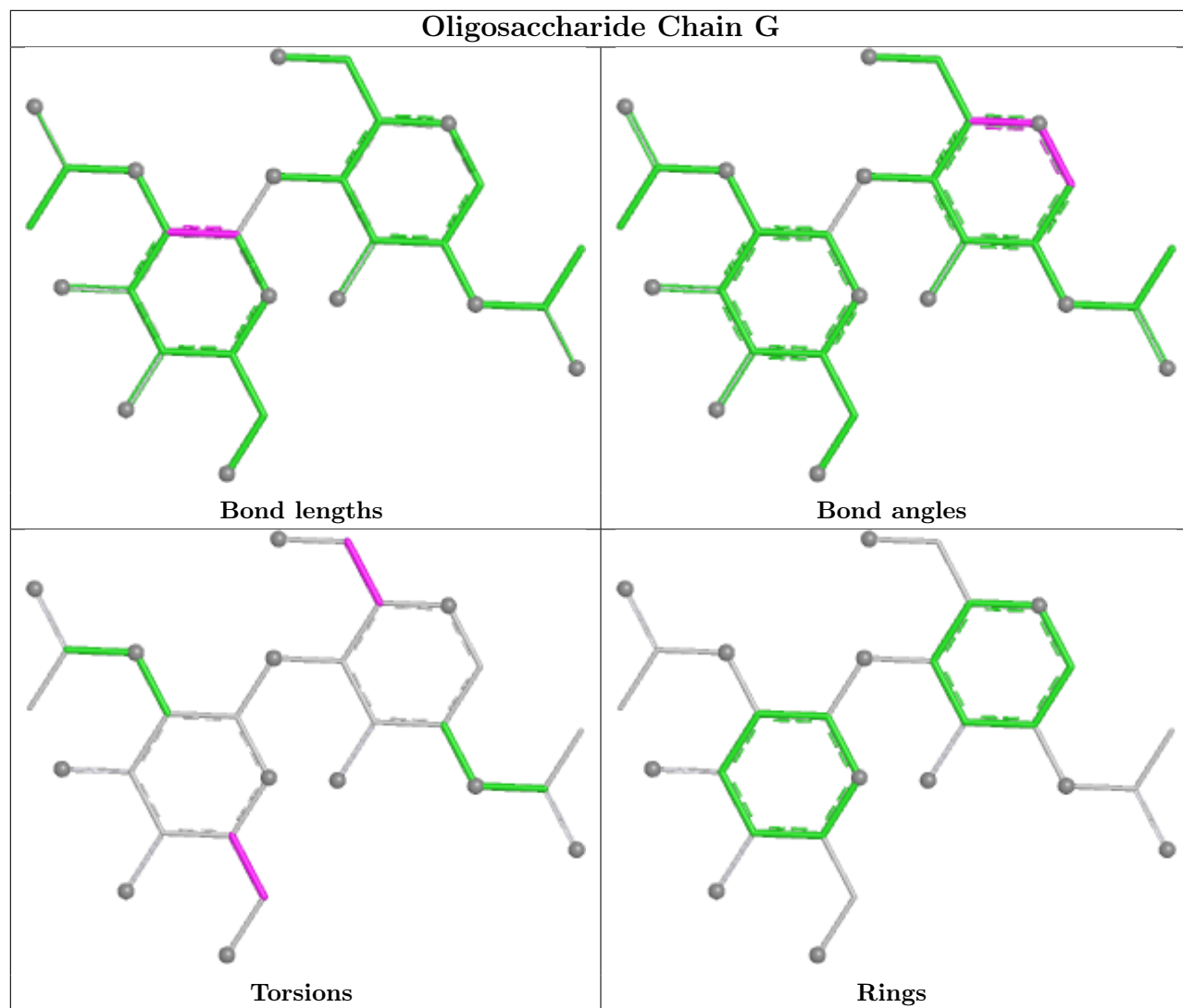
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	E	1	NAG	1	0
3	G	1	NAG	1	0
3	I	1	NAG	2	0

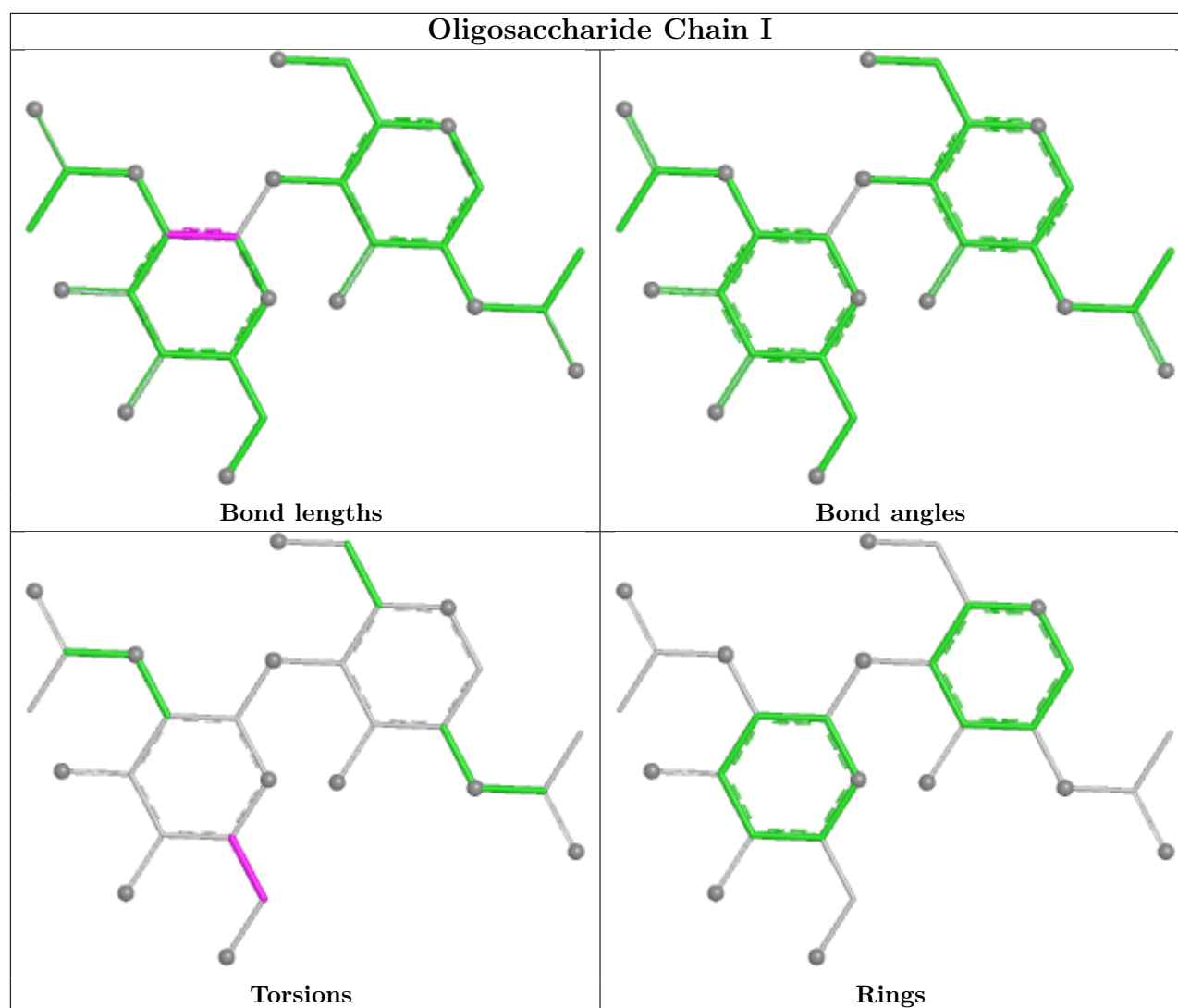
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.











5.6 Ligand geometry [i](#)

4 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	NAG	B	1001	1	14,14,15	0.58	0	17,19,21	0.59	0
4	NAG	B	1002	1	14,14,15	0.57	0	17,19,21	0.60	0
4	NAG	A	1001	1	14,14,15	0.39	0	17,19,21	0.50	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	NAG	A	1002	1	14,14,15	0.61	0	17,19,21	0.69	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	B	1001	1	-	1/6/23/26	0/1/1/1
4	NAG	B	1002	1	-	2/6/23/26	0/1/1/1
4	NAG	A	1001	1	-	0/6/23/26	0/1/1/1
4	NAG	A	1002	1	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	1001	NAG	O5-C5-C6-O6
4	B	1002	NAG	O5-C5-C6-O6
4	B	1002	NAG	C4-C5-C6-O6

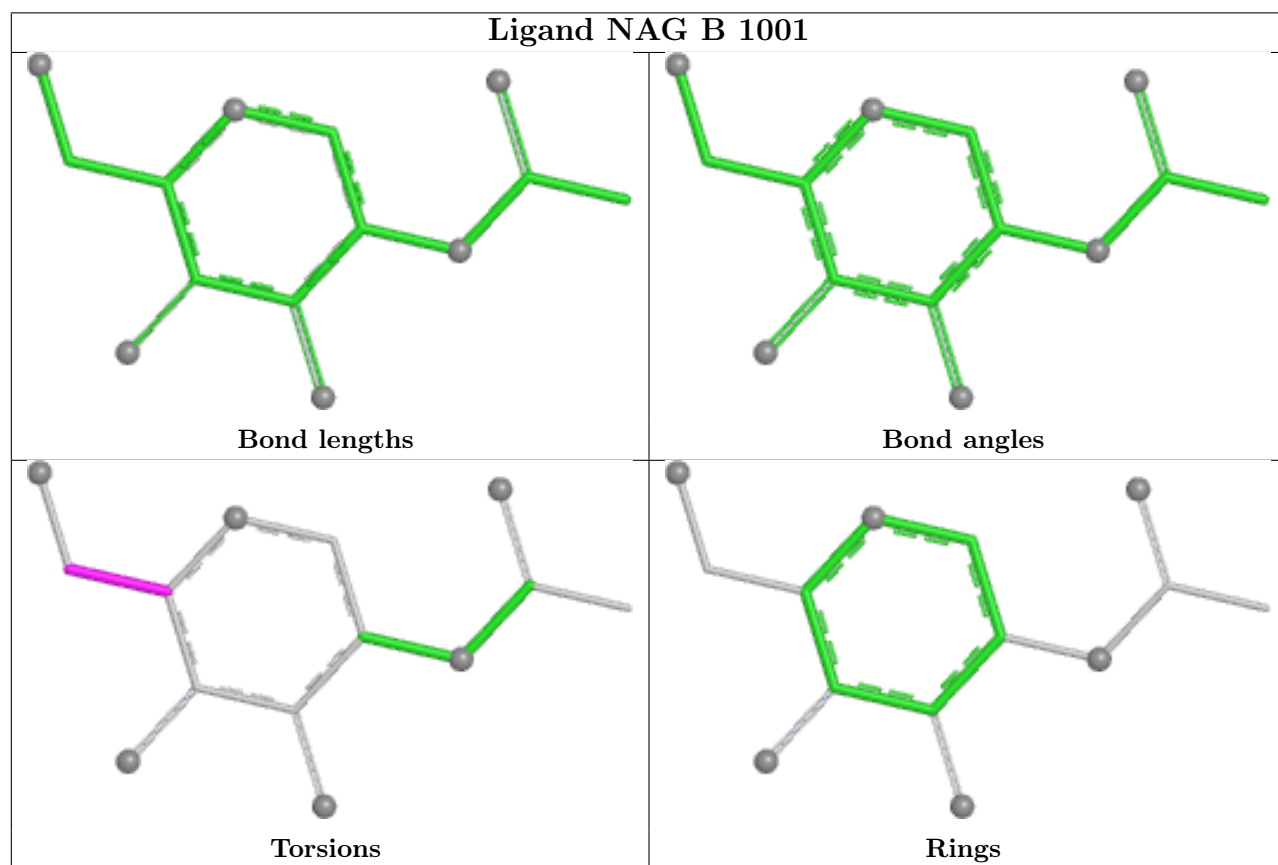
There are no ring outliers.

2 monomers are involved in 4 short contacts:

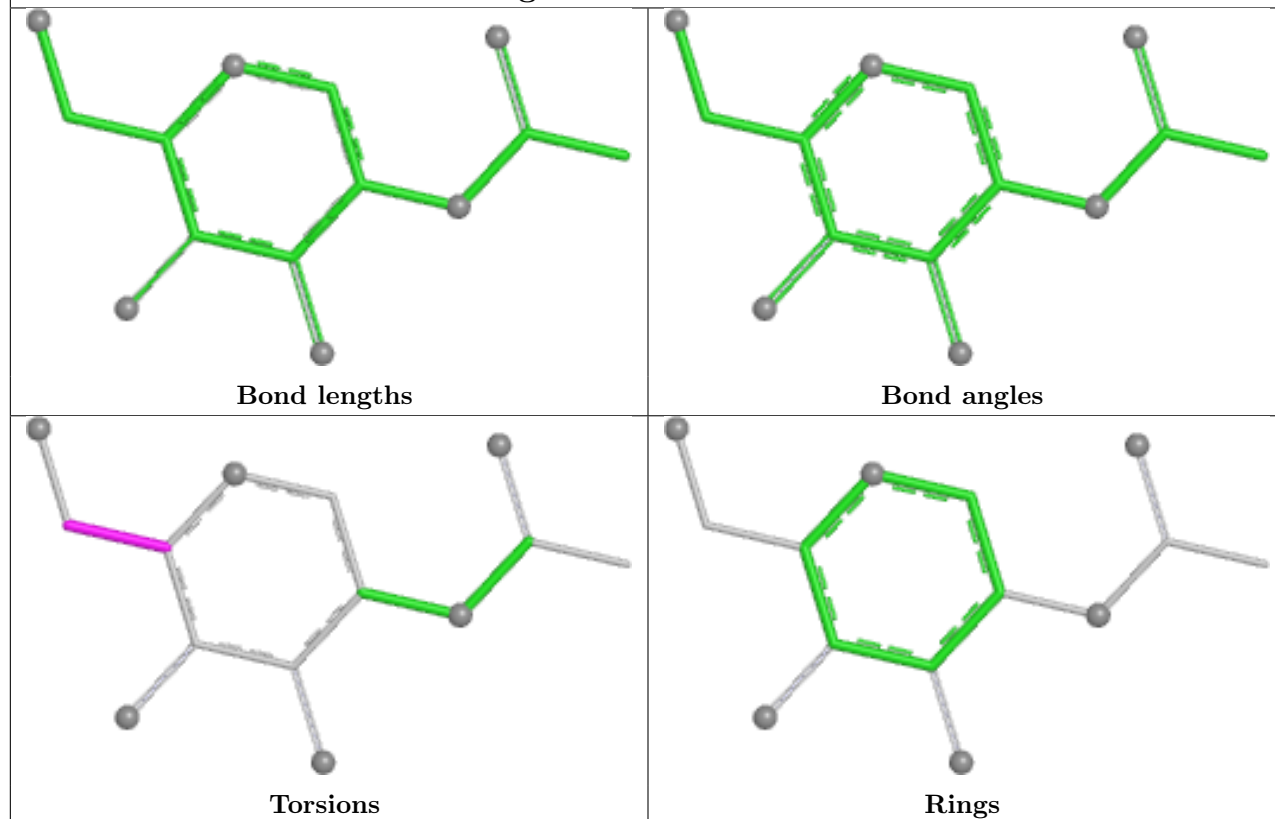
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	1001	NAG	2	0
4	A	1002	NAG	2	0

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

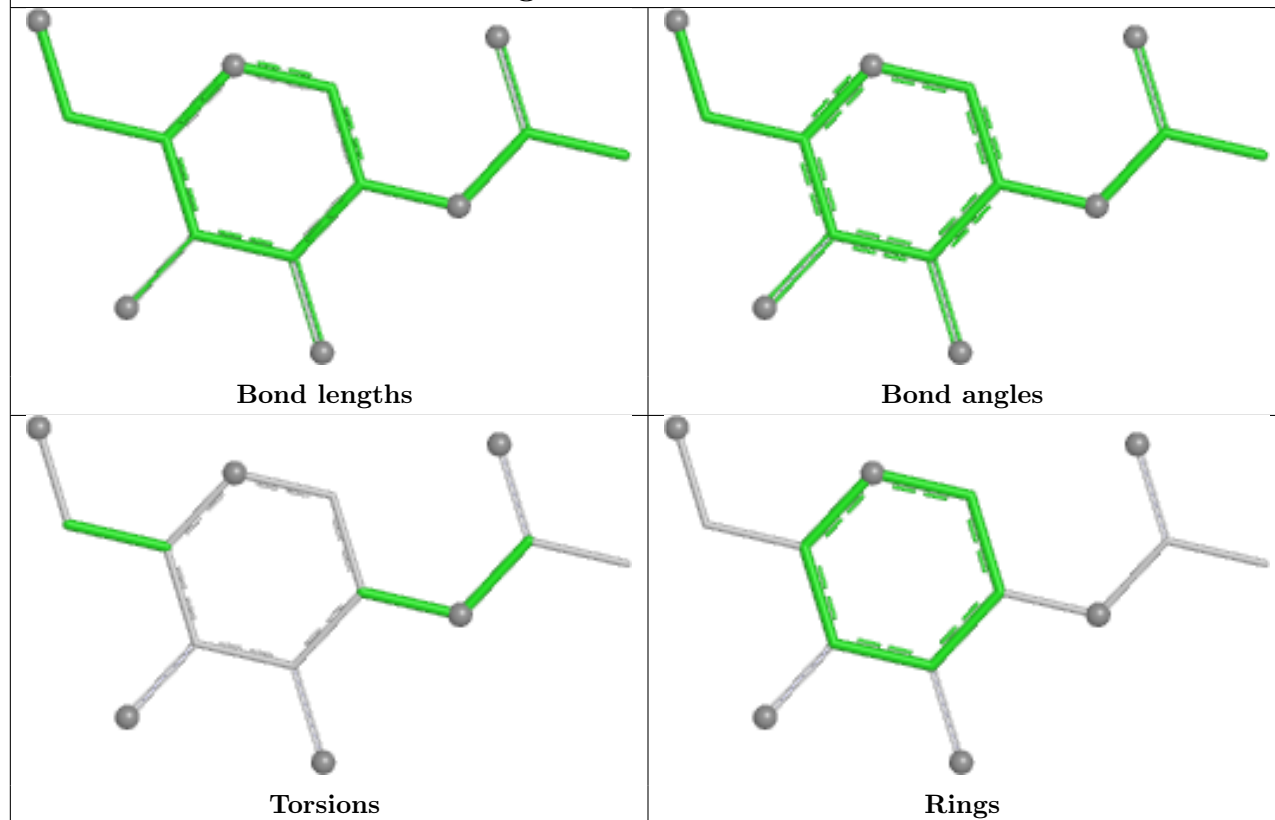
The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

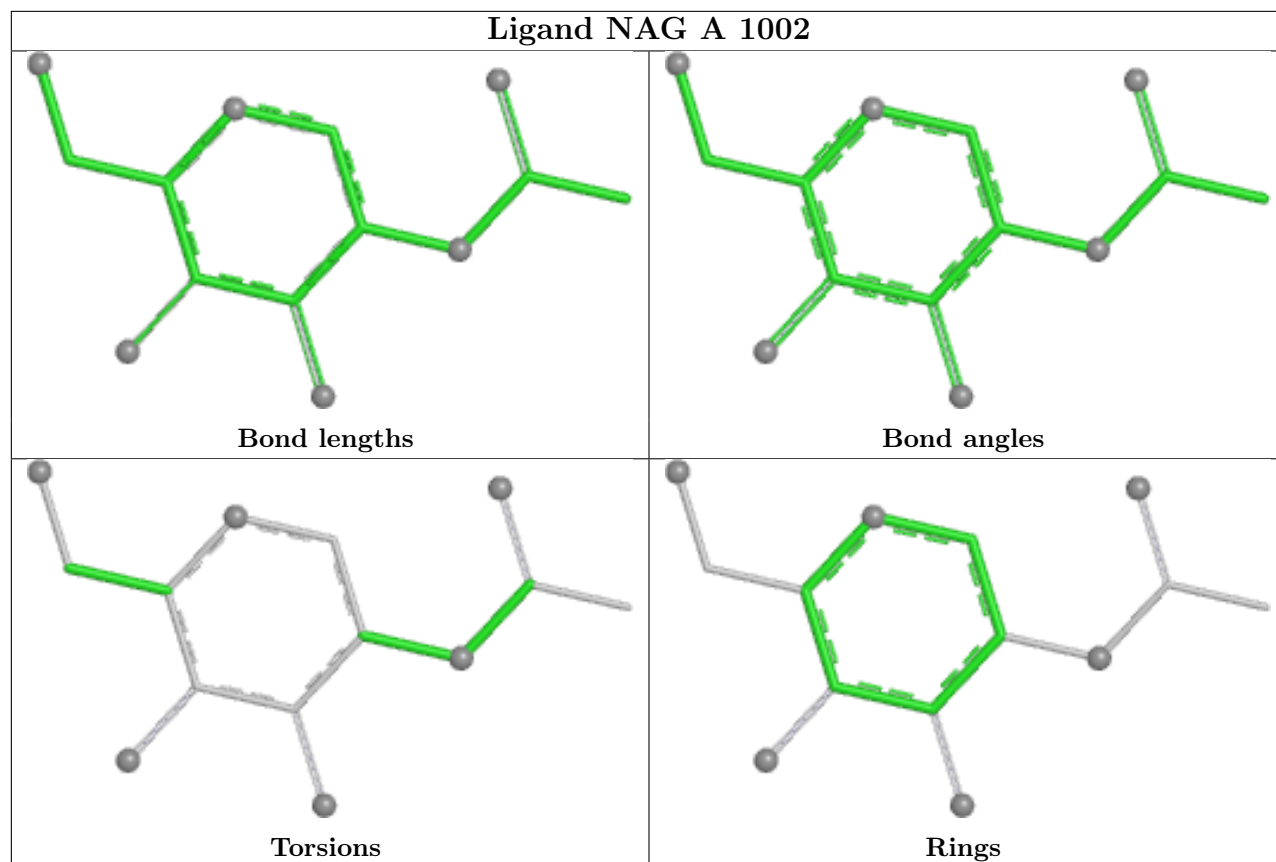


Ligand NAG B 1002



Ligand NAG A 1001





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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	670/694 (96%)	0.66	64 (9%)	13 17	146, 212, 261, 297	0
1	B	670/694 (96%)	0.56	47 (7%)	22 25	146, 207, 281, 324	0
2	C	525/552 (95%)	0.57	32 (6%)	27 28	130, 176, 234, 280	0
2	D	527/552 (95%)	0.53	37 (7%)	22 25	133, 180, 240, 263	0
All	All	2392/2492 (95%)	0.58	180 (7%)	20 23	130, 195, 265, 324	0

All (180) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	708	GLN	5.6
2	D	544	VAL	4.9
1	A	164	GLN	4.9
1	A	166	ALA	4.7
2	D	206	SER	4.6
1	A	165	GLU	4.5
2	C	544	VAL	4.5
1	A	500	TYR	4.4
2	D	46	GLY	4.3
2	C	484	ASP	4.3
1	A	167	GLY	4.2
1	B	192	ASP	4.2
1	A	131	GLY	4.1
1	A	488	ARG	4.1
1	A	112	ASP	4.1
2	D	554	GLU	4.1
2	D	484	ASP	4.0
1	A	469	GLY	3.9
2	C	554	GLU	3.8
1	B	500	TYR	3.8
2	D	468	GLU	3.7

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Mol	Chain	Res	Type	RSRZ
2	D	55	HIS	3.6
1	A	551	ASP	3.6
1	A	132	SER	3.6
1	B	581	PRO	3.6
1	A	58	GLU	3.6
1	A	537	LEU	3.5
1	A	111	THR	3.5
1	A	496	HIS	3.5
1	B	39	PRO	3.5
1	B	551	ASP	3.4
1	B	467	PRO	3.4
1	A	655	ASP	3.4
2	C	467	GLU	3.3
1	A	381	GLY	3.3
2	C	569	HIS	3.3
1	A	80	LEU	3.3
1	A	72	TYR	3.3
1	A	49	TRP	3.3
2	C	46	GLY	3.3
1	B	691	ALA	3.2
2	C	206	SER	3.2
2	C	478	SER	3.2
2	D	176	ALA	3.2
2	D	91	ILE	3.1
1	B	658	PHE	3.1
1	A	672	ASN	3.1
1	A	499	LEU	3.1
2	C	494	ARG	3.1
2	C	341	LYS	3.1
1	A	123	ALA	3.0
1	A	39	PRO	3.0
1	B	488	ARG	3.0
1	B	469	GLY	3.0
1	B	465	SER	3.0
1	B	470	ARG	3.0
1	B	167	GLY	3.0
1	A	116	LYS	3.0
1	A	238	LYS	3.0
1	A	684	HIS	2.9
1	B	708	GLN	2.9
2	D	495	ASP	2.9
2	C	55	HIS	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	605	GLU	2.9
1	B	553	PRO	2.8
2	C	164	ALA	2.8
1	B	196	GLU	2.8
2	D	189	ALA	2.8
2	D	569	HIS	2.8
1	A	169	MET	2.8
2	D	489	SER	2.8
1	A	52	THR	2.7
1	A	465	SER	2.7
2	D	204	ASP	2.7
1	A	494	PRO	2.7
1	B	672	ASN	2.7
1	B	489	ASP	2.6
2	D	426	VAL	2.6
2	D	298	GLN	2.6
1	B	213	ALA	2.6
1	A	441	ASP	2.6
1	A	656	PHE	2.6
1	B	154	HIS	2.6
1	B	208	ALA	2.6
1	A	592	LEU	2.6
1	A	168	SER	2.6
1	B	683	ARG	2.6
1	B	253	GLU	2.5
2	C	207	ALA	2.5
2	D	369	CYS	2.5
1	B	605	GLU	2.5
2	C	542	GLU	2.5
1	A	651	PHE	2.5
1	B	694	ALA	2.5
1	A	110	SER	2.5
1	B	319	ARG	2.5
2	C	534	GLY	2.5
2	D	129	ASN	2.5
2	D	478	SER	2.5
2	C	468	GLU	2.5
1	A	88	PRO	2.5
1	A	512	VAL	2.5
2	D	205	GLY	2.5
2	C	157	GLY	2.5
2	D	219	LYS	2.5

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Mol	Chain	Res	Type	RSRZ
2	C	500	TYR	2.4
1	B	241	ALA	2.4
1	A	70	ARG	2.4
1	B	479	VAL	2.4
1	A	501	ALA	2.4
2	D	192	LEU	2.4
2	D	307	ASN	2.4
1	A	581	PRO	2.4
1	B	600	PHE	2.4
1	A	519	THR	2.4
1	B	682	TYR	2.4
1	B	251	ARG	2.3
1	B	445	ARG	2.3
2	D	111	LYS	2.3
2	C	92	PRO	2.3
1	B	93	GLU	2.3
1	A	74	LEU	2.3
1	A	631	GLY	2.3
1	A	495	ASN	2.3
2	D	534	GLY	2.3
2	C	555	GLN	2.3
1	B	519	THR	2.3
2	D	490	LEU	2.3
1	A	130	CYS	2.3
1	B	417	GLY	2.3
1	A	557	ALA	2.3
1	A	59	GLN	2.2
2	D	130	ASP	2.2
2	D	236	PHE	2.2
2	C	245	ASN	2.2
1	A	103	SER	2.2
1	B	480	ALA	2.2
2	C	428	ARG	2.2
2	C	401	SER	2.2
1	B	164	GLN	2.2
1	A	489	ASP	2.2
1	A	270	PRO	2.2
2	D	62	MET	2.2
1	A	582	LEU	2.2
1	A	470	ARG	2.2
1	B	660	ASN	2.2
1	B	576	THR	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	695	PHE	2.2
2	D	225	HIS	2.2
1	B	131	GLY	2.2
2	C	246	ASN	2.2
1	A	658	PHE	2.2
1	A	458	ARG	2.2
2	D	128	ARG	2.2
1	A	50	GLY	2.1
1	B	116	LYS	2.1
1	B	474	ALA	2.1
1	B	411	PHE	2.1
1	A	161	SER	2.1
2	C	128	ARG	2.1
2	C	40	GLN	2.1
1	A	274	GLY	2.1
2	D	92	PRO	2.1
2	C	91	ILE	2.1
1	A	76	GLY	2.1
2	D	203	GLY	2.1
2	C	302	ASP	2.1
2	C	564	HIS	2.1
1	B	505	LYS	2.1
2	C	247	LEU	2.1
2	C	378	ALA	2.1
2	D	161	SER	2.1
2	D	522	LYS	2.1
1	A	457	ILE	2.1
2	C	465	LEU	2.0
1	A	150	GLY	2.0
2	D	75	ASP	2.0
2	D	457	PRO	2.0
1	B	155	ARG	2.0
1	B	655	ASP	2.0

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

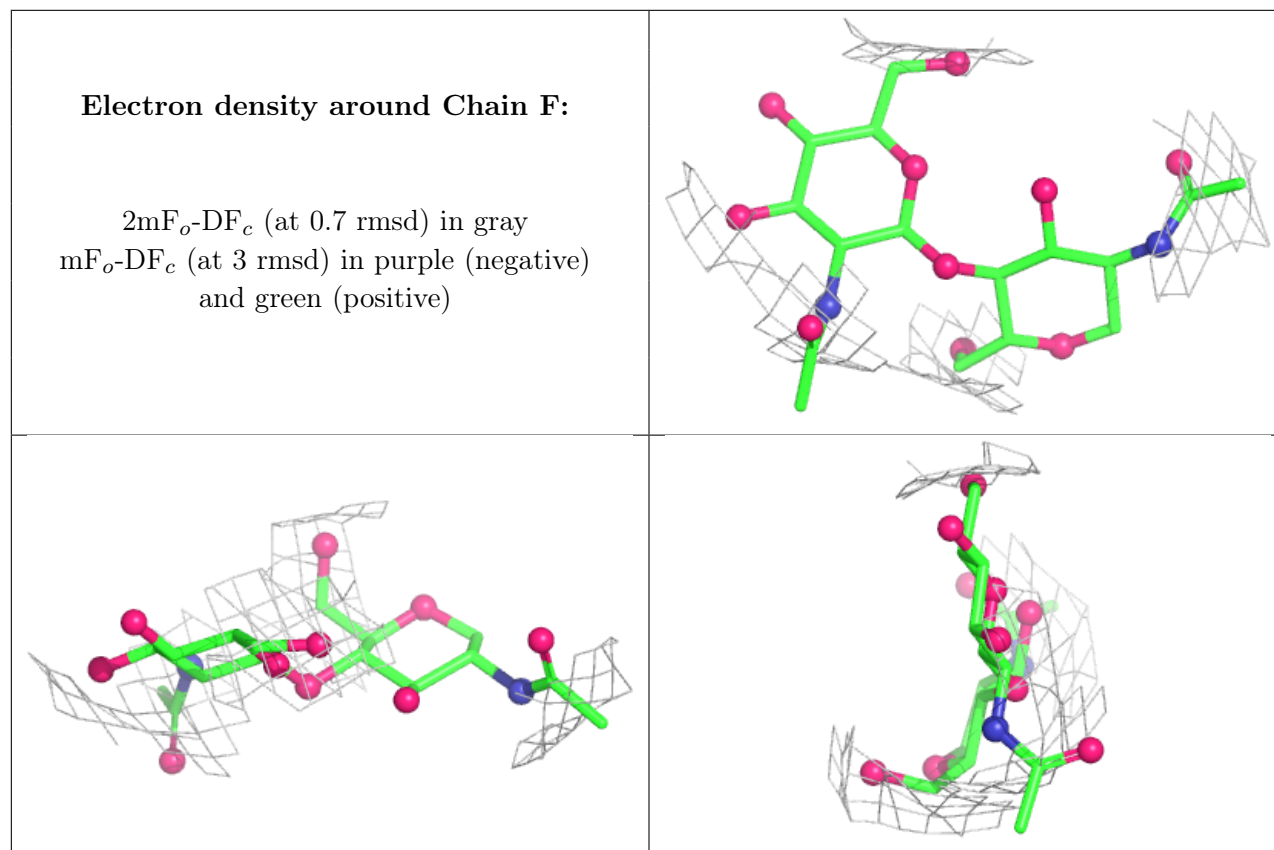
There are no non-standard protein/DNA/RNA residues in this entry.

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.

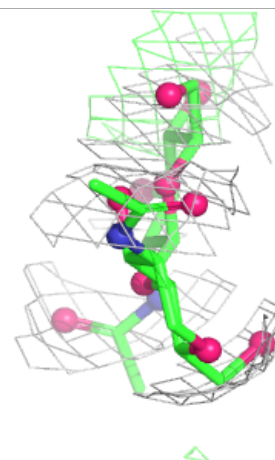
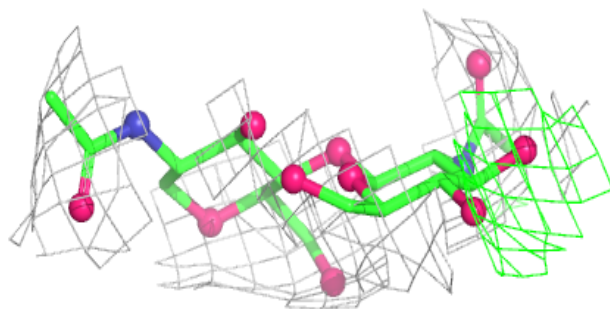
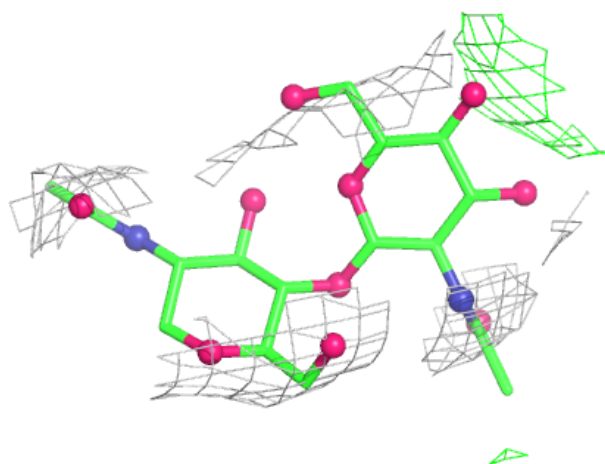
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	NAG	G	2	14/15	0.33	0.18	222,239,256,260	0
3	NAG	I	2	14/15	0.49	0.17	217,261,272,283	0
3	NAG	H	2	14/15	0.57	0.22	207,222,234,238	0
3	NAG	F	2	14/15	0.65	0.20	237,255,260,268	0
3	NAG	G	1	14/15	0.79	0.13	174,194,204,213	0
3	NAG	E	2	14/15	0.84	0.18	234,280,285,285	0
3	NAG	I	1	14/15	0.94	0.09	219,247,257,260	0
3	NAG	E	1	14/15	0.95	0.08	240,259,272,277	0
3	NAG	F	1	14/15	0.96	0.09	231,238,248,249	0
3	NAG	H	1	14/15	0.97	0.09	150,181,202,203	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.



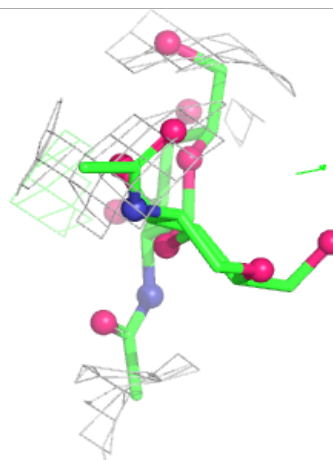
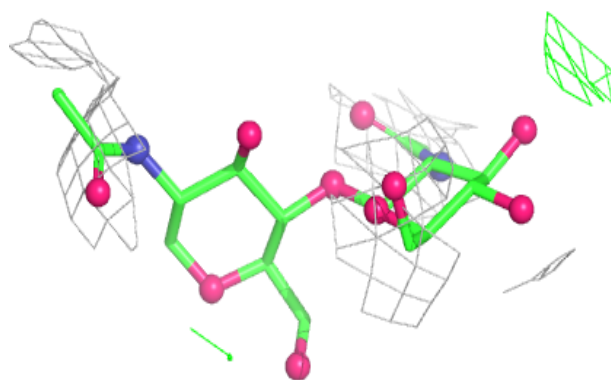
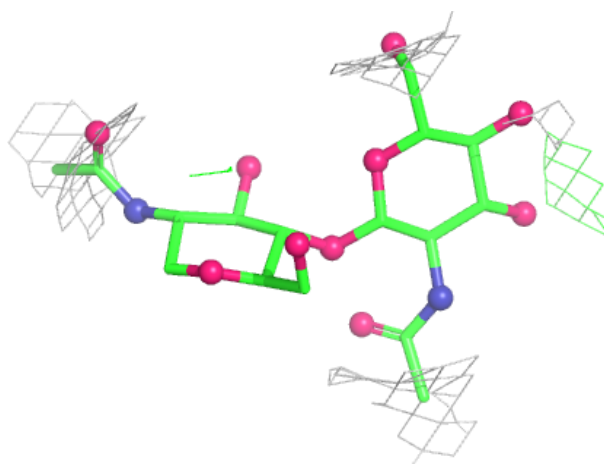
Electron density around Chain H:

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



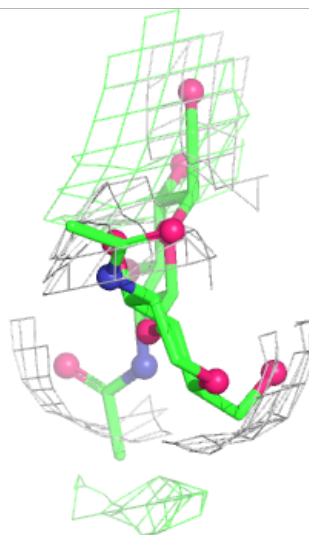
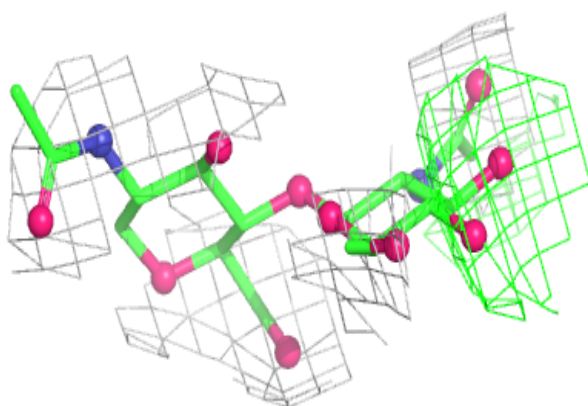
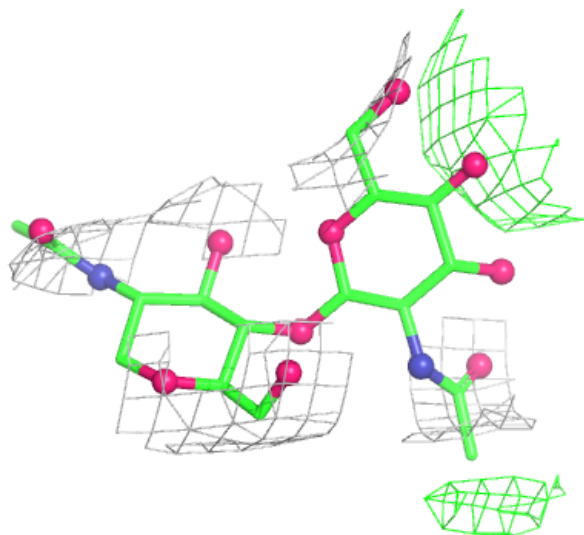
Electron density around Chain E:

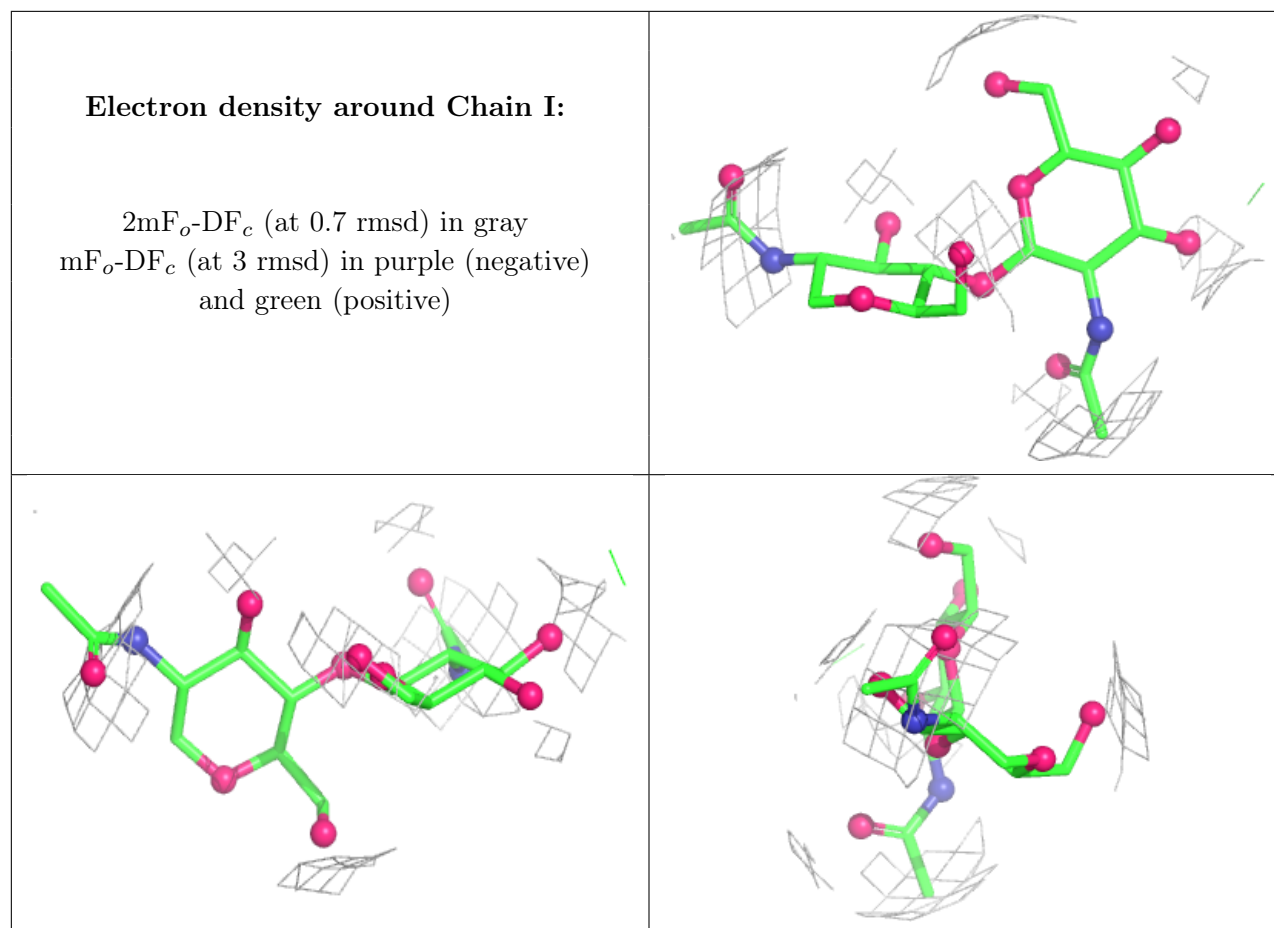
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain G:

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





6.4 Ligands [i](#)

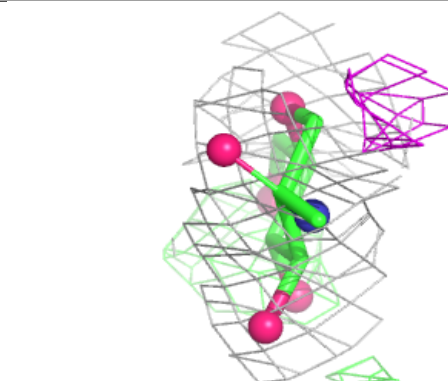
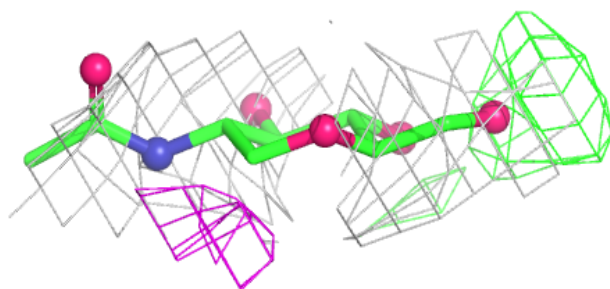
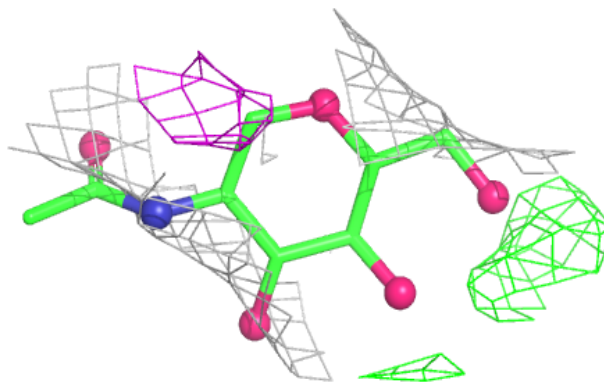
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	NAG	B	1001	14/15	0.59	0.25	224,253,263,265	0
4	NAG	A	1002	14/15	0.64	0.18	227,238,240,240	0
4	NAG	A	1001	14/15	0.74	0.21	254,269,272,273	0
4	NAG	B	1002	14/15	0.83	0.12	225,236,242,245	0

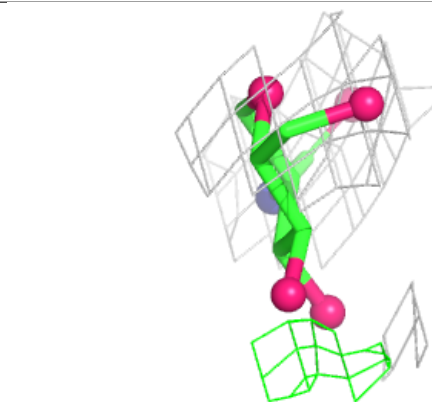
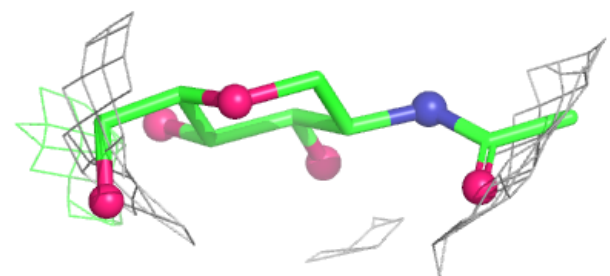
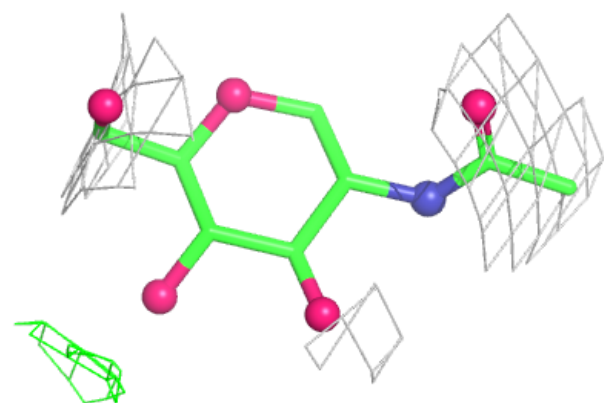
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around NAG B 1001:

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

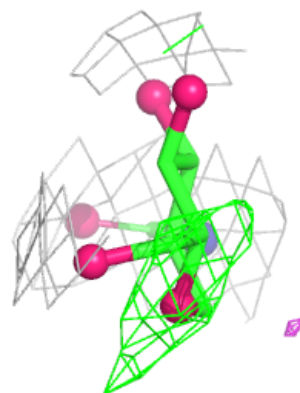
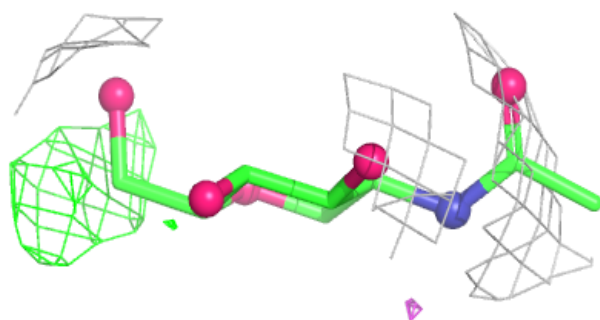
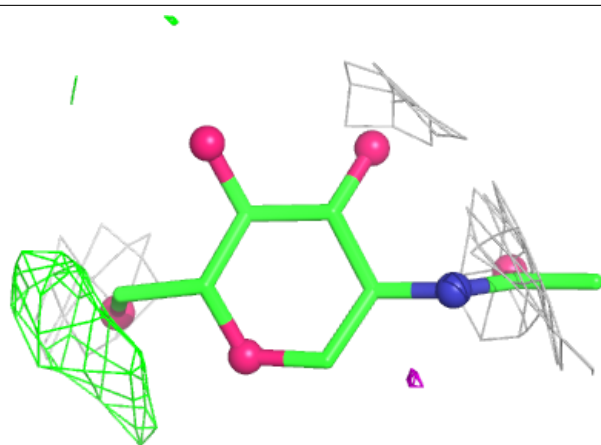
**Electron density around NAG A 1002:**

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

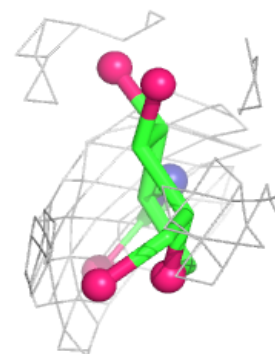
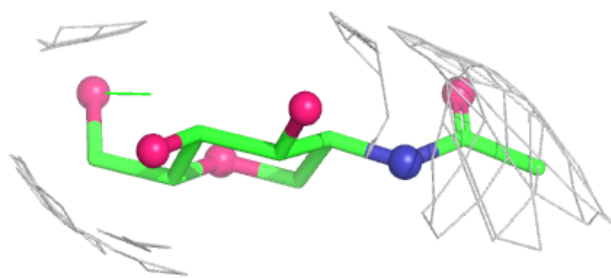
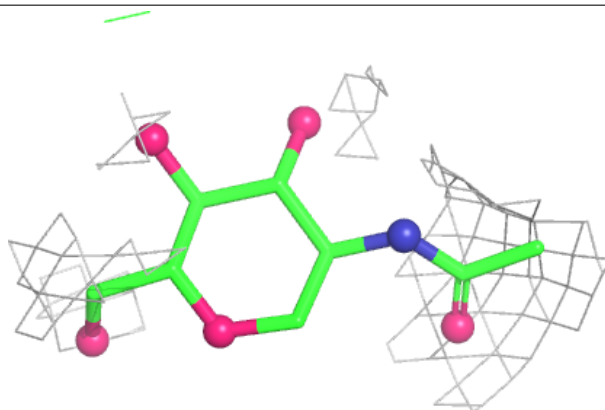


Electron density around NAG A 1001:

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

**Electron density around NAG B 1002:**

$2mF_o-DF_c$ (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.