



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:22 AM UTC

PDB ID : 8C5L / pdb_00008c5l
Title : NR2F6 ligand binding domain in complex with NSD1 peptide
Authors : Oerlemans, G.J.M.; van den Oetelaar, M.C.M.; Brunsveld, L.
Deposited on : 2023-01-09
Resolution : 2.60 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

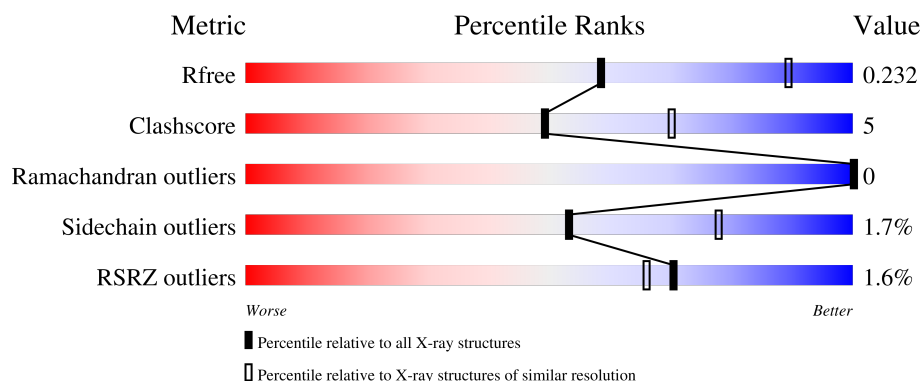
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

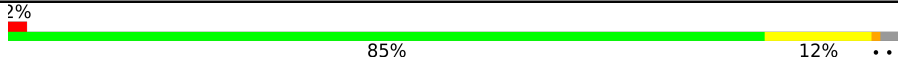
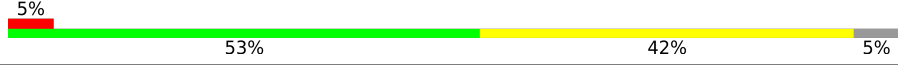
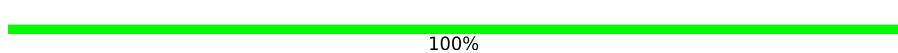
The reported resolution of this entry is 2.60 Å.

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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	568	
1	B	568	
2	C	19	
2	D	19	
3	F	2	

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Mol	Chain	Length	Quality of chain
3	G	2	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 8796 atoms, of which 44 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 Maltose/maltodextrin-binding periplasmic protein,Nuclear receptor subfamily 2 group F member 6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	556	Total	C	N	O	S	0	1	0
			4184	2707	681	784	12			
1	B	557	Total	C	N	O	S	0	0	0
			4177	2698	678	790	11			

There are 30 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	GLY	-	expression tag	UNP P0AEX9
A	83	ALA	ASP	engineered mutation	UNP P0AEX9
A	84	ALA	LYS	engineered mutation	UNP P0AEX9
A	173	ALA	GLU	engineered mutation	UNP P0AEX9
A	174	ALA	ASN	engineered mutation	UNP P0AEX9
A	240	ALA	LYS	engineered mutation	UNP P0AEX9
A	360	ALA	GLU	engineered mutation	UNP P0AEX9
A	363	ALA	LYS	engineered mutation	UNP P0AEX9
A	364	ALA	ASP	engineered mutation	UNP P0AEX9
A	368	ASN	-	linker	UNP P0AEX9
A	369	ALA	-	linker	UNP P0AEX9
A	370	ALA	-	linker	UNP P0AEX9
A	371	ALA	-	linker	UNP P0AEX9
A	372	GLU	-	linker	UNP P0AEX9
A	373	PHE	-	linker	UNP P0AEX9
B	1	GLY	-	expression tag	UNP P0AEX9
B	83	ALA	ASP	engineered mutation	UNP P0AEX9
B	84	ALA	LYS	engineered mutation	UNP P0AEX9
B	173	ALA	GLU	engineered mutation	UNP P0AEX9
B	174	ALA	ASN	engineered mutation	UNP P0AEX9
B	240	ALA	LYS	engineered mutation	UNP P0AEX9
B	360	ALA	GLU	engineered mutation	UNP P0AEX9
B	363	ALA	LYS	engineered mutation	UNP P0AEX9
B	364	ALA	ASP	engineered mutation	UNP P0AEX9

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Chain	Residue	Modelled	Actual	Comment	Reference
B	368	ASN	-	linker	UNP P0AEX9
B	369	ALA	-	linker	UNP P0AEX9
B	370	ALA	-	linker	UNP P0AEX9
B	371	ALA	-	linker	UNP P0AEX9
B	372	GLU	-	linker	UNP P0AEX9
B	373	PHE	-	linker	UNP P0AEX9

- Molecule 2 is a protein called Histone-lysine N-methyltransferase, H3 lysine-36 specific.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	18	Total 145	C 94	N 20	O 28	S 3	0	0	0
2	D	18	Total 144	C 93	N 20	O 28	S 3	0	0	0

- Molecule 3 is an oligosaccharide called alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	F	2	Total	C	H	O	0	0	0
			45	12	22	11			
3	G	2	Total	C	H	O	0	0	0
			45	12	22	11			

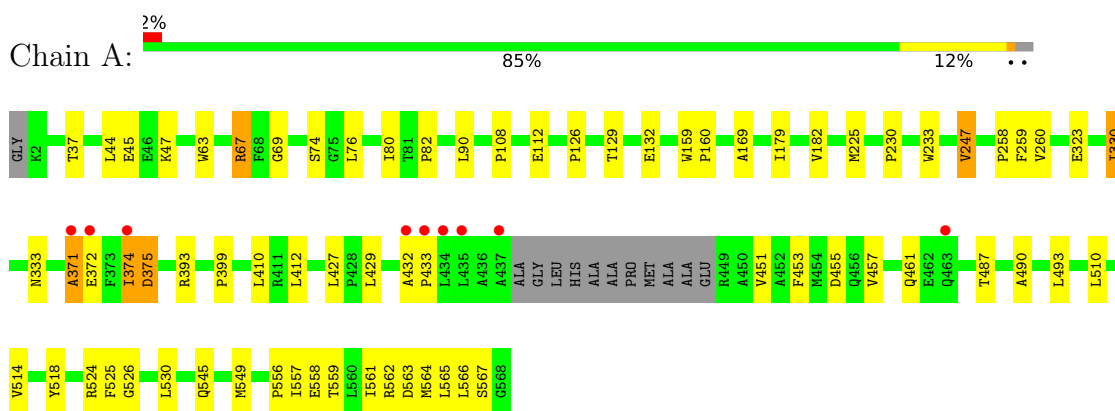
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	21	Total	O	0	0
			21	21		
4	B	35	Total	O	0	0
			35	35		

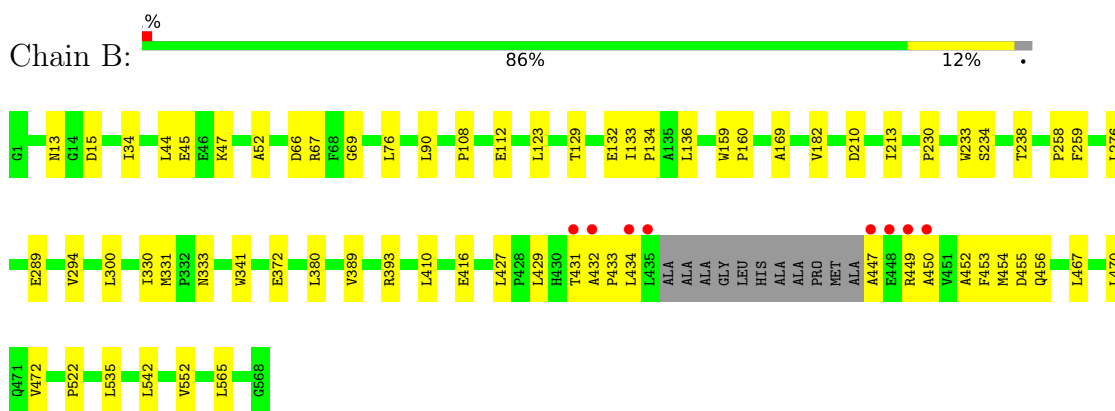
3 Residue-property plots

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 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: Maltose/maltodextrin-binding periplasmic protein,Nuclear receptor subfamily 2 group F member 6



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- Molecule 2: Histone-lysine N-methyltransferase, H3 lysine-36 specific



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- Molecule 3: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose



- Molecule 3: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose



4 Data and refinement statistics

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, α , β , γ	170.38Å 170.38Å 83.67Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	55.34 – 2.60 55.34 – 2.60	Depositor EDS
% Data completeness (in resolution range)	100.0 (55.34-2.60) 100.0 (55.34-2.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.36 (at 2.61Å)	Xtriage
Refinement program	REFMAC 5.8.0.352, PDB-REDO, PHENIX 1.19.1_4122	Depositor
R, R_{free}	0.200 , 0.233 0.203 , 0.232	Depositor DCC
R_{free} test set	2212 reflections (5.17%)	wwPDB-VP
Wilson B-factor (Å ²)	76.1	Xtriage
Anisotropy	0.014	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 39.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.027 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	8796	wwPDB-VP
Average B, all atoms (Å ²)	73.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.18% 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: GLC

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	0.40	1/4284 (0.0%)	0.60	1/5848 (0.0%)
1	B	0.36	1/4277 (0.0%)	0.56	1/5846 (0.0%)
2	C	0.57	0/147	0.87	0/195
2	D	0.44	0/146	0.87	0/194
All	All	0.38	2/8854 (0.0%)	0.59	2/12083 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	126	PRO	CA-C	7.10	1.56	1.52
1	B	522	PRO	C-N	-6.18	1.25	1.33

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	371	ALA	CA-C-O	-5.70	115.66	122.27
1	B	372	GLU	N-CA-C	-5.46	106.50	112.72

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4184	0	4058	45	0
1	B	4177	0	4017	38	0
2	C	145	0	134	6	0
2	D	144	0	132	6	0
3	F	23	22	21	0	0
3	G	23	22	21	1	0
4	A	21	0	0	2	0
4	B	35	0	0	0	0
All	All	8752	44	8383	93	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 5.

All (93) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:380:LEU:HD23	1:B:427:LEU:HD12	1.72	0.71
1:A:562:ARG:O	1:A:566:LEU:HG	1.90	0.70
1:A:90:LEU:HD23	1:A:108:PRO:HG2	1.77	0.67
1:A:510:LEU:HD22	1:A:525:PHE:HD1	1.58	0.67
1:B:45:GLU:HB2	1:B:67:ARG:HD2	1.77	0.67
1:A:371:ALA:O	1:A:375:ASP:N	2.28	0.66
1:A:557:ILE:HG22	1:A:561:ILE:HG13	1.77	0.65
2:C:2:TYR:O	2:C:6:THR:HG23	1.96	0.65
1:A:457:VAL:O	1:A:461:GLN:HG3	1.99	0.62
1:A:80:ILE:HG13	1:A:82:PRO:HD3	1.82	0.61
1:B:410:LEU:HD12	1:B:565:LEU:HD11	1.81	0.61
2:D:17:SER:O	2:D:18:LYS:C	2.44	0.61
1:A:545:GLN:HA	1:A:549:MET:HG2	1.84	0.60
1:B:210:ASP:OD1	1:B:213:ILE:HG12	2.02	0.60
1:B:447:ALA:HB3	1:B:450:ALA:HB2	1.84	0.58
1:B:52:ALA:HB3	1:B:76:LEU:HD13	1.85	0.57
1:B:169:ALA:O	1:B:182:VAL:HA	2.04	0.57
1:A:490:ALA:HB3	1:A:493:LEU:HD11	1.86	0.57
1:A:556:PRO:HG2	1:A:559:THR:HG23	1.87	0.56
1:B:123:LEU:HD11	1:B:136:LEU:HD21	1.86	0.56
2:D:16:ASP:O	2:D:18:LYS:HG3	2.06	0.55
1:B:389:VAL:O	1:B:393:ARG:HG3	2.07	0.55
2:D:6:THR:O	2:D:10:MET:HG3	2.07	0.55
1:B:433:PRO:HB3	1:B:454:MET:SD	2.49	0.52
1:A:74:SER:HB2	1:A:76:LEU:HG	1.92	0.52
1:B:452:ALA:HA	1:B:455:ASP:OD2	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:289:GLU:H	1:B:289:GLU:CD	2.18	0.51
1:B:416:GLU:HG2	1:B:535:LEU:HB3	1.91	0.51
1:B:552:VAL:HA	2:D:12:LYS:HG2	1.92	0.51
1:A:514:VAL:HG11	1:A:524:ARG:HB3	1.93	0.51
1:B:449:ARG:O	1:B:452:ALA:N	2.44	0.50
1:A:45:GLU:HB2	1:A:67:ARG:HD2	1.92	0.50
1:A:230:PRO:HA	1:A:233:TRP:CE2	2.47	0.50
1:B:66:ASP:HB3	1:B:331:MET:HE3	1.94	0.50
1:A:374:ILE:HG22	2:C:2:TYR:OH	2.13	0.49
2:D:9:MET:O	2:D:12:LYS:HB3	2.13	0.49
1:A:410:LEU:HD12	1:A:565:LEU:HD22	1.94	0.49
1:A:427:LEU:O	1:A:429:LEU:HD12	2.13	0.49
1:B:44:LEU:HA	1:B:47:LYS:HB2	1.95	0.49
1:B:13:ASN:ND2	1:B:15:ASP:OD1	2.43	0.48
1:B:453:PHE:HA	1:B:456:GLN:NE2	2.28	0.48
1:A:129:THR:OG1	1:A:132:GLU:HG3	2.14	0.48
1:A:179:ILE:HD12	1:A:374:ILE:HG13	1.95	0.48
1:B:230:PRO:HA	1:B:233:TRP:CE2	2.48	0.48
1:A:461:GLN:HB3	4:A:606:HOH:O	2.14	0.47
1:B:129:THR:OG1	1:B:132:GLU:HG3	2.13	0.47
2:C:16:ASP:C	2:C:18:LYS:N	2.72	0.47
1:B:160:PRO:HG3	1:B:258:PRO:HA	1.97	0.47
1:A:371:ALA:O	1:A:374:ILE:N	2.48	0.47
1:B:433:PRO:O	1:B:434:LEU:C	2.57	0.46
2:C:16:ASP:C	2:C:18:LYS:H	2.22	0.46
1:A:453:PHE:O	1:A:457:VAL:HG23	2.15	0.46
1:B:427:LEU:HD23	1:B:429:LEU:HD21	1.97	0.46
1:A:69:GLY:HA3	1:A:333:ASN:O	2.17	0.45
1:A:451:VAL:O	1:A:455:ASP:N	2.47	0.45
2:D:4:PHE:HA	2:D:7:LEU:HB2	1.99	0.45
1:A:563:ASP:O	1:A:567:SER:N	2.50	0.45
1:A:564:MET:HE2	1:A:564:MET:HB3	1.88	0.44
2:C:12:LYS:HG3	2:C:13:ASP:N	2.32	0.44
1:A:412:LEU:HD23	1:A:487:THR:HG21	1.99	0.44
1:B:34:ILE:HG21	1:B:276:LEU:HD13	1.99	0.44
1:B:431:THR:HG22	1:B:432:ALA:N	2.33	0.43
1:B:449:ARG:O	1:B:450:ALA:C	2.61	0.43
1:B:90:LEU:HD23	1:B:108:PRO:HG2	2.00	0.43
1:A:44:LEU:HD21	1:A:63:TRP:CZ3	2.54	0.43
1:A:371:ALA:O	1:A:372:GLU:C	2.61	0.43
1:A:518:TYR:OH	4:A:601:HOH:O	2.21	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:467:LEU:O	1:B:472:VAL:HG23	2.19	0.43
1:B:159:TRP:O	1:B:160:PRO:C	2.62	0.43
1:A:169:ALA:O	1:A:182:VAL:HA	2.19	0.42
1:B:69:GLY:HA3	1:B:333:ASN:O	2.18	0.42
1:B:294:VAL:HG11	1:B:300:LEU:HD21	2.02	0.42
1:A:557:ILE:O	1:A:558:GLU:C	2.62	0.42
1:B:234:SER:O	1:B:238:THR:HG23	2.20	0.42
1:A:393:ARG:HG3	1:A:565:LEU:CD1	2.49	0.42
1:B:341:TRP:CD1	3:G:2:GLC:H4	2.55	0.42
1:B:456:GLN:HB3	1:B:542:LEU:HD22	2.01	0.41
1:A:432:ALA:O	1:A:433:PRO:C	2.64	0.41
1:B:453:PHE:HA	1:B:456:GLN:HE21	1.85	0.41
2:C:8:LEU:HD12	2:C:8:LEU:HA	1.89	0.41
1:A:526:GLY:O	1:A:530:LEU:HG	2.20	0.41
1:B:431:THR:HG22	1:B:432:ALA:H	1.84	0.41
1:A:169:ALA:O	1:A:182:VAL:HG13	2.21	0.41
1:A:225:MET:HE3	1:A:225:MET:HB2	1.95	0.41
1:B:133:ILE:HB	1:B:134:PRO:HD3	2.03	0.41
1:A:159:TRP:O	1:A:160:PRO:C	2.63	0.41
1:A:247:VAL:HG22	1:A:323:GLU:OE1	2.21	0.41
1:A:410:LEU:HD23	1:A:410:LEU:HA	1.84	0.41
1:A:514:VAL:CG1	1:A:524:ARG:HB3	2.51	0.41
1:A:160:PRO:HG3	1:A:258:PRO:HA	2.02	0.41
1:A:44:LEU:HA	1:A:47:LYS:HB3	2.03	0.40
1:A:393:ARG:O	1:A:399:PRO:HG3	2.21	0.40
1:A:260:VAL:HB	1:A:330:ILE:HA	2.03	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	553/568 (97%)	537 (97%)	16 (3%)	0	100	100
1	B	553/568 (97%)	540 (98%)	13 (2%)	0	100	100
2	C	16/19 (84%)	15 (94%)	1 (6%)	0	100	100
2	D	16/19 (84%)	15 (94%)	1 (6%)	0	100	100
All	All	1138/1174 (97%)	1107 (97%)	31 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	408/444 (92%)	400 (98%)	8 (2%)	48	74
1	B	404/444 (91%)	400 (99%)	4 (1%)	68	86
2	C	15/19 (79%)	14 (93%)	1 (7%)	15	33
2	D	15/19 (79%)	14 (93%)	1 (7%)	15	33
All	All	842/926 (91%)	828 (98%)	14 (2%)	53	78

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

Mol	Chain	Res	Type
1	A	37	THR
1	A	67	ARG
1	A	112	GLU
1	A	247	VAL
1	A	259	PHE
1	A	330	ILE
1	A	374	ILE
1	A	375	ASP
1	B	112	GLU
1	B	259	PHE
1	B	330	ILE
1	B	470	LEU
2	C	7	LEU

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Mol	Chain	Res	Type
2	D	7	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	19	ASN
1	A	242	ASN
1	A	283	ASN
1	A	461	GLN
1	A	523	GLN
1	B	40	HIS
1	B	202	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

4 monosaccharides are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	GLC	F	1	3	12,12,12	0.26	0	17,17,17	0.46	0
3	GLC	F	2	3	11,11,12	0.48	0	15,15,17	0.48	0
3	GLC	G	1	3	12,12,12	0.26	0	17,17,17	0.46	0
3	GLC	G	2	3	11,11,12	0.50	0	15,15,17	0.38	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	GLC	F	1	3	-	0/2/22/22	0/1/1/1
3	GLC	F	2	3	-	2/2/19/22	0/1/1/1
3	GLC	G	1	3	-	0/2/22/22	0/1/1/1
3	GLC	G	2	3	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

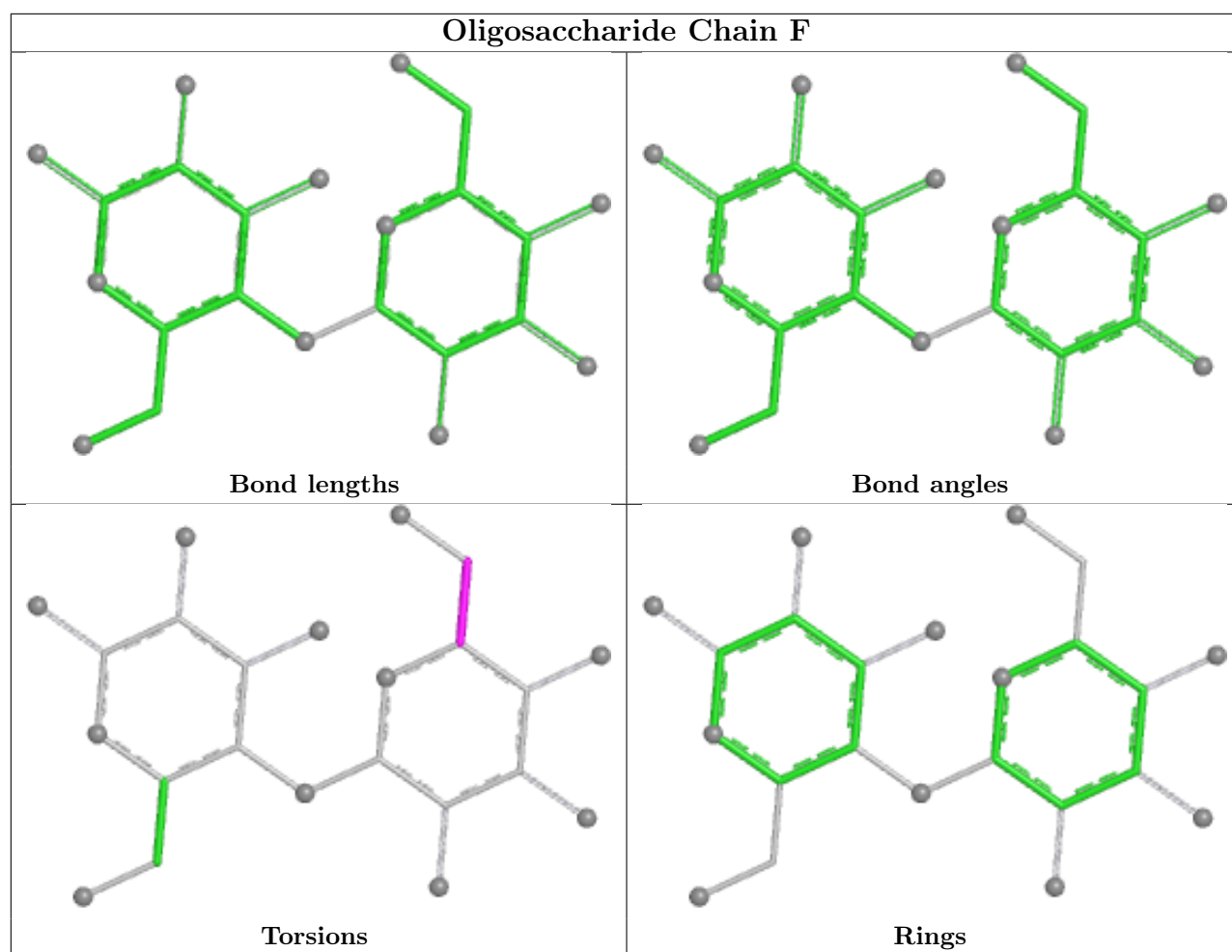
Mol	Chain	Res	Type	Atoms
3	F	2	GLC	C4-C5-C6-O6
3	F	2	GLC	O5-C5-C6-O6

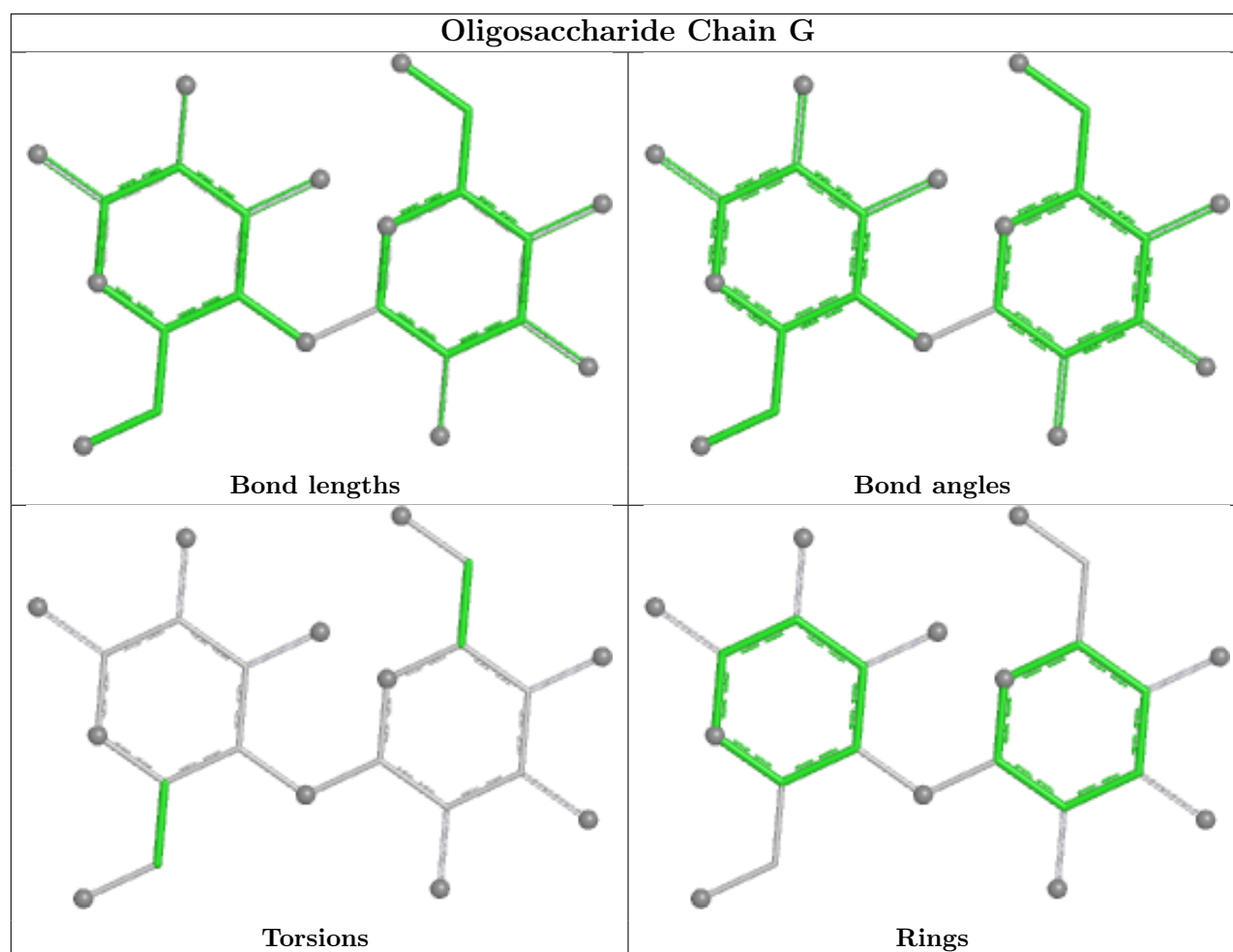
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	G	2	GLC	1	0

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





5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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	556/568 (97%)	-0.26	9 (1%) 70 66	33, 70, 91, 118	1 (0%)
1	B	557/568 (98%)	-0.27	8 (1%) 73 69	55, 72, 99, 134	0
2	C	18/19 (94%)	0.25	1 (5%) 30 24	72, 84, 105, 106	0
2	D	18/19 (94%)	0.17	0 100 100	75, 88, 110, 111	0
All	All	1149/1174 (97%)	-0.25	18 (1%) 70 66	33, 71, 97, 134	1 (0%)

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	374	ILE	4.2
1	A	433	PRO	3.8
1	A	432	ALA	3.6
1	A	437	ALA	3.4
1	A	372	GLU	3.3
1	A	371	ALA	3.3
1	B	447	ALA	3.2
1	A	435	LEU	3.1
1	B	450	ALA	2.9
1	B	448	GLU	2.8
1	B	431	THR	2.7
1	B	435	LEU	2.7
1	A	434	LEU	2.7
1	A	463	GLN	2.6
1	B	434	LEU	2.6
1	B	432	ALA	2.6
2	C	18	LYS	2.2
1	B	449	ARG	2.2

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

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

6.3 Carbohydrates [i](#)

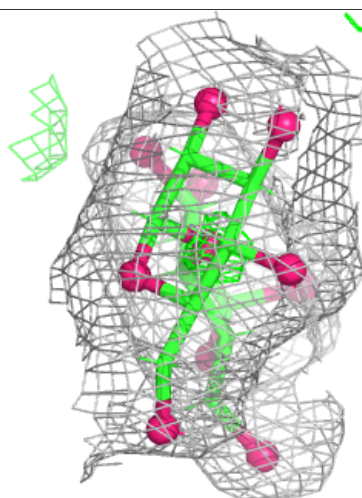
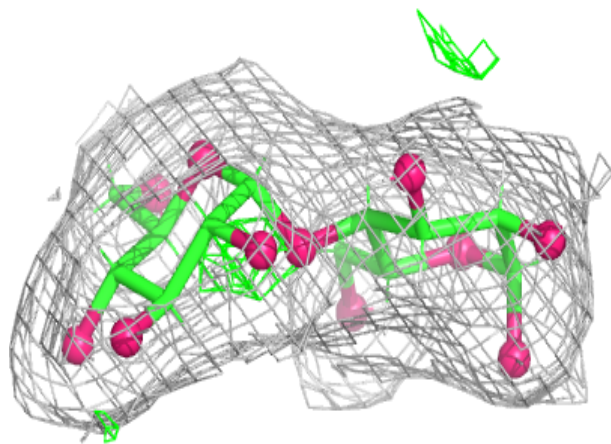
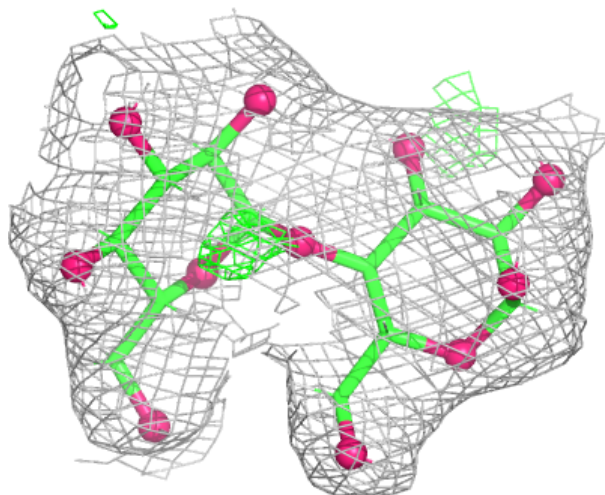
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

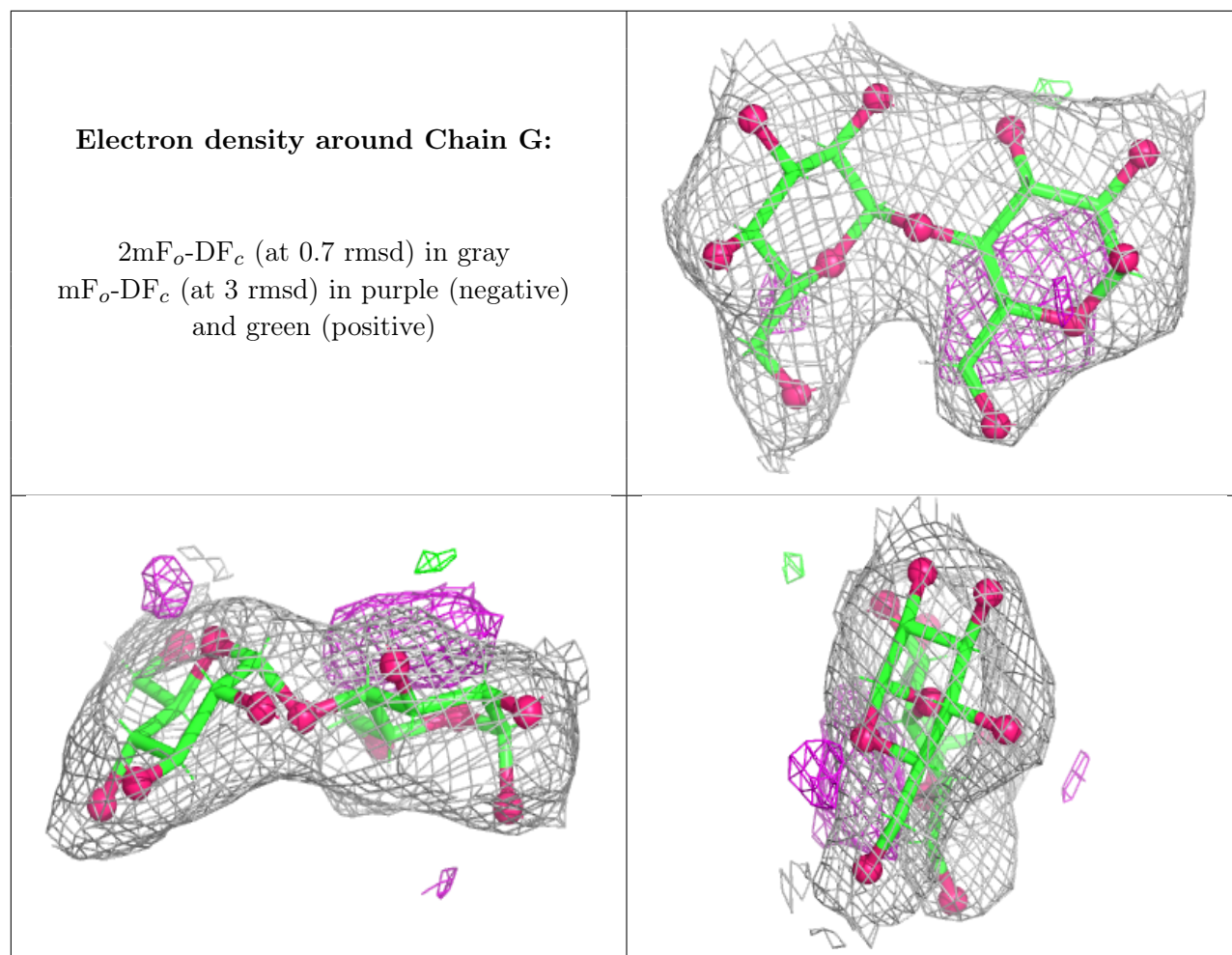
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	GLC	F	1	12/12	0.88	0.10	64,74,88,89	0
3	GLC	F	2	11/12	0.95	0.06	59,70,80,84	0
3	GLC	G	1	12/12	-	-	65,75,90,92	0
3	GLC	G	2	11/12	-	-	62,72,81,87	0

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 F:

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





6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.