



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 02:57 AM UTC

PDB ID : 1JCQ / pdb_00001jcq
Title : CRYSTAL STRUCTURE OF HUMAN PROTEIN FARNESYLTRANSFERASE COMPLEXED WITH FARNESYL DIPHOSPHATE AND THE PEPTIDOMIMETIC INHIBITOR L-739,750
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Deposited on : 2001-06-11
Resolution : 2.30 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

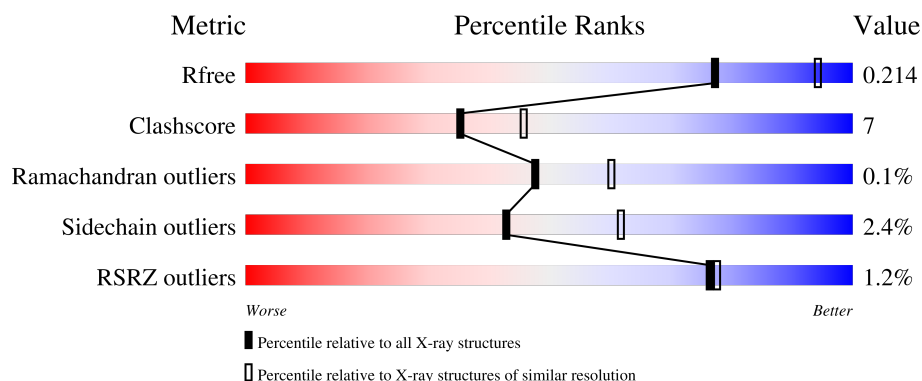
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	382	 66% 14% 18%
2	B	437	 75% 18% 6%
3	C	2	 100%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	ACY	A	3002	-	-	X	-

2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 6318 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 FARNESYLTRANSFERASE, ALPHA SUBUNIT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	313	Total	C	N	O	S	0	0	0
			2670	1704	465	496	5			

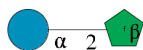
There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	380	GLU	-	insertion	UNP P49354
A	381	GLU	-	insertion	UNP P49354
A	382	PHE	-	insertion	UNP P49354

- Molecule 2 is a protein called PROTEIN FARNESYLTRANSFERASE, BETA SUBUNIT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	410	Total	C	N	O	S	0	0	0
			3228	2065	552	589	22			

- Molecule 3 is an oligosaccharide called beta-D-fructofuranose-(2-1)-alpha-D-glucopyranose.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
3	C	2	Total	C	O	0	0	0
			23	12	11			

- Molecule 4 is ACETIC ACID (CCD ID: ACY) (formula: C₂H₄O₂).

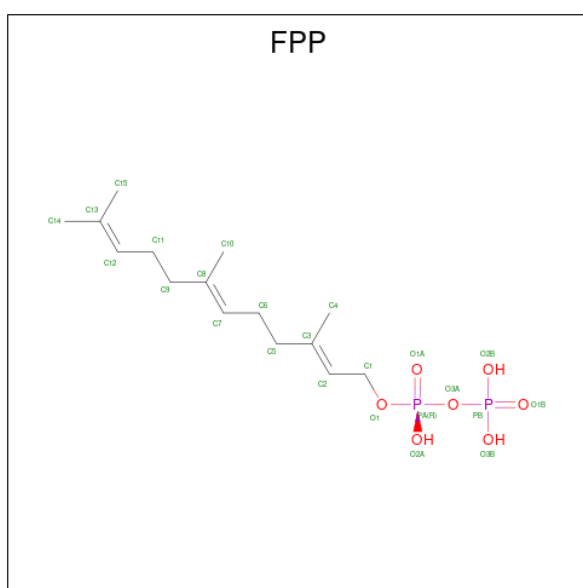


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			4	2	2		

- Molecule 5 is ZINC ION (CCD ID: ZN) (formula: Zn).

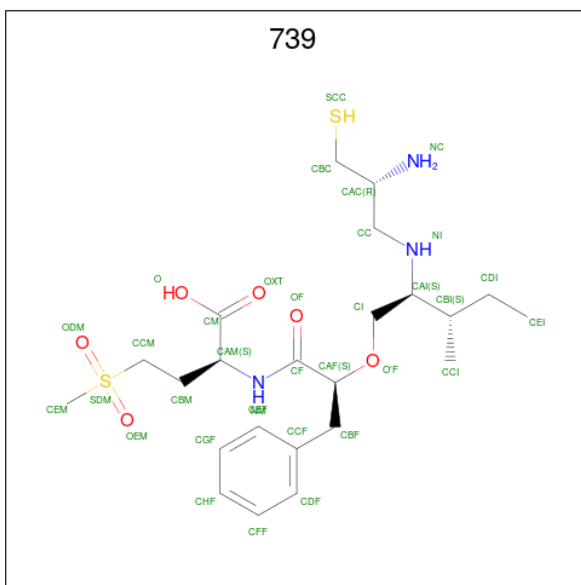
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	1	Total	Zn	0	0
			1	1		

- Molecule 6 is FARNESYL DIPHOSPHATE (CCD ID: FPP) (formula: C₁₅H₂₈O₇P₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	B	1	Total	C	O	P	0	0
			24	15	7	2		

- Molecule 7 is 2(S)-{2(S)-[2(R)-AMINO-3-MERCAPTO]PROPYLAMINO-3(S)-METHYL}PENTYLOXY-3-PHENYLPROPIONYLMETHIONINE SULFONE (CCD ID: 739) (formula: C₂₃H₃₉N₃O₆S₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
7	B	1	Total	C	N	O	S	0	0
			34	23	3	6	2		

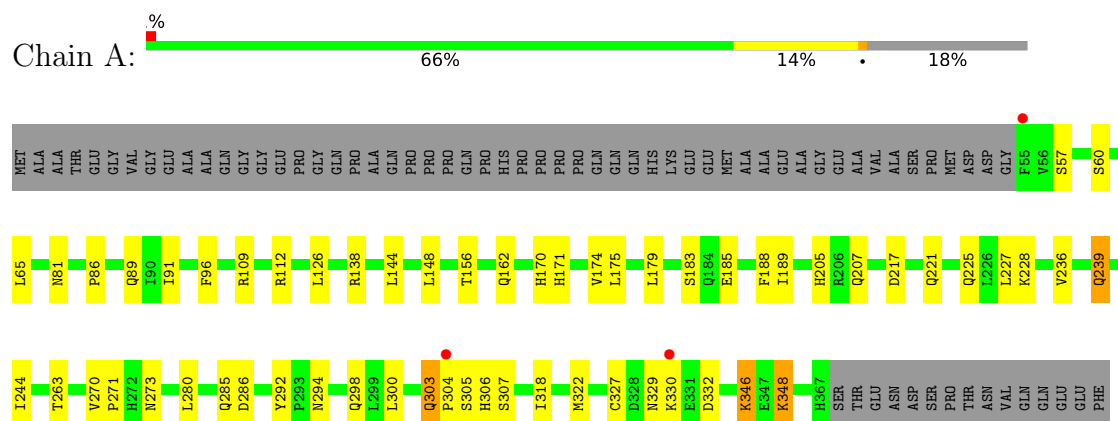
- Molecule 8 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	156	Total O 156 156	0	0
8	B	178	Total O 178 178	0	0

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: PROTEIN FARNESYLTRANSFERASE, ALPHA SUBUNIT



• Molecule 2: PROTEIN FARNESYLTRANSFERASE, BETA SUBUNIT



• Molecule 3: beta-D-fructofuranose-(2-1)-alpha-D-glucopyranose



4 Data and refinement statistics

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, α , β , γ	178.48 Å 178.48 Å 64.84 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	50.00 – 2.30 50.00 – 2.30	Depositor EDS
% Data completeness (in resolution range)	89.0 (50.00-2.30) 96.0 (50.00-2.30)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.76 (at 2.29 Å)	Xtriage
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.179 , 0.209 0.184 , 0.214	Depositor DCC
R_{free} test set	2534 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	38.7	Xtriage
Anisotropy	0.054	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 53.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.026 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6318	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.98% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, FRU, 739, ACY, FPP, 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	0.51	0/2737	0.91	10/3717 (0.3%)
2	B	0.55	0/3317	0.98	11/4508 (0.2%)
All	All	0.53	0/6054	0.95	21/8225 (0.3%)

There are no bond length outliers.

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	185	GLU	N-CA-C	8.10	120.91	111.02
1	A	348	LYS	N-CA-C	7.39	122.12	113.18
2	B	351	LEU	N-CA-C	6.55	118.10	108.86
2	B	174	ARG	N-CA-C	6.55	118.50	111.36
2	B	92	ALA	N-CA-C	-6.50	104.61	112.54
2	B	403	ILE	N-CA-C	-6.03	98.90	108.90
1	A	285	GLN	N-CA-C	6.02	117.84	111.28
2	B	375	HIS	N-CA-C	6.00	118.68	108.90
1	A	270	VAL	CA-C-N	5.96	126.39	119.47
1	A	270	VAL	C-N-CA	5.96	126.39	119.47
2	B	151	ALA	CA-C-N	-5.96	113.55	119.56
2	B	151	ALA	C-N-CA	-5.96	113.55	119.56
2	B	295	LEU	N-CA-C	5.91	119.27	110.52
2	B	164	GLY	N-CA-C	5.88	123.17	115.47
2	B	116	GLU	N-CA-C	-5.81	102.46	109.72
2	B	301	SER	N-CA-C	-5.51	105.38	111.71
1	A	217	ASP	N-CA-C	5.32	122.13	110.80
1	A	244	ILE	N-CA-C	5.30	115.51	110.42
1	A	292	TYR	CA-C-N	5.20	125.00	119.28
1	A	292	TYR	C-N-CA	5.20	125.00	119.28
1	A	183	SER	N-CA-C	5.03	118.19	111.75

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	2670	0	2589	41	0
2	B	3228	0	3147	48	0
3	C	23	0	21	0	0
4	A	4	0	3	6	0
5	B	1	0	0	0	0
6	B	24	0	25	1	0
7	B	34	0	37	2	0
8	A	156	0	0	7	0
8	B	178	0	0	5	0
All	All	6318	0	5822	87	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 7.

All (87) 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:303:GLN:O	1:A:307:SER:HB2	1.81	0.80
2:B:131:GLU:HG3	8:B:1471:HOH:O	1.82	0.80
1:A:81:ASN:HB3	8:A:1361:HOH:O	1.81	0.78
8:A:1213:HOH:O	2:B:275:GLN:HG2	1.84	0.76
2:B:152:PRO:HD3	7:B:3012:739:OEM	1.86	0.75
2:B:134:GLN:HE22	2:B:173:ASN:H	1.34	0.72
2:B:417:GLN:HG3	8:B:1223:HOH:O	1.89	0.71
1:A:300:LEU:O	1:A:303:GLN:HB2	1.91	0.71
1:A:170:HIS:HD2	4:A:3002:ACY:O	1.74	0.69
2:B:280:ARG:HE	2:B:289:GLN:HE21	1.41	0.68
2:B:74:GLN:H	2:B:344:GLN:HE22	1.45	0.63
2:B:135:SER:HB3	2:B:148:PRO:HG3	1.80	0.63
1:A:239:GLN:HE21	1:A:239:GLN:HA	1.64	0.62
1:A:346:LYS:NZ	1:A:346:LYS:HB2	2.14	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:135:SER:CB	2:B:148:PRO:HG3	2.30	0.62
1:A:189:ILE:HD11	1:A:205:HIS:HD2	1.64	0.61
2:B:301:SER:O	2:B:305:ALA:HB3	2.02	0.60
2:B:78:HIS:HD1	2:B:349:GLY:H	1.51	0.59
2:B:135:SER:OG	2:B:137:GLU:HG2	2.03	0.58
2:B:280:ARG:HE	2:B:289:GLN:NE2	2.00	0.58
1:A:170:HIS:HA	4:A:3002:ACY:CH3	2.34	0.57
2:B:253:PHE:HA	2:B:307:LEU:HD21	1.86	0.56
1:A:207:GLN:HE21	1:A:239:GLN:NE2	2.04	0.55
2:B:124:THR:HG22	2:B:167:GLU:OE1	2.06	0.55
2:B:312:HIS:ND1	8:B:1669:HOH:O	2.34	0.55
1:A:148:LEU:HB2	1:A:179:LEU:HD21	1.89	0.54
1:A:304:PRO:HG2	1:A:305:SER:H	1.73	0.54
2:B:239:ILE:HB	2:B:252:THR:HA	1.90	0.54
1:A:271:PRO:HG3	1:A:306:HIS:HD2	1.73	0.53
2:B:377:GLY:HA2	2:B:381:MET:O	2.09	0.53
2:B:77:LYS:HD3	2:B:346:PRO:O	2.09	0.53
2:B:86:LEU:HB2	2:B:107:ILE:HG21	1.89	0.52
1:A:294:ASN:O	1:A:298:GLN:HG3	2.09	0.52
1:A:170:HIS:CD2	4:A:3002:ACY:O	2.59	0.52
2:B:325:MET:HG2	8:B:1495:HOH:O	2.11	0.51
1:A:318:ILE:O	1:A:322:MET:HG3	2.11	0.50
2:B:338:TYR:CE2	2:B:343:CYS:SG	3.04	0.50
1:A:57:SER:HB3	1:A:60:SER:OG	2.11	0.50
1:A:329:ASN:HB3	1:A:332:ASP:HB3	1.93	0.50
1:A:346:LYS:HB2	1:A:346:LYS:HZ2	1.75	0.50
2:B:78:HIS:HD1	2:B:349:GLY:N	2.09	0.50
1:A:170:HIS:HA	4:A:3002:ACY:H3	1.93	0.50
2:B:93:TYR:CE2	2:B:96:LEU:HD12	2.48	0.49
2:B:93:TYR:CD2	2:B:96:LEU:HD12	2.48	0.49
2:B:148:PRO:HB2	2:B:180:TYR:CZ	2.47	0.49
2:B:306:GLY:O	2:B:309:PRO:HD2	2.12	0.48
1:A:239:GLN:HE21	1:A:239:GLN:CA	2.27	0.48
1:A:138:ARG:HD2	1:A:174:VAL:HG11	1.94	0.48
2:B:308:LEU:HD13	2:B:330:PHE:HB3	1.95	0.47
1:A:239:GLN:HA	1:A:239:GLN:NE2	2.29	0.47
2:B:423:PHE:CE2	2:B:424:GLU:HG2	2.49	0.47
1:A:91:ILE:O	2:B:91:ASP:HB3	2.14	0.46
2:B:175:GLU:HA	2:B:421:PRO:HB3	1.98	0.46
2:B:332:GLN:O	2:B:336:GLN:HG3	2.16	0.46
1:A:170:HIS:HA	4:A:3002:ACY:H2	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:271:PRO:HD2	2:B:38:ASP:CG	2.42	0.45
2:B:350:LEU:HB2	2:B:363:THR:HA	1.98	0.45
1:A:221:GLN:HG3	8:A:1180:HOH:O	2.16	0.45
2:B:148:PRO:HB2	2:B:180:TYR:CE1	2.52	0.45
1:A:170:HIS:HE1	2:B:196:GLY:O	1.99	0.45
1:A:86:PRO:HG2	1:A:89:GLN:NE2	2.32	0.44
1:A:170:HIS:CA	4:A:3002:ACY:H2	2.48	0.44
2:B:238:GLY:HA3	2:B:247:ALA:HB1	1.99	0.44
1:A:286:ASP:HB2	8:A:1601:HOH:O	2.17	0.44
2:B:134:GLN:HB2	2:B:140:PHE:CE2	2.53	0.44
1:A:273:ASN:HB3	8:A:1205:HOH:O	2.17	0.43
2:B:370:LEU:HD23	2:B:394:LEU:HD11	2.00	0.43
1:A:227:LEU:HG	1:A:236:VAL:HG11	2.00	0.43
2:B:99:SER:O	2:B:102:TRP:HB2	2.19	0.43
2:B:202:ARG:HD2	6:B:3011:FPP:H142	2.01	0.43
1:A:170:HIS:CE1	8:A:1345:HOH:O	2.72	0.42
2:B:345:CYS:HB3	2:B:348:GLY:O	2.20	0.42
1:A:227:LEU:HG	1:A:236:VAL:CG1	2.49	0.42
1:A:96:PHE:HA	1:A:126:LEU:HD13	2.01	0.42
1:A:156:THR:HG23	1:A:188:PHE:CD1	2.55	0.42
1:A:171:HIS:O	1:A:175:LEU:HG	2.21	0.41
2:B:192:LEU:HD23	2:B:199:VAL:CG2	2.50	0.41
2:B:211:ALA:HA	2:B:216:ILE:HG12	2.01	0.41
2:B:260:VAL:HG22	2:B:265:GLU:HG2	2.02	0.41
1:A:327:CYS:O	1:A:330:LYS:HG3	2.21	0.41
1:A:225:GLN:HG2	8:A:1286:HOH:O	2.21	0.41
2:B:271:LYS:HD3	8:B:1574:HOH:O	2.20	0.40
2:B:140:PHE:HB2	2:B:153:THR:HA	2.03	0.40
1:A:263:THR:HG21	1:A:280:LEU:HB2	2.02	0.40
1:A:112:ARG:O	1:A:144:LEU:HD21	2.22	0.40
2:B:378:SER:O	2:B:381:MET:HB2	2.21	0.40
7:B:3012:739:HDI2	7:B:3012:739:O'F	2.21	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	311/382 (81%)	294 (94%)	17 (6%)	0	100	100
2	B	408/437 (93%)	395 (97%)	12 (3%)	1 (0%)	43	55
All	All	719/819 (88%)	689 (96%)	29 (4%)	1 (0%)	48	60

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	74	GLN

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	292/344 (85%)	284 (97%)	8 (3%)	39	58
2	B	346/370 (94%)	339 (98%)	7 (2%)	48	67
All	All	638/714 (89%)	623 (98%)	15 (2%)	43	62

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	65	LEU
1	A	109	ARG
1	A	162	GLN
1	A	228	LYS
1	A	239	GLN
1	A	303	GLN
1	A	346	LYS
1	A	348	LYS
2	B	15	SER
2	B	75	ARG
2	B	87	ARG

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Mol	Chain	Res	Type
2	B	90	THR
2	B	99	SER
2	B	223	GLU
2	B	351	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	89	GLN
1	A	135	HIS
1	A	162	GLN
1	A	170	HIS
1	A	218	ASN
1	A	221	GLN
1	A	239	GLN
1	A	261	GLN
1	A	303	GLN
1	A	306	HIS
1	A	367	HIS
2	B	30	HIS
2	B	56	GLN
2	B	74	GLN
2	B	134	GLN
2	B	289	GLN
2	B	344	GLN
2	B	375	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

2 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	C	1	3	11,11,12	2.55	1 (9%)	15,15,17	1.51	3 (20%)
3	FRU	C	2	3	11,12,12	1.58	2 (18%)	10,18,18	1.04	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	C	1	3	-	0/2/19/22	0/1/1/1
3	FRU	C	2	3	-	0/5/24/24	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	1	GLC	C2-C3	7.78	1.64	1.52
3	C	2	FRU	O2-C2	3.46	1.46	1.40
3	C	2	FRU	C1-C2	2.77	1.58	1.52

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	1	GLC	C1-O5-C5	3.69	117.13	112.19
3	C	1	GLC	C1-C2-C3	-2.83	105.52	109.64
3	C	1	GLC	O3-C3-C2	-2.18	105.60	110.05

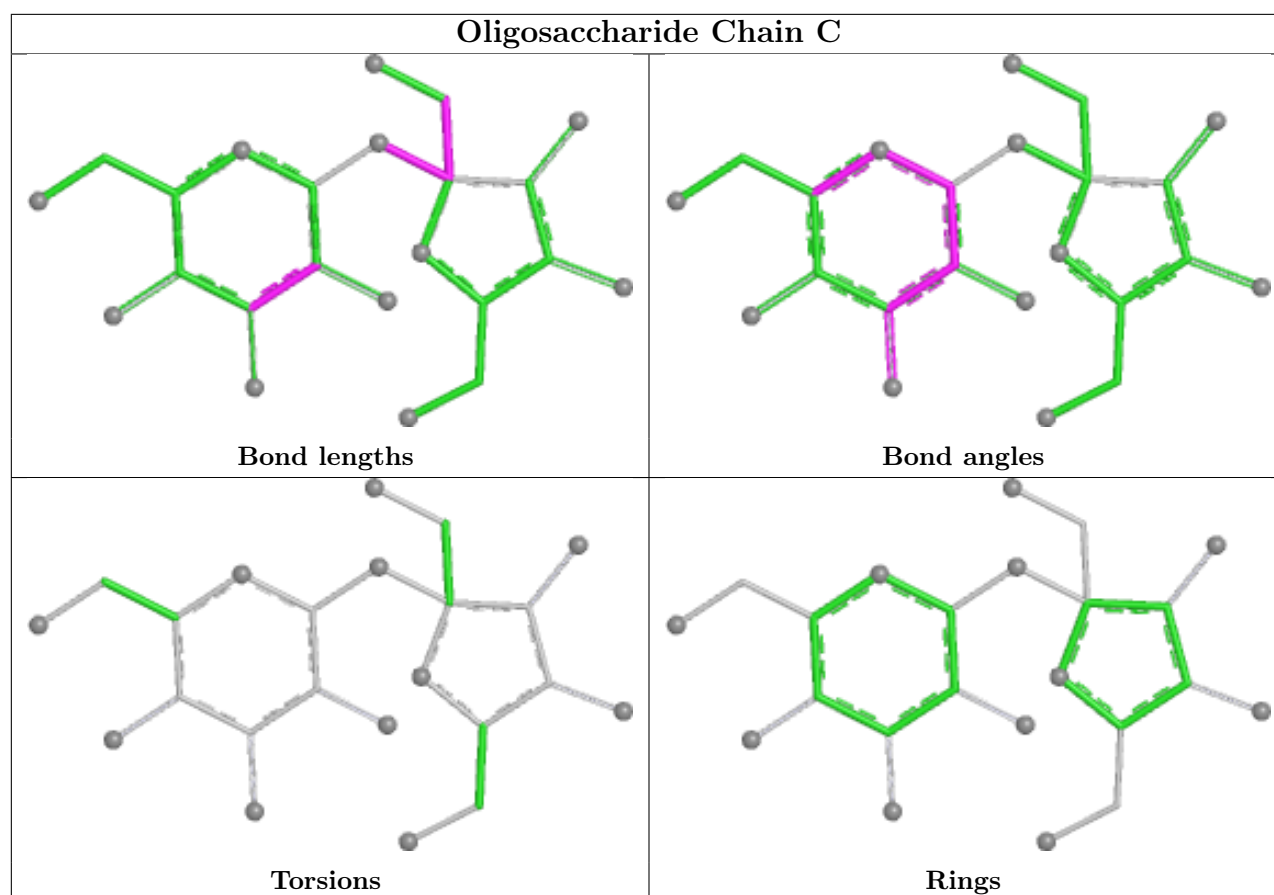
There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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](#)

Of 4 ligands modelled in this entry, 1 is monoatomic - leaving 3 for Mogul analysis.

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	$\# Z > 2$
7	739	B	3012	5	34,34,34	2.67	6 (17%)	36,45,45	1.81	6 (16%)
4	ACY	A	3002	-	3,3,3	1.26	1 (33%)	3,3,3	1.65	1 (33%)
6	FPP	B	3011	-	22,23,23	0.62	0	27,31,31	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
7	739	B	3012	5	-	4/38/40/40	0/1/1/1
6	FPP	B	3011	-	-	4/25/25/25	-

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	B	3012	739	OEM-SDM	11.57	1.66	1.44
7	B	3012	739	CCM-SDM	-6.08	1.70	1.78
7	B	3012	739	ODM-SDM	4.95	1.54	1.44
7	B	3012	739	O'F-CI	2.95	1.48	1.43
7	B	3012	739	CC-NI	-2.82	1.40	1.47
7	B	3012	739	CBI-CAI	2.11	1.59	1.53
4	A	3002	ACY	O-C	2.01	1.31	1.22

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	B	3012	739	ODM-SDM-CEM	-5.87	103.42	108.87
7	B	3012	739	OEM-SDM-CCM	5.23	112.10	108.33
7	B	3012	739	OEM-SDM-CEM	3.88	112.48	108.87
7	B	3012	739	CBI-CAI-NI	-2.77	105.67	111.64
4	A	3002	ACY	O-C-CH3	-2.24	113.36	122.53
7	B	3012	739	O'F-CAF-CF	-2.22	107.12	111.56
7	B	3012	739	ODM-SDM-CCM	2.12	109.86	108.33

There are no chirality outliers.

All (8) torsion outliers are listed below:

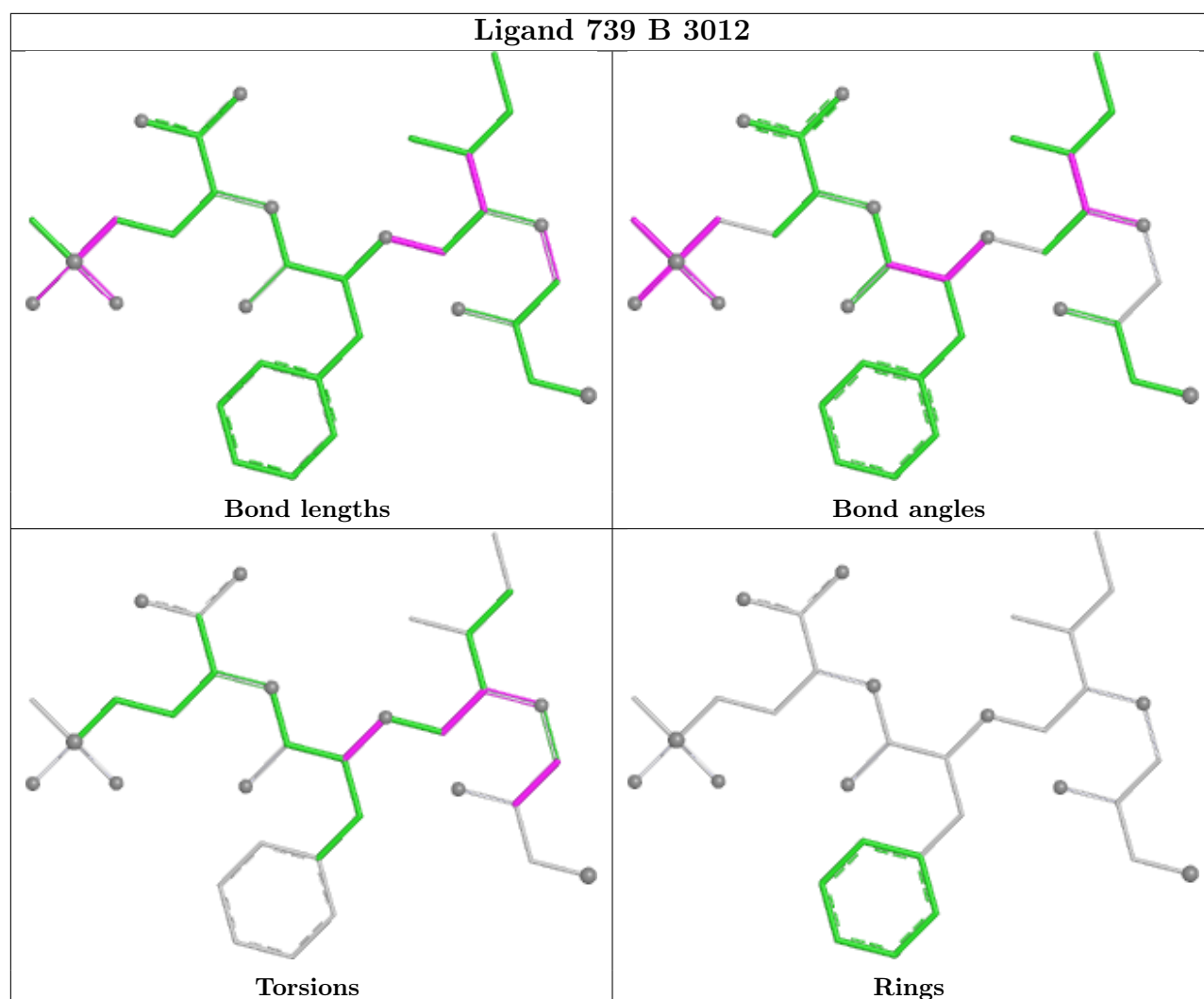
Mol	Chain	Res	Type	Atoms
7	B	3012	739	NI-CAI-CI-O'F
7	B	3012	739	NC-CAC-CC-NI
7	B	3012	739	CI-CAI-NI-CC
6	B	3011	FPP	PB-O3A-PA-O1A
7	B	3012	739	CBF-CAF-O'F-CI
6	B	3011	FPP	C10-C8-C9-C11
6	B	3011	FPP	C9-C11-C12-C13
6	B	3011	FPP	PB-O3A-PA-O2A

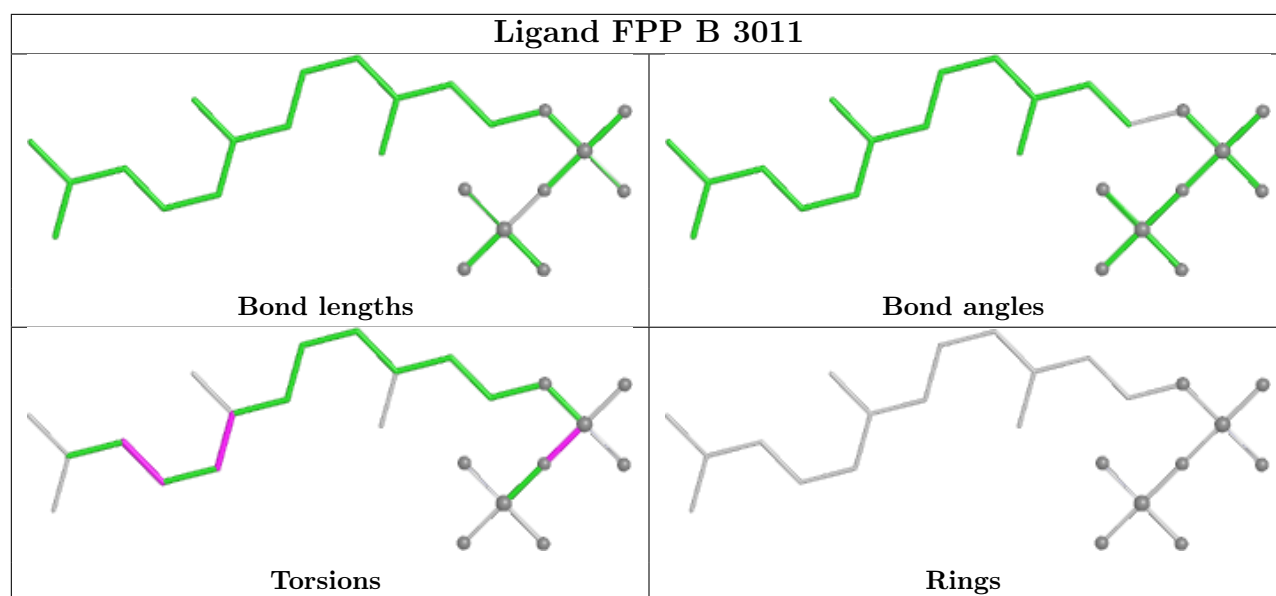
There are no ring outliers.

3 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	B	3012	739	2	0
4	A	3002	ACY	6	0
6	B	3011	FPP	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 all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [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	313/382 (81%)	-0.04	3 (0%) 79 80	26, 44, 65, 83	0
2	B	410/437 (93%)	-0.24	6 (1%) 72 73	25, 37, 60, 86	0
All	All	723/819 (88%)	-0.15	9 (1%) 76 77	25, 40, 62, 86	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	55	PHE	5.8
2	B	424	GLU	3.8
2	B	15	SER	3.1
1	A	304	PRO	2.9
2	B	118	ILE	2.7
1	A	330	LYS	2.4
2	B	121	ILE	2.4
2	B	120	GLN	2.2
2	B	381	MET	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.

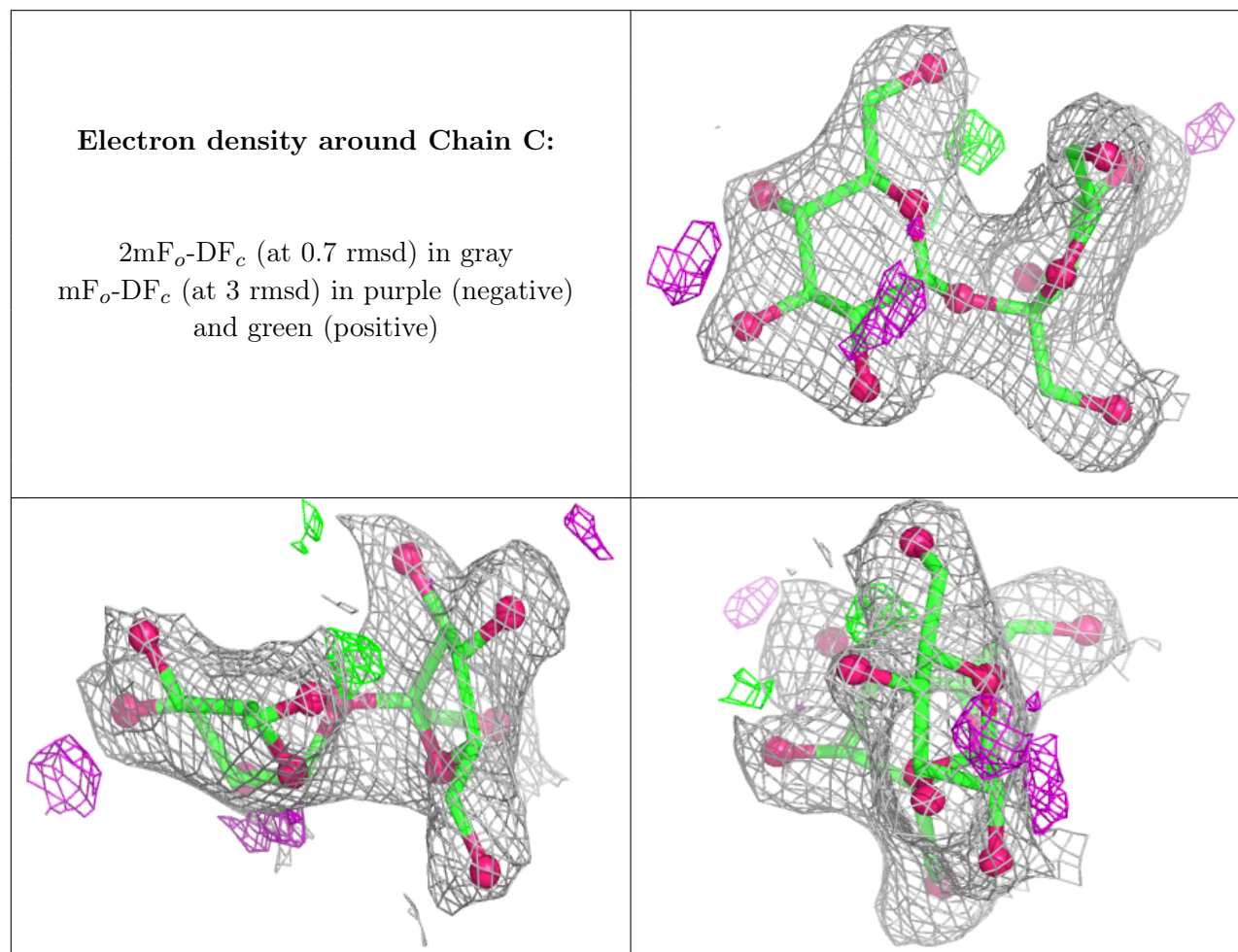
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	GLC	C	1	11/12	0.90	0.11	53,57,59,59	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	FRU	C	2	12/12	0.92	0.10	49,52,55,56	0

The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.



6.4 Ligands ⓘ

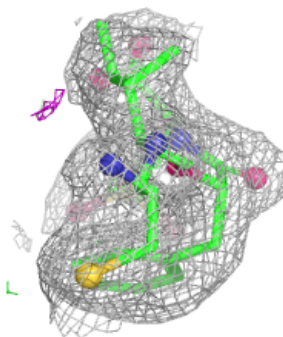
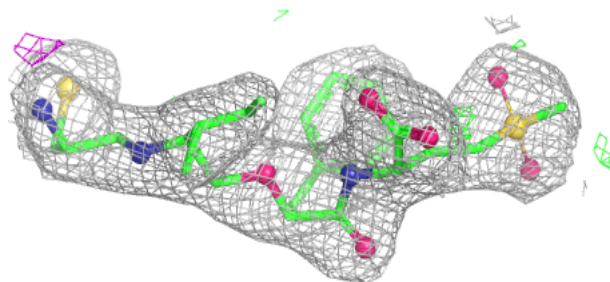
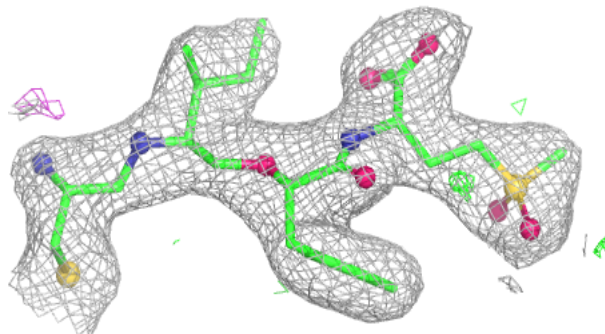
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

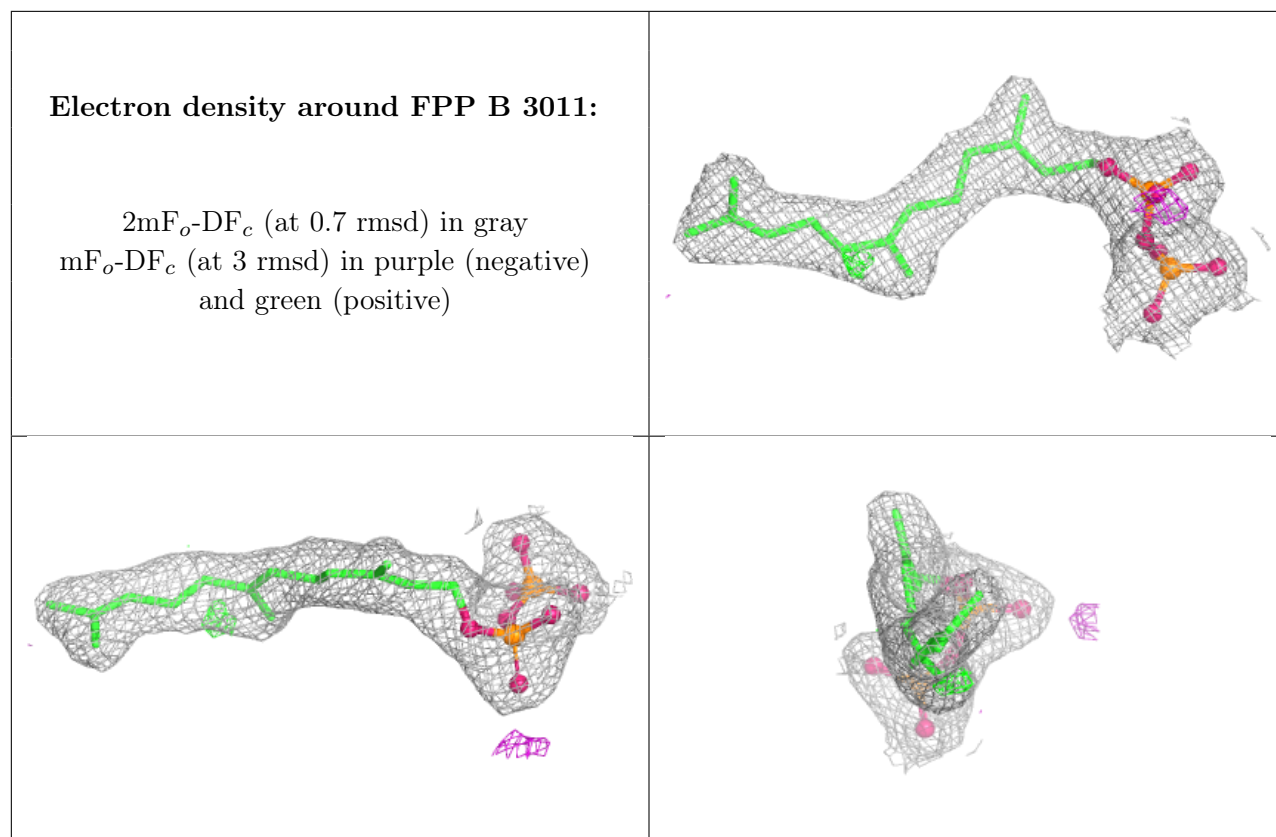
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	ACY	A	3002	4/4	0.90	0.18	53,54,55,57	0
7	739	B	3012	34/34	0.97	0.07	33,36,41,42	0
6	FPP	B	3011	24/24	0.98	0.06	22,31,33,34	0
5	ZN	B	1001	1/1	1.00	0.03	37,37,37,37	0

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 739 B 3012:

$2mF_o-DF_c$ (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.