Full wwPDB X-ray Structure Validation Report (i)

Aug 7, 2020 – 10:46 PM BST

PDB ID	:	4DG6
Title	:	Crystal structure of domains 1 and 2 of LRP6
Authors	:	Ceska, T.A.; Doyle, C.; Slocombe, P.
Deposited on	:	2012-01-25
Resolution	:	2.90 Å(reported)

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

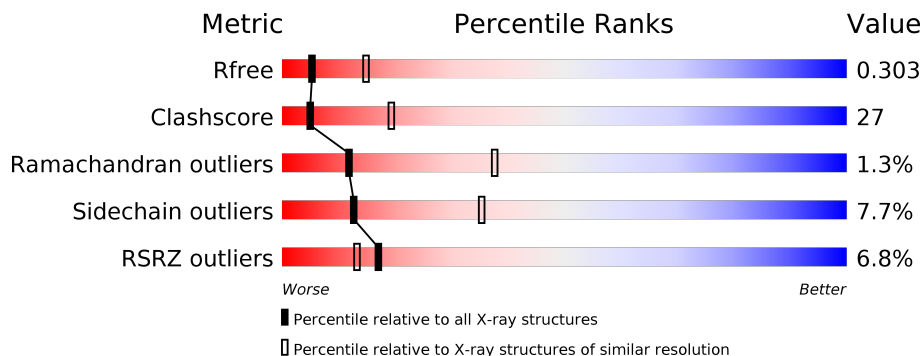|		
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.13.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.13.1

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 2.90 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.

Metric	Percentile Ranks	Value
Rfree		0.303
Clashscore		27
Ramachandran outliers		1.3%
Sidechain outliers		7.7%
RSRZ outliers		6.8%

Worse   *Better*

▮ Percentile relative to all X-ray structures

▯ Percentile relative to X-ray structures of similar resolution

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (2.90-2.90)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on 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 <=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	616	7% 52% 43% 6%
2	B	2	100%
2	C	2	50% 50%
2	D	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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	NAG	B	1	-	-	-	X
2	NAG	B	2	-	-	-	X

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 5005 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 Low-density lipoprotein receptor-related protein 6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	616	Total 4898	C 3107	N 836	O 931	S 24	0	0	0

- Molecule 2 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
2	B	2	Total 28	C 16	N 2	O 10	0	0	0
2	C	2	Total 28	C 16	N 2	O 10	0	0	0
2	D	2	Total 28	C 16	N 2	O 10	0	0	0

- Molecule 3 is PHOSPHATE ION (three-letter code: PO4) (formula: O_4P).

PO4

O3

O⁻

O2 O⁻ — P = O O1

P

O⁻

O4

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total 5	O 4	P 1	0	0
3	A	1	Total 5	O 4	P 1	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	13	Total 13	O 13	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: Low-density lipoprotein receptor-related protein 6
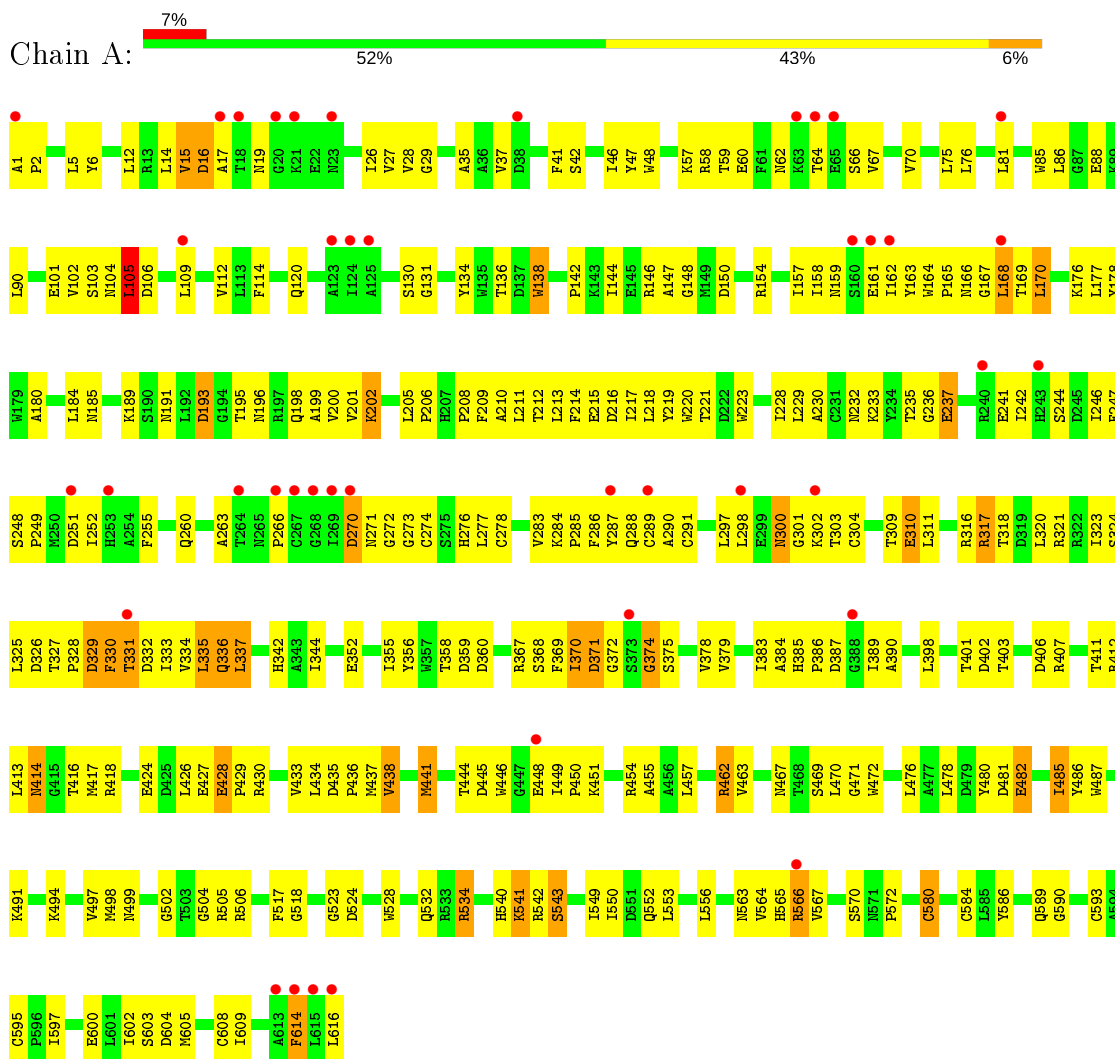
- Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

NAG1
NAG2

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

Chain C:

50%   50%

NAG1
NAG2

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

Chain D:

100%

NAG1
NAG2

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4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α, β, γ	109.36Å 109.36Å 130.88Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	32.09 – 2.90 32.09 – 2.90	Depositor EDS
% Data completeness (in resolution range)	99.7 (32.09-2.90) 99.7 (32.09-2.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$< I/\sigma(I) >$ [1]	2.89 (at 2.90Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.257 , 0.313 0.246 , 0.303	Depositor DCC
R_{free} test set	2010 reflections (9.82%)	wwPDB-VP
Wilson B-factor (Å^2)	82.5	Xtriage
Anisotropy	0.402	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 70.7	EDS
L-test for twinning[2]	$< |L| > = 0.50, < L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.026 for -h,-k,l	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	5005	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 5.80% of the height of the origin peak. No significant pseudotranslation is detected.*

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[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: PO4, 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	0.40	0/5015	0.57	0/6823

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	2

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	15	VAL	Peptide
1	A	374	GLY	Peptide

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	4898	0	4768	261	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	28	0	25	1	0
2	C	28	0	25	2	0
2	D	28	0	25	1	0
3	A	10	0	0	0	0
4	A	13	0	0	0	0
All	All	5005	0	4843	263	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 27.

All (263) 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:336:GLN:NE2	1:A:337:LEU:H	1.40	1.17
1:A:389:ILE:HD11	1:A:398:LEU:HD11	1.49	0.93
1:A:336:GLN:NE2	1:A:337:LEU:N	2.21	0.86
1:A:162:ILE:HB	1:A:165:PRO:HG3	1.56	0.85
1:A:549:ILE:HG13	1:A:550:ILE:HG22	1.61	0.82
1:A:597:ILE:HD12	1:A:597:ILE:H	1.45	0.81
1:A:193:ASP:HB2	1:A:195:THR:HG23	1.63	0.81
1:A:335:LEU:HD23	1:A:335:LEU:H	1.49	0.78
1:A:580:CYS:HG	1:A:593:CYS:HG	0.90	0.78
1:A:273:GLY:HA3	1:A:278:CYS:HB2	1.65	0.78
1:A:288:GLN:NE2	1:A:289:CYS:H	1.82	0.77
1:A:157:ILE:HG13	1:A:158:ILE:H	1.50	0.77
1:A:321:ARG:HG2	1:A:335:LEU:HB3	1.66	0.76
1:A:274:CYS:HG	1:A:289:CYS:HG	1.35	0.73
1:A:170:LEU:HD23	1:A:170:LEU:H	1.55	0.71
1:A:336:GLN:CD	1:A:337:LEU:H	1.93	0.71
1:A:228:ILE:HD12	1:A:244:SER:HB3	1.72	0.71
1:A:454:ARG:HG2	1:A:455:ALA:N	2.04	0.70
1:A:580:CYS:HB3	1:A:593:CYS:HB3	1.73	0.70
1:A:136:THR:HG22	1:A:144:ILE:HG12	1.72	0.70
1:A:330:PHE:O	1:A:332:ASP:N	2.25	0.69
1:A:310:GLU:HB2	1:A:325:LEU:HB2	1.74	0.69
2:D:1:NAG:O3	2:D:2:NAG:H2	1.92	0.69
1:A:370:ILE:HG22	1:A:371:ASP:H	1.58	0.69
1:A:266:PRO:HB2	1:A:287:TYR:HB3	1.74	0.68
1:A:218:LEU:O	1:A:230:ALA:HA	1.92	0.68
1:A:288:GLN:HE21	1:A:289:CYS:H	1.40	0.67
1:A:497:VAL:HG12	1:A:498:MET:H	1.57	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1:ALA:HB1	1:A:17:ALA:O	1.96	0.65
1:A:291:CYS:HG	1:A:304:CYS:HG	1.12	0.65
1:A:211:LEU:HD12	1:A:219:TYR:O	1.97	0.65
1:A:336:GLN:O	1:A:337:LEU:HB2	1.98	0.64
1:A:523:GLY:O	1:A:524:ASP:HB3	1.97	0.64
1:A:427:GLU:HG3	1:A:451:LYS:HE3	1.80	0.64
1:A:411:THR:HG22	1:A:418:ARG:HA	1.79	0.63
1:A:437:MET:O	1:A:438:VAL:HB	1.98	0.63
1:A:589:GLN:HG2	1:A:590:GLY:H	1.63	0.63
1:A:191:ASN:HB2	1:A:195:THR:OG1	1.98	0.63
1:A:335:LEU:CD2	1:A:335:LEU:H	2.12	0.62
1:A:131:GLY:O	1:A:148:GLY:HA2	1.99	0.62
1:A:37:VAL:HA	1:A:47:TYR:O	1.99	0.62
1:A:235:THR:HG23	1:A:237:GLU:H	1.64	0.62
1:A:370:ILE:HG22	1:A:371:ASP:N	2.14	0.62
1:A:263:ALA:O	2:B:1:NAG:H62	1.98	0.62
1:A:416:THR:HG22	1:A:417:MET:HG3	1.80	0.62
1:A:157:ILE:HG13	1:A:158:ILE:N	2.14	0.62
1:A:220:TRP:NE1	1:A:229:LEU:HD13	2.16	0.60
1:A:540:HIS:HB3	1:A:543:SER:HB3	1.83	0.60
1:A:60:GLU:HG2	1:A:66:SER:HB3	1.82	0.60
1:A:498:MET:SD	1:A:502:GLY:HA2	2.42	0.60
1:A:161:GLU:O	1:A:161:GLU:HG3	2.03	0.59
1:A:454:ARG:HH21	1:A:463:VAL:HG21	1.68	0.59
1:A:532:GLN:O	1:A:534:ARG:HG2	2.03	0.59
1:A:291:CYS:HG	1:A:304:CYS:CB	2.16	0.58
1:A:441:MET:HB3	1:A:457:LEU:HD23	1.86	0.58
1:A:429:PRO:HA	1:A:445:ASP:HB3	1.85	0.58
1:A:524:ASP:HA	1:A:541:LYS:NZ	2.19	0.58
1:A:166:ASN:CG	1:A:167:GLY:H	2.07	0.57
1:A:358:THR:HB	1:A:386:PRO:HB2	1.87	0.57
1:A:176:LYS:NZ	1:A:198:GLN:HE22	2.03	0.57
1:A:291:CYS:CB	1:A:304:CYS:HG	2.18	0.57
1:A:372:GLY:C	1:A:374:GLY:H	2.08	0.57
1:A:602:ILE:HG22	1:A:603:SER:N	2.20	0.57
1:A:159:ASN:HA	1:A:162:ILE:HD11	1.85	0.57
1:A:385:HIS:HB2	1:A:403:THR:HB	1.87	0.57
1:A:499:ASN:ND2	1:A:504:GLY:N	2.53	0.57
1:A:158:ILE:N	1:A:158:ILE:HD12	2.20	0.56
1:A:199:ALA:HB1	1:A:202:LYS:HB3	1.87	0.56
1:A:130:SER:OG	1:A:131:GLY:N	2.39	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:370:ILE:CG2	1:A:371:ASP:H	2.15	0.56
1:A:241:GLU:O	1:A:241:GLU:HG3	2.05	0.56
1:A:147:ALA:HB2	1:A:154:ARG:HA	1.88	0.56
1:A:291:CYS:SG	1:A:297:LEU:HD13	2.45	0.56
1:A:176:LYS:HZ1	1:A:198:GLN:HE22	1.55	0.55
1:A:166:ASN:HB3	1:A:180:ALA:HB3	1.87	0.55
1:A:101:GLU:CD	1:A:112:VAL:HG22	2.26	0.55
1:A:271:ASN:ND2	1:A:273:GLY:O	2.40	0.55
1:A:300:ASN:N	1:A:300:ASN:HD22	2.03	0.54
1:A:580:CYS:SG	1:A:584:CYS:HB2	2.47	0.54
1:A:170:LEU:N	1:A:170:LEU:HD23	2.21	0.54
1:A:446:TRP:CZ2	1:A:472:TRP:HB3	2.42	0.54
1:A:15:VAL:O	1:A:16:ASP:HB2	2.07	0.54
1:A:298:LEU:HB2	1:A:300:ASN:ND2	2.23	0.54
1:A:2:PRO:HB3	1:A:214:PHE:CD2	2.43	0.54
1:A:329:ASP:O	1:A:331:THR:N	2.40	0.53
1:A:273:GLY:HA3	1:A:278:CYS:CB	2.37	0.53
1:A:498:MET:HG2	1:A:499:ASN:O	2.09	0.53
1:A:524:ASP:HA	1:A:541:LYS:HD2	1.91	0.53
1:A:335:LEU:N	1:A:335:LEU:CD2	2.71	0.53
1:A:602:ILE:HD11	1:A:609:ILE:HG12	1.91	0.53
1:A:316:ARG:NH1	1:A:321:ARG:HH21	2.06	0.53
1:A:471:GLY:HA3	1:A:491:LYS:HB3	1.91	0.53
1:A:289:CYS:SG	1:A:302:LYS:HG2	2.48	0.53
1:A:478:LEU:HD23	1:A:480:TYR:CE1	2.44	0.53
1:A:485:ILE:HG23	1:A:498:MET:HB3	1.91	0.53
1:A:580:CYS:CB	1:A:593:CYS:HG	2.21	0.52
1:A:602:ILE:HG22	1:A:603:SER:H	1.74	0.52
1:A:216:ASP:O	1:A:233:LYS:HB2	2.09	0.52
1:A:284:LYS:HB2	1:A:285:PRO:HD3	1.90	0.52
1:A:344:ILE:HG21	1:A:387:ASP:HA	1.91	0.52
1:A:178:TYR:CD1	1:A:189:LYS:HB2	2.45	0.52
1:A:321:ARG:HG2	1:A:335:LEU:CB	2.39	0.52
1:A:320:LEU:O	1:A:335:LEU:HB2	2.10	0.51
1:A:327:THR:HB	1:A:328:PRO:HD2	1.90	0.51
1:A:540:HIS:CB	1:A:543:SER:HB3	2.41	0.51
1:A:60:GLU:HB2	1:A:64:THR:HG21	1.93	0.51
1:A:426:LEU:HD13	1:A:448:GLU:OE1	2.11	0.51
1:A:595:CYS:HG	1:A:608:CYS:HG	0.52	0.51
1:A:144:ILE:HD13	1:A:168:LEU:HD11	1.92	0.51
1:A:235:THR:HG23	1:A:237:GLU:HB2	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:220:TRP:CE2	1:A:229:LEU:HB2	2.46	0.51
1:A:46:ILE:O	1:A:58:ARG:HA	2.11	0.51
1:A:164:TRP:N	1:A:165:PRO:HD3	2.26	0.51
1:A:26:ILE:O	1:A:26:ILE:HG22	2.10	0.51
1:A:384:ALA:O	1:A:386:PRO:HD3	2.11	0.50
1:A:170:LEU:HB3	1:A:177:LEU:HD13	1.93	0.50
1:A:486:TYR:OH	1:A:541:LYS:HG3	2.11	0.50
1:A:553:LEU:HD23	1:A:556:LEU:HD21	1.93	0.50
1:A:136:THR:HG23	1:A:168:LEU:HD22	1.93	0.50
1:A:390:ALA:HB1	1:A:434:LEU:HG	1.93	0.50
1:A:524:ASP:HA	1:A:541:LYS:CD	2.41	0.50
1:A:383:ILE:HG23	1:A:386:PRO:HG3	1.93	0.49
1:A:517:PHE:CE1	1:A:528:TRP:HB2	2.47	0.49
1:A:58:ARG:NH1	1:A:105:LEU:O	2.45	0.49
1:A:534:ARG:HA	1:A:553:LEU:O	2.13	0.49
1:A:457:LEU:HB2	1:A:586:TYR:O	2.13	0.49
1:A:104:ASN:HD21	1:A:109:LEU:HD12	1.78	0.49
1:A:330:PHE:C	1:A:332:ASP:H	2.15	0.49
1:A:138:TRP:HB3	1:A:165:PRO:HD2	1.94	0.49
1:A:436:PRO:HA	1:A:457:LEU:HD11	1.94	0.49
1:A:563:ASN:HB3	1:A:566:ARG:HB3	1.94	0.49
1:A:572:PRO:HD2	1:A:586:TYR:CE2	2.48	0.49
1:A:19:ASN:HD22	1:A:242:ILE:HG23	1.78	0.48
1:A:330:PHE:C	1:A:332:ASP:N	2.66	0.48
1:A:57:LYS:HA	1:A:70:VAL:HG23	1.94	0.48
1:A:298:LEU:HB2	1:A:300:ASN:HD21	1.77	0.48
1:A:283:VAL:HG12	1:A:285:PRO:HD2	1.95	0.48
1:A:270:ASP:O	1:A:272:GLY:N	2.45	0.48
1:A:270:ASP:C	1:A:272:GLY:H	2.16	0.48
1:A:277:LEU:HB2	1:A:290:ALA:HB3	1.95	0.48
1:A:406:ASP:HA	1:A:429:PRO:HD2	1.95	0.48
1:A:469:SER:O	1:A:470:LEU:HD12	2.14	0.48
1:A:246:ILE:O	1:A:249:PRO:HD3	2.13	0.48
1:A:355:ILE:O	1:A:367:ARG:HA	2.13	0.48
1:A:14:LEU:HD21	1:A:255:PHE:CD1	2.49	0.48
1:A:232:ASN:HB2	1:A:235:THR:O	2.13	0.48
1:A:335:LEU:N	1:A:335:LEU:HD23	2.18	0.48
1:A:428:GLU:HA	1:A:429:PRO:HD3	1.71	0.48
1:A:426:LEU:HB3	1:A:448:GLU:OE1	2.15	0.47
1:A:138:TRP:CD1	1:A:138:TRP:C	2.87	0.47
1:A:367:ARG:O	1:A:375:SER:HA	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:505:ARG:O	1:A:506:ARG:HB3	2.14	0.47
1:A:414:ASN:OD1	1:A:414:ASN:N	2.45	0.47
1:A:450:PRO:HG2	1:A:470:LEU:O	2.15	0.47
1:A:14:LEU:HD11	1:A:255:PHE:CE1	2.49	0.47
1:A:27:VAL:HG11	1:A:46:ILE:HD12	1.97	0.47
1:A:467:ASN:HB2	2:C:1:NAG:N2	2.30	0.47
1:A:134:TYR:CE1	1:A:146:ARG:HG3	2.50	0.47
1:A:316:ARG:HH11	1:A:321:ARG:HH21	1.61	0.47
1:A:201:VAL:HG22	1:A:236:GLY:HA2	1.96	0.47
1:A:401:THR:HB	1:A:429:PRO:HB2	1.97	0.47
1:A:589:GLN:HG2	1:A:590:GLY:N	2.28	0.47
1:A:166:ASN:ND2	1:A:208:PRO:O	2.48	0.46
1:A:356:TYR:OH	1:A:413:LEU:HB3	2.15	0.46
1:A:163:TYR:HB3	1:A:184:LEU:HG	1.98	0.46
1:A:213:LEU:C	1:A:213:LEU:HD12	2.35	0.46
1:A:542:ARG:O	1:A:543:SER:CB	2.63	0.46
1:A:216:ASP:C	1:A:233:LYS:HB2	2.36	0.46
1:A:300:ASN:ND2	1:A:301:GLY:N	2.63	0.46
1:A:497:VAL:HG12	1:A:498:MET:N	2.28	0.46
1:A:75:LEU:O	1:A:76:LEU:HD23	2.16	0.46
1:A:148:GLY:C	1:A:150:ASP:H	2.19	0.45
1:A:114:PHE:CD1	1:A:154:ARG:HD2	2.51	0.45
1:A:437:MET:O	1:A:438:VAL:CB	2.59	0.45
1:A:120:GLN:HB2	1:A:138:TRP:O	2.16	0.45
1:A:138:TRP:HA	1:A:142:PRO:HA	1.99	0.45
1:A:215:GLU:O	1:A:217:ILE:HG13	2.17	0.45
1:A:369:PHE:N	1:A:374:GLY:HA3	2.31	0.45
1:A:383:ILE:CG2	1:A:386:PRO:HG3	2.46	0.45
1:A:499:ASN:ND2	1:A:504:GLY:CA	2.79	0.45
1:A:81:LEU:HD12	1:A:81:LEU:HA	1.76	0.45
1:A:221:THR:HG21	1:A:249:PRO:HB2	1.99	0.45
1:A:321:ARG:CZ	1:A:335:LEU:HD13	2.46	0.45
1:A:435:ASP:O	1:A:437:MET:O	2.34	0.45
1:A:59:THR:HG22	1:A:67:VAL:HA	1.98	0.45
1:A:317:ARG:HG3	1:A:342:HIS:CD2	2.52	0.45
1:A:59:THR:HG22	1:A:66:SER:O	2.17	0.45
1:A:210:ALA:HB3	1:A:252:ILE:HG22	1.99	0.45
1:A:481:ASP:C	1:A:482:GLU:HG2	2.37	0.45
1:A:336:GLN:HB3	1:A:337:LEU:HD22	1.99	0.44
1:A:5:LEU:HD13	1:A:14:LEU:HG	2.00	0.44
1:A:285:PRO:O	1:A:286:PHE:HB2	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:369:PHE:H	1:A:374:GLY:HA3	1.82	0.44
1:A:352:GLU:HA	1:A:352:GLU:OE1	2.18	0.44
1:A:476:LEU:HD11	1:A:485:ILE:HD11	2.00	0.44
1:A:437:MET:HE3	1:A:570:SER:H	1.83	0.44
1:A:300:ASN:H	1:A:300:ASN:HD22	1.64	0.44
1:A:553:LEU:O	1:A:556:LEU:HG	2.18	0.44
1:A:158:ILE:CD1	1:A:158:ILE:N	2.80	0.44
1:A:389:ILE:HG12	1:A:390:ALA:N	2.31	0.44
1:A:427:GLU:HG3	1:A:451:LYS:CE	2.45	0.44
1:A:35:ALA:O	1:A:251:ASP:HB2	2.17	0.44
1:A:291:CYS:SG	1:A:304:CYS:CB	3.05	0.43
1:A:102:VAL:HG22	1:A:103:SER:N	2.32	0.43
1:A:342:HIS:O	1:A:359:ASP:HA	2.19	0.43
1:A:62:ASN:O	1:A:64:THR:HG22	2.19	0.43
1:A:157:ILE:C	1:A:158:ILE:HD12	2.39	0.43
1:A:205:LEU:HB3	1:A:206:PRO:HD3	2.01	0.43
1:A:19:ASN:HD21	1:A:242:ILE:HD13	1.83	0.43
1:A:372:GLY:C	1:A:374:GLY:N	2.71	0.43
1:A:5:LEU:O	1:A:252:ILE:HA	2.19	0.43
1:A:297:LEU:HD12	1:A:303:THR:O	2.19	0.43
1:A:316:ARG:HD3	1:A:321:ARG:HE	1.83	0.43
1:A:424:GLU:HG3	1:A:462:ARG:HH22	1.84	0.43
1:A:223:TRP:CD1	1:A:248:SER:HA	2.54	0.43
1:A:291:CYS:CB	1:A:304:CYS:SG	3.05	0.43
1:A:430:ARG:HE	1:A:430:ARG:HB3	1.68	0.43
1:A:214:PHE:O	1:A:217:ILE:HD12	2.19	0.43
1:A:323:ILE:HG22	1:A:324:SER:O	2.18	0.43
1:A:604:ASP:O	1:A:605:MET:HB2	2.19	0.43
1:A:499:ASN:HB2	1:A:504:GLY:H	1.84	0.43
1:A:518:GLY:O	1:A:528:TRP:HA	2.19	0.43
1:A:229:LEU:N	1:A:229:LEU:HD12	2.34	0.42
1:A:169:THR:O	1:A:177:LEU:HD12	2.18	0.42
1:A:200:VAL:HG12	1:A:201:VAL:HG23	2.01	0.42
1:A:317:ARG:HB3	1:A:318:THR:HG23	2.01	0.42
1:A:384:ALA:HB3	1:A:402:ASP:OD1	2.18	0.42
1:A:90:LEU:O	1:A:102:VAL:HA	2.19	0.42
1:A:205:LEU:HA	1:A:205:LEU:HD12	1.87	0.42
1:A:310:GLU:HB2	1:A:325:LEU:HD12	2.02	0.42
1:A:228:ILE:H	1:A:244:SER:CB	2.32	0.42
1:A:212:THR:CG2	1:A:219:TYR:HB2	2.49	0.42
1:A:277:LEU:CB	1:A:290:ALA:HB3	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:379:VAL:CG1	1:A:383:ILE:HD11	2.50	0.42
1:A:438:VAL:HG12	1:A:438:VAL:O	2.20	0.42
1:A:517:PHE:CZ	1:A:528:TRP:HB2	2.55	0.42
1:A:309:THR:H	1:A:326:ASP:CG	2.23	0.42
1:A:6:TYR:CE1	1:A:246:ILE:HG21	2.53	0.42
1:A:378:VAL:HG12	1:A:379:VAL:HG23	2.02	0.42
1:A:614:PHE:CD1	1:A:616:LEU:HG	2.55	0.42
1:A:12:LEU:HD11	1:A:48:TRP:HB3	2.02	0.41
1:A:168:LEU:HA	1:A:168:LEU:HD12	1.79	0.41
1:A:276:HIS:ND1	1:A:304:CYS:HB2	2.34	0.41
1:A:138:TRP:CD2	1:A:164:TRP:HA	2.55	0.41
1:A:28:VAL:HG12	1:A:29:GLY:H	1.85	0.41
1:A:104:ASN:C	1:A:106:ASP:H	2.24	0.41
1:A:311:LEU:HD12	1:A:324:SER:HA	2.02	0.41
1:A:564:VAL:O	1:A:565:HIS:HB2	2.19	0.41
1:A:600:GLU:O	1:A:608:CYS:HA	2.20	0.41
2:C:1:NAG:O3	2:C:1:NAG:H83	2.21	0.41
1:A:180:ALA:HB1	1:A:208:PRO:HG2	2.02	0.41
1:A:85:TRP:CD2	1:A:260:GLN:HB3	2.56	0.41
1:A:247:PHE:O	1:A:248:SER:HB3	2.20	0.41
1:A:166:ASN:HD21	1:A:209:PHE:C	2.23	0.41
1:A:16:ASP:OD2	1:A:19:ASN:HA	2.21	0.41
1:A:449:ILE:O	1:A:449:ILE:HG23	2.21	0.41
1:A:208:PRO:HA	1:A:221:THR:O	2.20	0.41
1:A:86:LEU:O	1:A:88:GLU:HG2	2.21	0.41
1:A:597:ILE:CD1	1:A:597:ILE:H	2.19	0.40
1:A:487:TRP:HZ3	1:A:498:MET:HB2	1.85	0.40
1:A:178:TYR:HD1	1:A:189:LYS:HB2	1.84	0.40
1:A:310:GLU:HG2	1:A:310:GLU:H	1.73	0.40
1:A:220:TRP:CZ2	1:A:229:LEU:HD22	2.56	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	614/616 (100%)	508 (83%)	98 (16%)	8 (1%)	12	37

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	438	VAL
1	A	543	SER
1	A	16	ASP
1	A	330	PHE
1	A	337	LEU
1	A	105	LEU
1	A	331	THR
1	A	370	ILE

5.3.2 Protein sidechains (i)

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	531/531 (100%)	490 (92%)	41 (8%)	13	35

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	41	PHE
1	A	42	SER
1	A	105	LEU
1	A	138	TRP
1	A	168	LEU
1	A	170	LEU
1	A	185	ASN
1	A	193	ASP
1	A	196	ASN
1	A	202	LYS
1	A	237	GLU

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Mol	Chain	Res	Type
1	A	270	ASP
1	A	300	ASN
1	A	310	GLU
1	A	317	ARG
1	A	329	ASP
1	A	333	ILE
1	A	334	VAL
1	A	335	LEU
1	A	336	GLN
1	A	360	ASP
1	A	368	SER
1	A	371	ASP
1	A	407	ARG
1	A	412	ARG
1	A	414	ASN
1	A	428	GLU
1	A	433	VAL
1	A	441	MET
1	A	444	THR
1	A	462	ARG
1	A	482	GLU
1	A	485	ILE
1	A	494	LYS
1	A	534	ARG
1	A	541	LYS
1	A	552	GLN
1	A	566	ARG
1	A	567	VAL
1	A	580	CYS
1	A	614	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	19	ASN
1	A	166	ASN
1	A	185	ASN
1	A	188	HIS
1	A	196	ASN
1	A	198	GLN
1	A	207	HIS
1	A	288	GLN

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Mol	Chain	Res	Type
1	A	300	ASN
1	A	336	GLN
1	A	499	ASN

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)

6 monosaccharides are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	#$|Z| > 2$	Counts	RMSZ	#$|Z| > 2$
2	NAG	B	1	1,2	14,14,15	0.74	0	17,19,21	0.79	0
2	NAG	B	2	2	14,14,15	0.68	0	17,19,21	0.65	1 (5%)
2	NAG	C	1	1,2	14,14,15	0.67	0	17,19,21	0.65	0
2	NAG	C	2	2	14,14,15	0.81	0	17,19,21	0.63	0
2	NAG	D	1	1,2	14,14,15	0.82	0	17,19,21	1.18	2 (11%)
2	NAG	D	2	2	14,14,15	0.80	1 (7%)	17,19,21	0.63	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	B	1	1,2	-	2/6/23/26	0/1/1/1

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

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	2	NAG	C1-C2	2.26	1.55	1.52

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	D	1	NAG	C4-C3-C2	3.68	116.41	111.02
2	B	2	NAG	C2-N2-C7	-2.12	119.88	122.90
2	D	1	NAG	C2-N2-C7	-2.09	119.93	122.90

There are no chirality outliers.

All (23) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	1	NAG	C8-C7-N2-C2
2	C	1	NAG	O7-C7-N2-C2
2	B	1	NAG	C8-C7-N2-C2
2	B	1	NAG	O7-C7-N2-C2
2	D	1	NAG	C3-C2-N2-C7
2	D	1	NAG	C8-C7-N2-C2
2	D	1	NAG	O7-C7-N2-C2
2	B	2	NAG	C1-C2-N2-C7
2	B	2	NAG	C8-C7-N2-C2
2	B	2	NAG	O7-C7-N2-C2
2	C	2	NAG	C1-C2-N2-C7
2	C	2	NAG	C8-C7-N2-C2
2	C	2	NAG	O7-C7-N2-C2
2	B	2	NAG	C4-C5-C6-O6
2	B	2	NAG	O5-C5-C6-O6
2	D	2	NAG	C8-C7-N2-C2
2	D	2	NAG	O7-C7-N2-C2
2	D	1	NAG	O5-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
2	D	2	NAG	C4-C5-C6-O6
2	D	2	NAG	O5-C5-C6-O6
2	D	1	NAG	C4-C5-C6-O6
2	B	2	NAG	C3-C2-N2-C7
2	C	2	NAG	C3-C2-N2-C7

There are no ring outliers.

4 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	1	NAG	2	0
2	B	1	NAG	1	0
2	D	2	NAG	1	0
2	D	1	NAG	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.

Oligosaccharide Chain B

Bond lengths

Bond angles

Torsions

Rings

Oligosaccharide Chain C

Bond lengths

Bond angles

Torsions

Rings
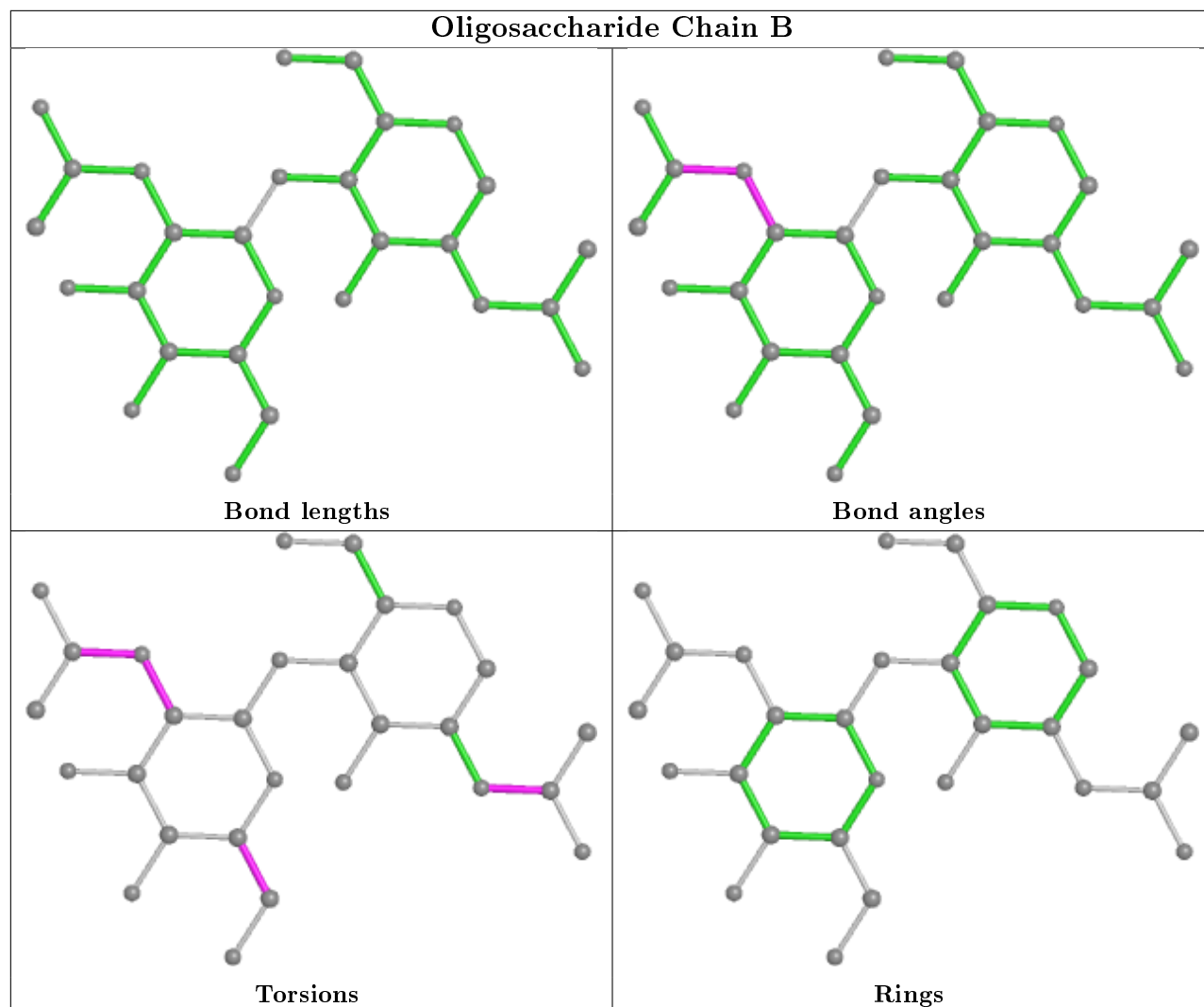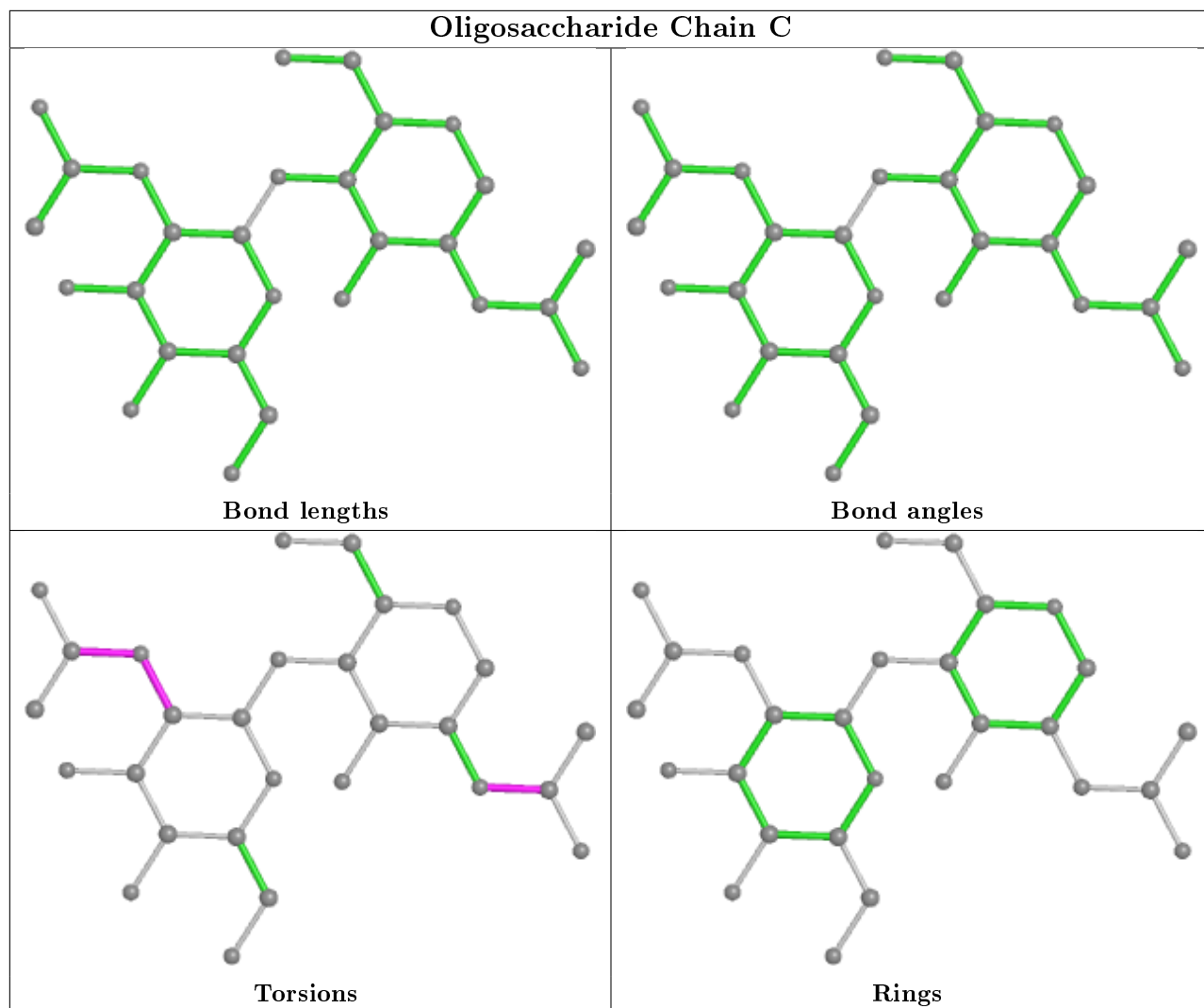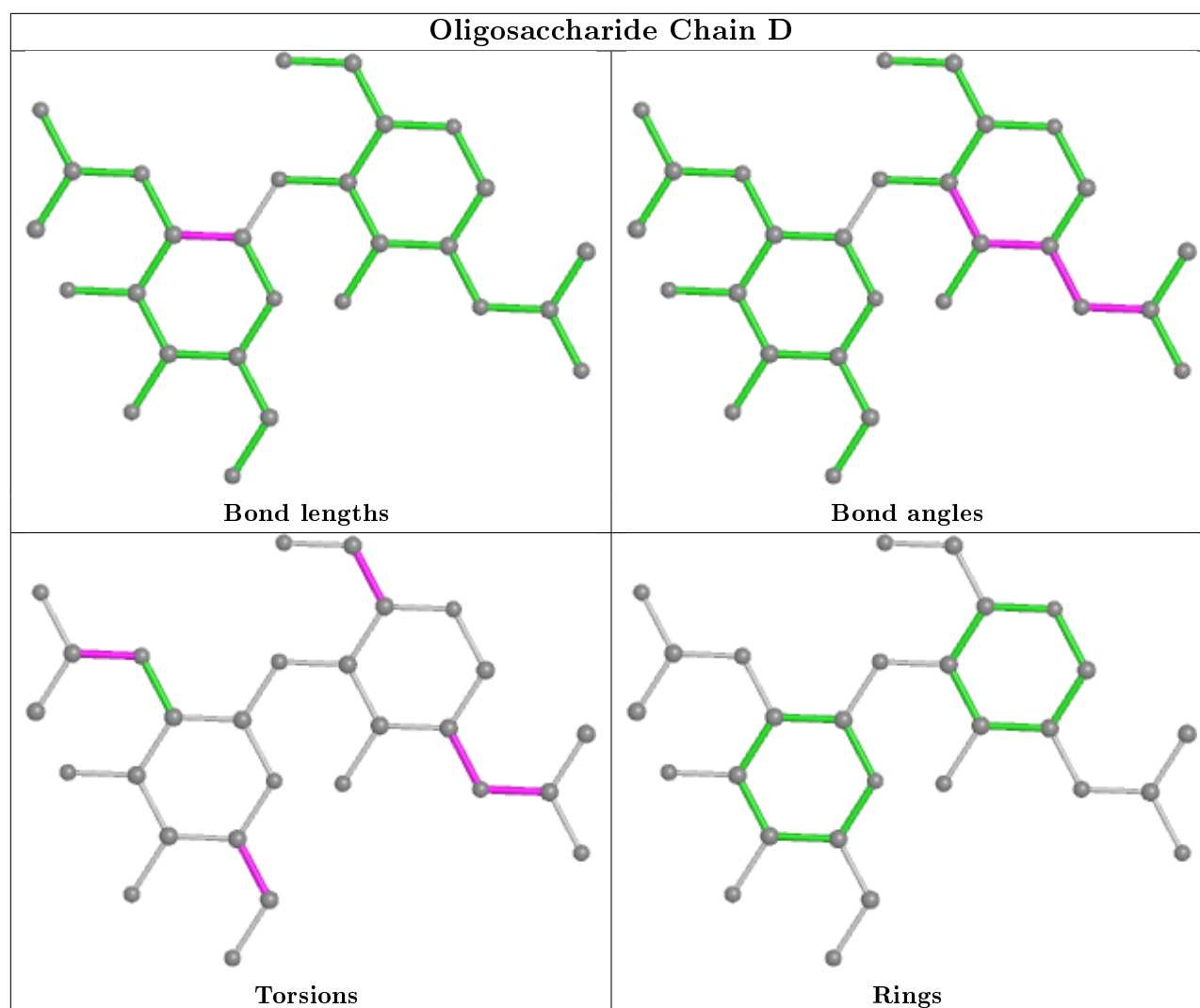
5.6 Ligand geometry (i)

2 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	$\#|Z| > 2$
3	PO4	A	708	-	4,4,4	1.59	0	6,6,6	0.45	0
3	PO4	A	707	-	4,4,4	1.52	0	6,6,6	0.44	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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(Å²)	Q<0.9
1	A	616/616 (100%)	0.32	42 (6%)	17	13	32, 73, 101, 118	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1	ALA	7.7
1	A	613	ALA	5.7
1	A	615	LEU	5.4
1	A	614	PHE	4.5
1	A	18	THR	4.5
1	A	240	ARG	4.2
1	A	331	THR	4.1
1	A	21	LYS	4.0
1	A	243	HIS	3.8
1	A	269	ILE	3.7
1	A	448	GLU	3.6
1	A	63	LYS	3.6
1	A	264	THR	3.6
1	A	266	PRO	3.5
1	A	160	SER	3.4
1	A	616	LEU	3.4
1	A	268	GLY	3.4
1	A	162	ILE	3.1
1	A	125	ALA	3.0
1	A	124	ILE	2.8
1	A	287	TYR	2.8
1	A	109	LEU	2.7
1	A	123	ALA	2.7
1	A	64	THR	2.7
1	A	566	ARG	2.6
1	A	267	CYS	2.6
1	A	17	ALA	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	38	ASP	2.6
1	A	65	GLU	2.5
1	A	81	LEU	2.5
1	A	251	ASP	2.5
1	A	168	LEU	2.4
1	A	289	CYS	2.4
1	A	388	GLY	2.4
1	A	20	GLY	2.3
1	A	270	ASP	2.3
1	A	23	ASN	2.3
1	A	298	LEU	2.2
1	A	253	HIS	2.2
1	A	373	SER	2.1
1	A	161	GLU	2.1
1	A	302	LYS	2.0

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
2	NAG	C	2	14/15	0.53	0.37	116,120,123,126	0
2	NAG	B	1	14/15	0.57	0.52	111,113,114,116	0
2	NAG	B	2	14/15	0.65	0.69	117,119,120,120	0
2	NAG	D	1	14/15	0.73	0.27	92,97,98,101	0
2	NAG	D	2	14/15	0.78	0.29	98,102,104,104	0
2	NAG	C	1	14/15	0.84	0.28	109,112,114,116	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.
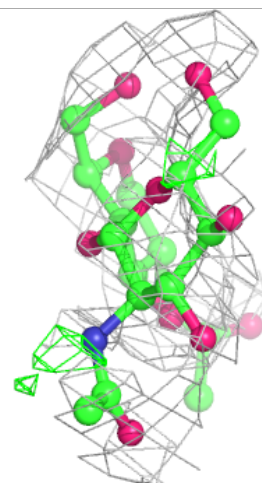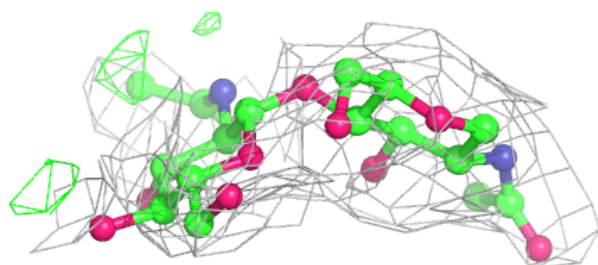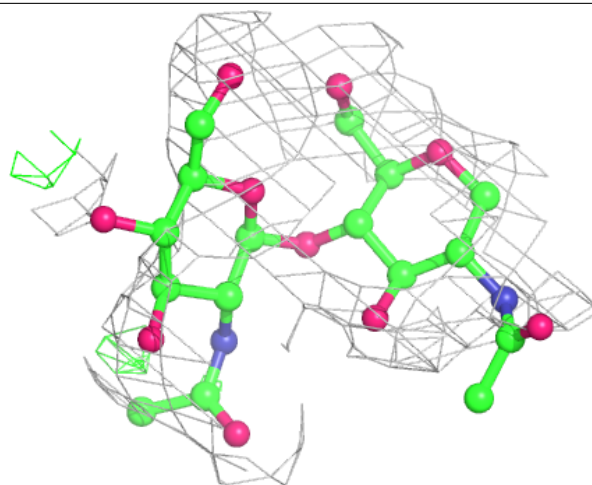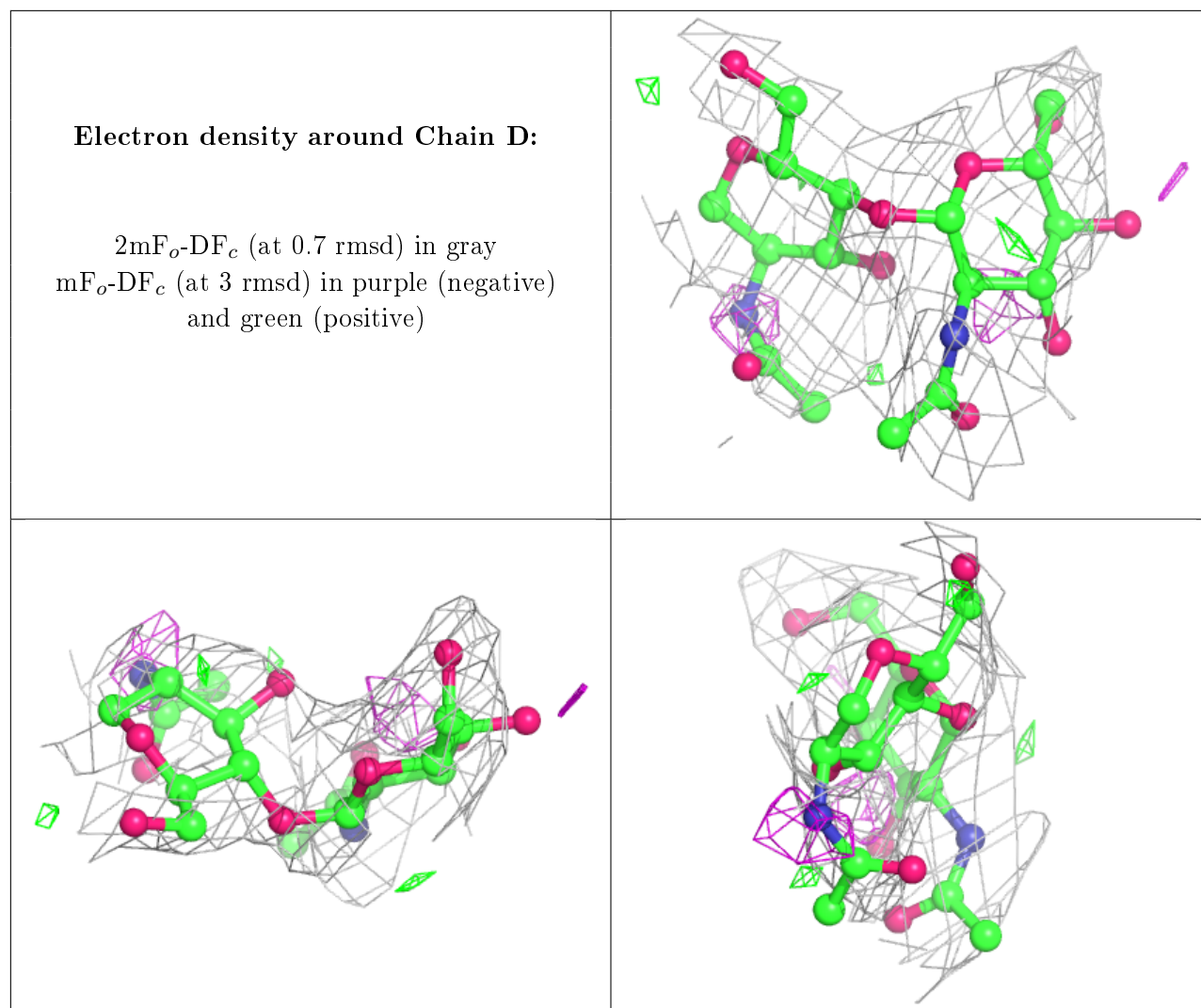
Electron density around Chain B:

2mF_o-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)
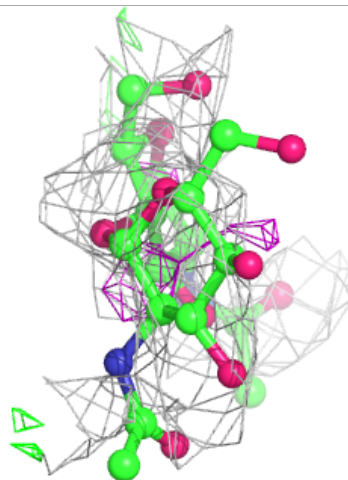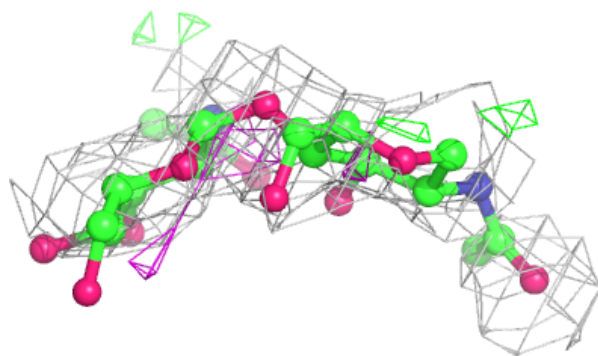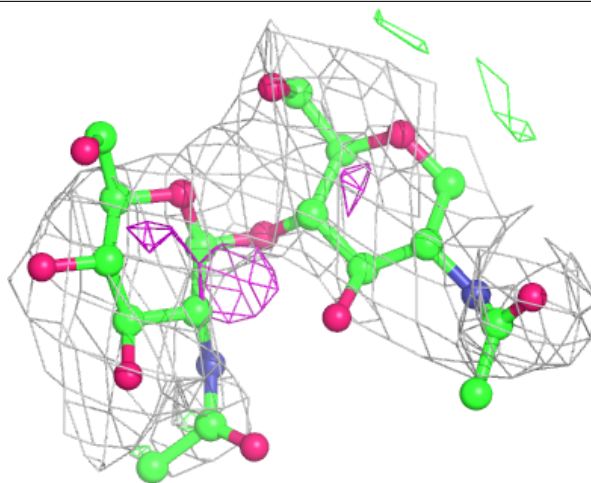
Electron density around Chain C:

$2mF_o$-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 D:

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

6.4 Ligands ⓘ

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	PO4	A	707	5/5	0.93	0.49	82,83,84,84	0
3	PO4	A	708	5/5	0.95	0.09	89,89,91,91	0

6.5 Other polymers ⓘ

There are no such residues in this entry.