Full wwPDB X-ray Structure Validation Report (i)

Jan 18, 2024 – 04:08 PM JST

PDB ID	:	6IQH
Title	:	X-ray crystal structure of covalent-bonded complex of Fc and peptide
Authors	:	Adachi, M.; Ito, Y.
Deposited on	:	2018-11-08
Resolution	:	3.00 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 (i) 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 (i)) were used in the production of this report:

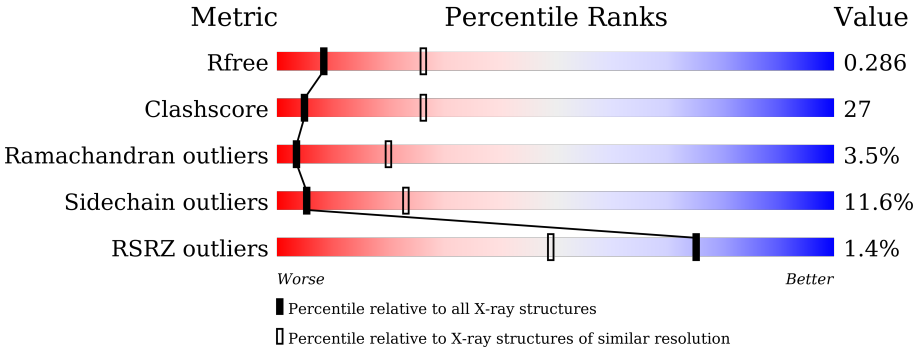
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 3.00 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}	130704	2092 (3.00-3.00)
Clashscore	141614	2416 (3.00-3.00)
Ramachandran outliers	138981	2333 (3.00-3.00)
Sidechain outliers	138945	2336 (3.00-3.00)
RSRZ outliers	127900	1990 (3.00-3.00)

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 <=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	210	.% 46% 43% 9% .
1	B	210	2% 45% 46% 8% .
2	C	17	35% 29% 12% 24%
2	D	17	47% 18% 12% 24%
3	E	8	75% 25%
3	F	8	25% 50% 25%

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
3	FUC	F	8	-	-	-	X

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 3738 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.

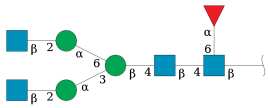
- Molecule 1 is a protein called Immunoglobulin gamma-1 heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	207	1658	1056	279	317	6	0	0	0
1	B	207	1658	1056	279	317	6	0	0	0

- Molecule 2 is a protein called 17-mer peptide (GPDCAYHKGELVWCTFH).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	C	13	112	72	17	21	2	0	0	0
2	D	13	112	72	17	21	2	0	0	0

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



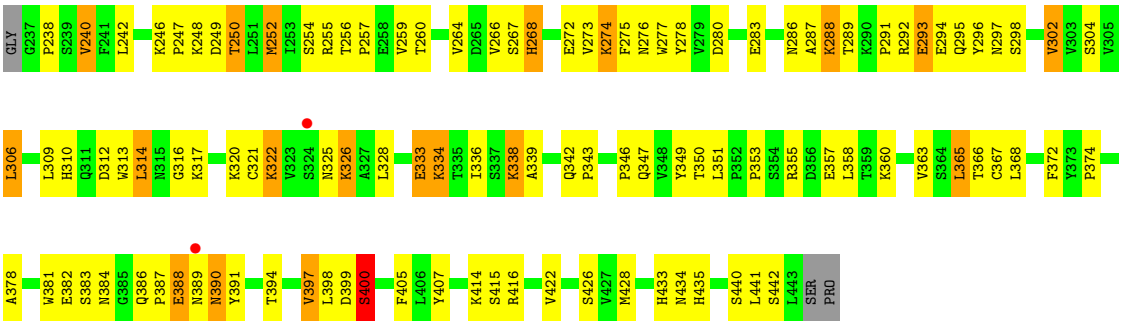
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
3	E	8	99	56	4	39	0	0	0
3	F	8	99	56	4	39	0	0	0

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.

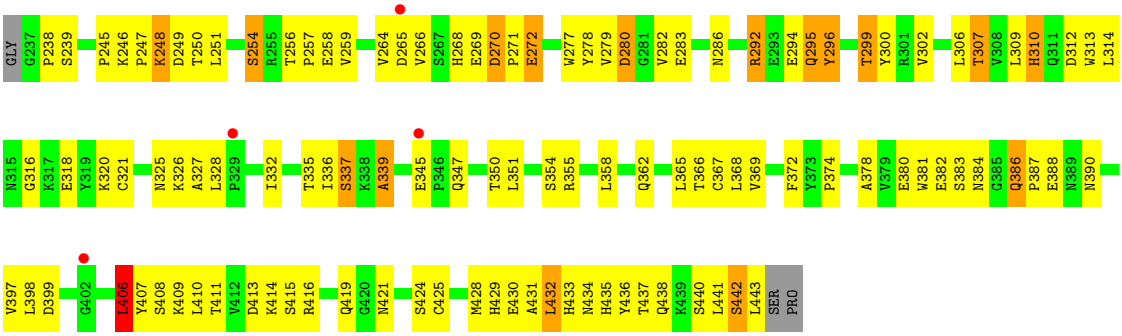
• Molecule 1: Immunoglobulin gamma-1 heavy chain

Chain A:



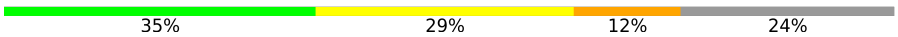
• Molecule 1: Immunoglobulin gamma-1 heavy chain

Chain B:



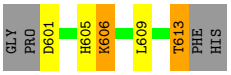
• Molecule 2: 17-mer peptide (GPDCAYHKGELVWCTFH)

Chain C:



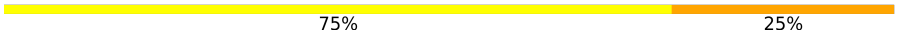
• Molecule 2: 17-mer peptide (GPDCAYHKGELVWCTFH)

Chain D:



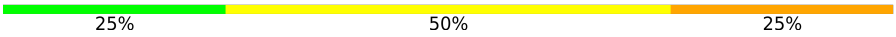
• Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain E:



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

Chain F:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	61.08Å 66.99Å 78.34Å 90.00° 106.41° 90.00°	Depositor
Resolution (Å)	44.10 – 3.00 44.10 – 2.99	Depositor EDS
% Data completeness (in resolution range)	98.1 (44.10-3.00) 93.0 (44.10-2.99)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.59 (at 3.01Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.223 , 0.293 0.227 , 0.286	Depositor DCC
R_{free} test set	610 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å^2)	45.2	Xtriage
Anisotropy	0.494	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 32.4	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.89	EDS
Total number of atoms	3738	wwPDB-VP
Average B, all atoms (Å^2)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 20.83 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 7.9869e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[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: A1LWV, NAG, BMA, FUC, MAN

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	0.67	0/1704	0.77	0/2322
1	B	0.65	0/1704	0.77	1/2322 (0.0%)
2	C	0.76	0/98	0.85	0/132
2	D	0.70	0/98	0.70	0/132
All	All	0.67	0/3604	0.77	1/4908 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	406	LEU	CA-CB-CG	5.29	127.47	115.30

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	1658	0	1623	98	1
1	B	1658	0	1623	98	0
2	C	112	0	78	7	0
2	D	112	0	78	6	0
3	E	99	0	85	4	0
3	F	99	0	85	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	3738	0	3572	200	1

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 (200) 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:333:GLU:OE1	1:A:334:LYS:N	1.66	1.26
1:A:383:SER:O	1:A:386:GLN:HG2	1.44	1.15
1:B:279:VAL:O	1:B:280:ASP:HB2	1.47	1.09
1:B:309:LEU:HB2	1:B:312:ASP:OD1	1.56	1.06
1:A:321:CYS:C	1:A:333:GLU:OE2	1.96	1.03
1:A:247:PRO:O	1:A:250:THR:HG23	1.66	0.95
1:B:419:GLN:O	1:B:419:GLN:HG2	1.68	0.92
1:B:269:GLU:O	1:B:270:ASP:HB2	1.68	0.92
1:B:294:GLU:O	1:B:295:GLN:HB2	1.67	0.92
1:B:358:LEU:O	1:B:414:LYS:HD3	1.70	0.90
1:A:399:ASP:O	1:A:400:SER:HB3	1.71	0.90
1:B:248:LYS:HD3	2:D:606:A1LWV:O1	1.72	0.87
1:B:336:ILE:O	1:B:337:SER:HB3	1.79	0.82
1:A:249:ASP:OD1	1:A:255:ARG:HD3	1.80	0.82
1:A:383:SER:O	1:A:386:GLN:CG	2.27	0.81
1:B:295:GLN:O	1:B:296:TYR:HB2	1.82	0.78
1:A:333:GLU:OE1	1:A:333:GLU:HA	1.84	0.78
1:B:249:ASP:O	1:B:257:PRO:HD3	1.83	0.77
1:B:380:GLU:OE1	2:D:605:HIS:NE2	2.17	0.77
1:B:248:LYS:CD	2:D:606:A1LWV:O1	2.35	0.75
1:A:248:LYS:O	1:A:252:MET:HG3	1.85	0.75
1:A:358:LEU:O	1:A:414:LYS:HE2	1.87	0.74
1:A:322:LYS:N	1:A:333:GLU:OE2	2.20	0.73
1:B:358:LEU:O	1:B:414:LYS:CD	2.37	0.73
1:B:429:HIS:CE1	1:B:431:ALA:H	2.06	0.73
1:A:333:GLU:OE1	1:A:333:GLU:CA	2.37	0.72
1:A:238:PRO:HD2	1:A:328:LEU:HD13	1.69	0.72
1:B:254:SER:OG	2:D:609:LEU:O	2.06	0.71
1:A:316:GLY:HA2	1:A:338:LYS:HD2	1.73	0.70
1:A:336:ILE:O	1:A:336:ILE:HG23	1.90	0.70
1:A:346:PRO:HB3	1:A:372:PHE:HB3	1.72	0.70
1:A:280:ASP:OD2	1:A:317:LYS:HD2	1.92	0.69
1:A:248:LYS:HE3	1:A:428:MET:HE1	1.74	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:295:GLN:O	1:B:296:TYR:CB	2.42	0.68
2:C:601:ASP:N	2:C:601:ASP:OD1	2.23	0.68
1:B:294:GLU:O	1:B:299:THR:O	2.11	0.67
1:B:295:GLN:OE1	3:F:1:NAG:H62	1.93	0.67
1:A:360:LYS:HE2	1:B:347:GLN:OE1	1.94	0.66
1:A:266:VAL:HG21	1:A:302:VAL:HG23	1.75	0.66
1:A:246:LYS:O	1:A:250:THR:HG22	1.95	0.66
1:A:357:GLU:O	1:A:363:VAL:HG12	1.95	0.66
1:B:442:SER:O	1:B:443:LEU:HG	1.96	0.65
1:A:249:ASP:O	1:A:257:PRO:HD3	1.96	0.64
1:A:295:GLN:O	3:E:8:FUC:H61	1.96	0.64
1:B:339:ALA:HB3	1:B:374:PRO:HB3	1.80	0.64
1:B:246:LYS:HD2	1:B:248:LYS:HB2	1.80	0.64
1:A:321:CYS:HB3	1:A:333:GLU:OE2	1.98	0.64
1:A:397:VAL:HG13	1:A:405:PHE:CE1	2.33	0.63
1:B:432:LEU:HD21	1:B:437:THR:HG22	1.80	0.63
3:F:4:MAN:H4	3:F:5:NAG:H83	1.80	0.63
1:B:294:GLU:O	1:B:295:GLN:CB	2.41	0.62
1:B:424:SER:OG	1:B:440:SER:CB	2.48	0.62
1:B:428:MET:HB2	1:B:435:HIS:O	1.99	0.62
1:A:275:PHE:CE2	1:A:304:SER:HB2	2.34	0.62
1:A:339:ALA:HB3	1:A:374:PRO:HB3	1.81	0.62
1:A:297:ASN:O	1:A:298:SER:OG	2.14	0.62
1:B:309:LEU:CB	1:B:312:ASP:OD1	2.42	0.61
1:B:351:LEU:HB2	1:B:366:THR:HB	1.82	0.61
1:B:279:VAL:O	1:B:279:VAL:HG12	2.00	0.61
1:A:322:LYS:HD2	1:A:333:GLU:HG2	1.82	0.61
1:B:269:GLU:O	1:B:270:ASP:CB	2.47	0.60
1:A:310:HIS:O	1:A:314:LEU:HD12	2.01	0.60
1:A:407:TYR:OH	1:B:366:THR:HG23	2.01	0.60
1:B:279:VAL:O	1:B:280:ASP:CB	2.31	0.60
1:B:386:GLN:HG3	1:B:387:PRO:HD2	1.84	0.60
1:A:389:ASN:O	1:A:391:TYR:N	2.33	0.60
1:B:442:SER:OG	1:B:443:LEU:N	2.33	0.59
1:A:333:GLU:OE1	1:A:333:GLU:C	2.38	0.58
1:A:399:ASP:O	1:A:400:SER:CB	2.46	0.58
3:F:4:MAN:O2	3:F:5:NAG:H83	2.04	0.58
1:B:367:CYS:HB2	1:B:381:TRP:CZ2	2.39	0.58
1:A:433:HIS:O	1:A:434:ASN:HB2	2.04	0.57
1:A:383:SER:O	1:A:384:ASN:C	2.42	0.57
1:A:273:VAL:O	1:A:274:LYS:HD2	2.04	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:247:PRO:O	1:A:250:THR:CG2	2.49	0.56
1:B:428:MET:HE2	1:B:436:TYR:HD2	1.69	0.56
1:B:247:PRO:O	1:B:251:LEU:HD12	2.06	0.55
1:B:278:TYR:HB2	1:B:320:LYS:HB3	1.88	0.55
1:A:242:LEU:HB2	1:A:334:LYS:NZ	2.21	0.55
1:A:254:SER:OG	2:C:609:LEU:O	2.16	0.55
1:A:287:ALA:HB2	1:A:306:LEU:HD12	1.89	0.55
1:B:382:GLU:OE2	1:B:438:GLN:OE1	2.24	0.55
1:B:270:ASP:H	1:B:271:PRO:HD3	1.71	0.55
1:A:275:PHE:O	1:A:276:ASN:OD1	2.25	0.54
1:B:362:GLN:HA	1:B:413:ASP:HA	1.89	0.54
1:A:367:CYS:HB2	1:A:381:TRP:CZ2	2.42	0.54
1:B:238:PRO:HD2	1:B:328:LEU:HD13	1.89	0.54
1:B:250:THR:HB	1:B:314:LEU:HD21	1.90	0.54
1:A:268:HIS:O	1:A:268:HIS:ND1	2.41	0.53
1:B:245:PRO:HB3	1:B:258:GLU:H	1.73	0.53
1:A:422:VAL:HA	1:A:442:SER:HB3	1.89	0.53
2:D:601:ASP:O	2:D:601:ASP:OD1	2.26	0.53
1:A:310:HIS:O	1:A:314:LEU:CD1	2.56	0.53
1:A:399:ASP:O	1:A:399:ASP:CG	2.48	0.52
1:B:266:VAL:HG21	1:B:302:VAL:HG23	1.92	0.52
1:B:309:LEU:HD23	1:B:312:ASP:OD1	2.09	0.52
1:A:434:ASN:HB3	2:C:603:ALA:HB2	1.92	0.51
1:B:419:GLN:O	1:B:419:GLN:CG	2.49	0.51
1:A:264:VAL:HG11	3:E:2:NAG:H2	1.91	0.51
1:B:318:GLU:O	1:B:318:GLU:HG2	2.11	0.51
1:B:259:VAL:HG13	1:B:336:ILE:HD11	1.91	0.51
1:A:368:LEU:HD13	1:A:407:TYR:CZ	2.46	0.51
1:A:312:ASP:OD1	1:A:312:ASP:C	2.49	0.51
1:A:394:THR:HG22	1:B:397:VAL:HG21	1.91	0.51
1:B:413:ASP:O	1:B:414:LYS:C	2.48	0.51
1:A:388:GLU:CD	1:A:416:ARG:HH12	2.14	0.50
1:B:424:SER:OG	1:B:440:SER:HB2	2.10	0.50
1:A:256:THR:HG22	1:A:256:THR:O	2.11	0.50
1:A:242:LEU:HB2	1:A:334:LYS:HZ1	1.76	0.50
1:A:295:GLN:O	1:A:296:TYR:HB2	2.12	0.50
1:B:271:PRO:HG2	1:B:272:GLU:OE1	2.12	0.49
1:B:369:VAL:O	1:B:372:PHE:HE2	1.95	0.49
1:A:249:ASP:HA	1:A:255:ARG:HD3	1.94	0.48
1:B:309:LEU:O	1:B:310:HIS:C	2.52	0.48
1:B:350:THR:HB	1:B:441:LEU:HD22	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:272:GLU:O	1:A:325:ASN:ND2	2.46	0.48
1:A:336:ILE:O	1:A:336:ILE:CG2	2.62	0.48
1:B:409:LYS:NZ	1:B:411:THR:OG1	2.47	0.48
1:A:252:MET:HB3	2:C:610:VAL:HA	1.96	0.48
1:B:424:SER:OG	1:B:440:SER:HB3	2.13	0.48
1:B:277:TRP:O	1:B:283:GLU:HA	2.13	0.48
1:B:325:ASN:O	1:B:327:ALA:N	2.47	0.47
1:A:428:MET:HA	1:A:435:HIS:O	2.14	0.47
1:B:264:VAL:O	1:B:265:ASP:HB2	2.15	0.47
1:A:278:TYR:CE2	1:A:283:GLU:HB2	2.50	0.47
1:A:249:ASP:OD1	1:A:255:ARG:CD	2.57	0.47
1:A:321:CYS:O	1:A:333:GLU:OE2	2.32	0.47
1:B:278:TYR:HA	1:B:282:VAL:O	2.15	0.47
1:B:383:SER:O	1:B:384:ASN:C	2.53	0.47
1:B:425:CYS:O	1:B:438:GLN:HA	2.15	0.47
1:A:415:SER:O	1:A:416:ARG:C	2.52	0.47
1:A:350:THR:HB	1:A:441:LEU:HD22	1.96	0.46
1:B:336:ILE:O	1:B:337:SER:CB	2.54	0.46
1:B:419:GLN:OE1	1:B:421:ASN:ND2	2.48	0.46
1:A:277:TRP:O	1:A:283:GLU:HA	2.16	0.46
1:A:321:CYS:CA	1:A:333:GLU:OE2	2.61	0.46
2:C:613:THR:O	2:C:613:THR:HG22	2.15	0.46
1:A:313:TRP:O	1:A:338:LYS:CD	2.64	0.46
1:A:353:PRO:HD3	1:A:365:LEU:HD12	1.97	0.46
1:B:432:LEU:HD21	1:B:437:THR:CG2	2.46	0.46
1:B:238:PRO:HG2	1:B:325:ASN:HD22	1.81	0.45
1:B:264:VAL:HG22	1:B:265:ASP:H	1.81	0.45
1:B:369:VAL:HB	1:B:406:LEU:CD2	2.47	0.45
1:A:293:GLU:O	1:A:293:GLU:HG2	2.17	0.45
1:A:321:CYS:CB	1:A:333:GLU:OE2	2.63	0.45
1:B:266:VAL:HB	1:B:300:TYR:HB2	1.98	0.45
1:B:354:SER:O	1:B:355:ARG:C	2.55	0.45
1:A:248:LYS:HD3	2:C:606:A1LWV:O1	2.17	0.45
1:A:347:GLN:HB3	1:A:349:TYR:CE2	2.51	0.45
1:B:432:LEU:HD12	1:B:433:HIS:H	1.81	0.45
1:B:365:LEU:O	1:B:409:LYS:HA	2.17	0.45
1:B:313:TRP:HD1	1:B:314:LEU:HD23	1.81	0.45
1:A:387:PRO:O	1:A:388:GLU:C	2.54	0.45
1:A:313:TRP:O	1:A:338:LYS:HD3	2.18	0.44
1:B:246:LYS:HG2	1:B:247:PRO:HD2	1.98	0.44
1:A:313:TRP:O	1:A:338:LYS:NZ	2.47	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:434:ASN:OD1	2:D:613:THR:OG1	2.31	0.44
1:A:306:LEU:HD23	1:A:306:LEU:O	2.17	0.44
1:B:443:LEU:HD23	1:B:443:LEU:HA	1.80	0.44
1:A:274:LYS:HD2	1:A:274:LYS:HA	1.89	0.43
1:A:297:ASN:C	1:A:298:SER:HG	2.18	0.43
1:B:398:LEU:HD12	1:B:399:ASP:H	1.83	0.43
1:A:398:LEU:HD23	1:A:399:ASP:H	1.84	0.43
1:B:368:LEU:HD13	1:B:407:TYR:CZ	2.54	0.43
1:A:346:PRO:CB	1:A:372:PHE:HB3	2.45	0.43
2:C:609:LEU:HD11	2:C:612:CYS:SG	2.59	0.43
1:B:292:ARG:HA	1:B:292:ARG:HD3	1.87	0.43
1:A:349:TYR:CD1	1:B:354:SER:HB3	2.54	0.43
1:A:321:CYS:O	1:A:322:LYS:HB2	2.18	0.43
1:B:320:LYS:HB2	1:B:320:LYS:HE3	1.78	0.43
1:A:295:GLN:O	1:A:296:TYR:CB	2.66	0.43
1:B:369:VAL:HB	1:B:406:LEU:HD23	2.00	0.42
1:B:414:LYS:O	1:B:415:SER:C	2.56	0.42
1:B:310:HIS:ND1	1:B:310:HIS:N	2.61	0.42
1:A:320:LYS:CG	1:A:333:GLU:HG3	2.50	0.42
1:A:355:ARG:H	1:A:355:ARG:HG3	1.57	0.42
1:A:248:LYS:HG3	1:A:378:ALA:CB	2.50	0.42
1:A:326:LYS:H	1:A:326:LYS:HG2	1.62	0.42
1:B:429:HIS:ND1	1:B:429:HIS:C	2.73	0.42
1:B:369:VAL:O	1:B:372:PHE:CE2	2.72	0.42
1:A:309:LEU:O	1:A:310:HIS:C	2.58	0.42
1:A:320:LYS:HG2	1:A:321:CYS:O	2.20	0.42
1:B:256:THR:HA	1:B:257:PRO:HD2	1.59	0.42
1:B:306:LEU:HG	1:B:307:THR:O	2.20	0.41
3:E:4:MAN:O2	3:E:5:NAG:H83	2.20	0.41
1:B:328:LEU:HD21	1:B:332:ILE:HD12	2.02	0.41
3:E:1:NAG:H82	3:E:1:NAG:H2	1.94	0.41
1:B:433:HIS:O	1:B:434:ASN:HB2	2.21	0.41
1:B:365:LEU:O	1:B:410:LEU:N	2.51	0.41
1:A:342:GLN:HE21	1:A:342:GLN:HB3	1.72	0.41
1:B:378:ALA:HB3	1:B:428:MET:HG2	2.02	0.41
1:A:246:LYS:HB3	1:A:246:LYS:HE3	1.82	0.41
1:B:264:VAL:HG21	3:F:1:NAG:O4	2.21	0.41
1:B:381:TRP:HE1	1:B:408:SER:CB	2.34	0.41
1:A:240:VAL:HG22	1:A:334:LYS:HE2	2.02	0.40
1:B:325:ASN:O	1:B:326:LYS:C	2.59	0.40
1:A:288:LYS:HG3	1:A:289:THR:N	2.34	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:334:LYS:HB2	1:A:334:LYS:HE3	1.66	0.40
1:A:351:LEU:HB2	1:A:366:THR:HB	2.03	0.40
1:B:271:PRO:HD2	1:B:272:GLU:OE1	2.22	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:294:GLU:N	1:A:342:GLN:OE1[2_647]	2.02	0.18

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	205/210 (98%)	180 (88%)	20 (10%)	5 (2%)	6	29
1	B	205/210 (98%)	164 (80%)	31 (15%)	10 (5%)	2	13
2	C	10/17 (59%)	7 (70%)	3 (30%)	0	100	100
2	D	10/17 (59%)	9 (90%)	1 (10%)	0	100	100
All	All	430/454 (95%)	360 (84%)	55 (13%)	15 (4%)	3	20

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	267	SER
1	B	270	ASP
1	B	280	ASP
1	B	337	SER
1	B	430	GLU
1	A	322	LYS
1	A	388	GLU

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Mol	Chain	Res	Type
1	A	390	ASN
1	B	295	GLN
1	B	296	TYR
1	A	400	SER
1	B	416	ARG
1	B	442	SER
1	B	339	ALA
1	B	316	GLY

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	193/195 (99%)	166 (86%)	27 (14%)	3	16
1	B	193/195 (99%)	175 (91%)	18 (9%)	9	33
2	C	10/13 (77%)	9 (90%)	1 (10%)	7	29
2	D	10/13 (77%)	9 (90%)	1 (10%)	7	29
All	All	406/416 (98%)	359 (88%)	47 (12%)	5	23

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	240	VAL
1	A	250	THR
1	A	252	MET
1	A	259	VAL
1	A	260	THR
1	A	268	HIS
1	A	274	LYS
1	A	286	ASN
1	A	288	LYS
1	A	291	PRO
1	A	292	ARG
1	A	293	GLU
1	A	302	VAL

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Mol	Chain	Res	Type
1	A	306	LEU
1	A	314	LEU
1	A	326	LYS
1	A	333	GLU
1	A	334	LYS
1	A	338	LYS
1	A	343	PRO
1	A	365	LEU
1	A	382	GLU
1	A	390	ASN
1	A	397	VAL
1	A	400	SER
1	A	426	SER
1	A	440	SER
2	C	601	ASP
1	B	239	SER
1	B	248	LYS
1	B	254	SER
1	B	268	HIS
1	B	272	GLU
1	B	286	ASN
1	B	292	ARG
1	B	299	THR
1	B	307	THR
1	B	310	HIS
1	B	321	CYS
1	B	335	THR
1	B	345	GLU
1	B	386	GLN
1	B	388	GLU
1	B	390	ASN
1	B	406	LEU
1	B	432	LEU
2	D	613	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	285	HIS
1	A	311	GLN
1	A	315	ASN
1	A	342	GLN

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Mol	Chain	Res	Type
1	A	386	GLN
1	A	434	ASN
1	B	325	ASN
1	B	433	HIS
1	B	435	HIS
1	B	438	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues 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	A1LWV	D	606	1,2	14,15,16	1.71	2 (14%)	11,16,18	2.36	3 (27%)
2	A1LWV	C	606	1,2	14,15,16	1.67	2 (14%)	11,16,18	2.47	3 (27%)

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	A1LWV	D	606	1,2	-	3/13/15/17	-
2	A1LWV	C	606	1,2	-	9/13/15/17	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	606	A1LWV	C5-NZ	5.27	1.45	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	606	A1LWV	C5-NZ	5.18	1.45	1.33
2	C	606	A1LWV	O3-C5	2.08	1.27	1.23
2	D	606	A1LWV	O3-C5	2.04	1.27	1.23

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	C	606	A1LWV	C4-C5-NZ	5.12	125.04	116.42
2	D	606	A1LWV	C4-C5-NZ	4.79	124.49	116.42
2	D	606	A1LWV	O3-C5-NZ	-4.28	114.93	123.01
2	C	606	A1LWV	O3-C5-NZ	-4.23	115.04	123.01
2	C	606	A1LWV	CE-NZ-C5	-4.20	115.04	122.84
2	D	606	A1LWV	CE-NZ-C5	-4.19	115.05	122.84

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	606	A1LWV	O-C-CA-CB
2	C	606	A1LWV	N-CA-CB-CG
2	C	606	A1LWV	C-CA-CB-CG
2	C	606	A1LWV	C4-C5-NZ-CE
2	D	606	A1LWV	C2-C3-C4-C5
2	C	606	A1LWV	O3-C5-NZ-CE
2	C	606	A1LWV	CG-CD-CE-NZ
2	C	606	A1LWV	CE-CD-CG-CB
2	C	606	A1LWV	C2-C3-C4-C5
2	C	606	A1LWV	C1-C2-C3-C4
2	D	606	A1LWV	CA-CB-CG-CD
2	D	606	A1LWV	CG-CD-CE-NZ

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	606	A1LWV	2	0
2	C	606	A1LWV	1	0

5.5 Carbohydrates (i)

16 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	1,3	14,14,15	0.35	0	17,19,21	0.66	0
3	NAG	E	2	3	14,14,15	0.31	0	17,19,21	0.50	0
3	BMA	E	3	3	11,11,12	1.14	1 (9%)	15,15,17	1.57	4 (26%)
3	MAN	E	4	3	11,11,12	0.95	1 (9%)	15,15,17	1.38	2 (13%)
3	NAG	E	5	3	14,14,15	0.48	0	17,19,21	0.48	0
3	MAN	E	6	3	11,11,12	1.30	2 (18%)	15,15,17	1.68	3 (20%)
3	NAG	E	7	3	14,14,15	0.32	0	17,19,21	0.90	1 (5%)
3	FUC	E	8	3	10,10,11	1.48	2 (20%)	14,14,16	2.05	6 (42%)
3	NAG	F	1	1,3	14,14,15	0.30	0	17,19,21	0.85	1 (5%)
3	NAG	F	2	3	14,14,15	0.33	0	17,19,21	0.75	0
3	BMA	F	3	3	11,11,12	1.39	1 (9%)	15,15,17	1.25	2 (13%)
3	MAN	F	4	3	11,11,12	0.71	0	15,15,17	1.35	3 (20%)
3	NAG	F	5	3	14,14,15	0.35	0	17,19,21	0.53	0
3	MAN	F	6	3	11,11,12	1.28	1 (9%)	15,15,17	1.56	1 (6%)
3	NAG	F	7	3	14,14,15	0.35	0	17,19,21	0.48	0
3	FUC	F	8	3	10,10,11	3.30	7 (70%)	14,14,16	1.93	4 (28%)

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	1,3	-	2/6/23/26	0/1/1/1
3	NAG	E	2	3	-	2/6/23/26	0/1/1/1
3	BMA	E	3	3	-	0/2/19/22	0/1/1/1
3	MAN	E	4	3	-	0/2/19/22	0/1/1/1
3	NAG	E	5	3	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	MAN	E	6	3	-	1/2/19/22	0/1/1/1
3	NAG	E	7	3	-	4/6/23/26	0/1/1/1
3	FUC	E	8	3	-	-	0/1/1/1
3	NAG	F	1	1,3	-	4/6/23/26	0/1/1/1
3	NAG	F	2	3	-	2/6/23/26	0/1/1/1
3	BMA	F	3	3	-	0/2/19/22	0/1/1/1
3	MAN	F	4	3	-	2/2/19/22	0/1/1/1
3	NAG	F	5	3	-	4/6/23/26	0/1/1/1
3	MAN	F	6	3	-	2/2/19/22	0/1/1/1
3	NAG	F	7	3	-	2/6/23/26	0/1/1/1
3	FUC	F	8	3	-	-	0/1/1/1

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	F	8	FUC	C1-C2	5.65	1.65	1.52
3	F	8	FUC	O5-C1	-5.06	1.35	1.43
3	F	8	FUC	C2-C3	4.85	1.59	1.52
3	F	3	BMA	C2-C3	3.44	1.57	1.52
3	F	8	FUC	C4-C5	3.33	1.60	1.52
3	E	8	FUC	O5-C5	2.95	1.49	1.43
3	F	8	FUC	O2-C2	2.70	1.49	1.43
3	F	6	MAN	C1-C2	2.56	1.58	1.52
3	E	6	MAN	O5-C1	-2.49	1.39	1.43
3	E	3	BMA	C4-C5	-2.33	1.48	1.53
3	E	6	MAN	C4-C5	2.13	1.57	1.53
3	F	8	FUC	O5-C5	2.08	1.48	1.43
3	E	8	FUC	C6-C5	2.05	1.56	1.51
3	E	4	MAN	O4-C4	-2.04	1.38	1.43
3	F	8	FUC	O3-C3	2.04	1.47	1.43

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	F	6	MAN	O2-C2-C3	-4.84	100.43	110.14
3	E	6	MAN	C3-C4-C5	-3.95	103.19	110.24
3	F	8	FUC	O2-C2-C1	3.85	117.04	109.15
3	E	8	FUC	C1-O5-C5	3.50	120.70	112.78
3	F	8	FUC	C1-C2-C3	3.23	113.63	109.67
3	E	8	FUC	C2-C3-C4	-3.13	105.48	110.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	E	4	MAN	C1-O5-C5	3.03	116.30	112.19
3	E	8	FUC	O2-C2-C1	3.02	115.34	109.15
3	E	4	MAN	O2-C2-C3	-2.98	104.16	110.14
3	F	4	MAN	C1-O5-C5	2.97	116.22	112.19
3	E	7	NAG	C1-O5-C5	2.95	116.18	112.19
3	F	8	FUC	O5-C1-C2	-2.93	106.25	110.77
3	E	8	FUC	C1-C2-C3	-2.91	106.09	109.67
3	E	6	MAN	C6-C5-C4	2.90	119.80	113.00
3	F	8	FUC	O5-C5-C4	2.77	114.48	109.52
3	F	1	NAG	C1-O5-C5	2.76	115.93	112.19
3	F	3	BMA	C1-C2-C3	2.72	113.01	109.67
3	E	8	FUC	O5-C5-C4	2.67	114.32	109.52
3	F	4	MAN	O2-C2-C3	-2.45	105.24	110.14
3	E	3	BMA	O5-C5-C4	-2.42	104.94	110.83
3	E	3	BMA	C1-C2-C3	2.38	112.59	109.67
3	E	8	FUC	O5-C1-C2	2.28	114.30	110.77
3	E	3	BMA	C1-O5-C5	2.22	115.20	112.19
3	F	4	MAN	O5-C1-C2	2.18	114.14	110.77
3	E	3	BMA	O2-C2-C3	-2.11	105.91	110.14
3	E	6	MAN	O4-C4-C5	2.10	114.52	109.30
3	F	3	BMA	C2-C3-C4	2.04	114.43	110.89

There are no chirality outliers.

All (27) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	E	2	NAG	O5-C5-C6-O6
3	F	5	NAG	O5-C5-C6-O6
3	E	2	NAG	C4-C5-C6-O6
3	E	7	NAG	O5-C5-C6-O6
3	F	5	NAG	C4-C5-C6-O6
3	E	1	NAG	C8-C7-N2-C2
3	E	1	NAG	O7-C7-N2-C2
3	E	5	NAG	C8-C7-N2-C2
3	E	5	NAG	O7-C7-N2-C2
3	E	7	NAG	C8-C7-N2-C2
3	E	7	NAG	O7-C7-N2-C2
3	F	1	NAG	C8-C7-N2-C2
3	F	1	NAG	O7-C7-N2-C2
3	F	5	NAG	C8-C7-N2-C2
3	F	5	NAG	O7-C7-N2-C2
3	F	7	NAG	C8-C7-N2-C2

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Mol	Chain	Res	Type	Atoms
3	F	7	NAG	O7-C7-N2-C2
3	F	2	NAG	O5-C5-C6-O6
3	E	7	NAG	C4-C5-C6-O6
3	F	1	NAG	O5-C5-C6-O6
3	F	1	NAG	C4-C5-C6-O6
3	F	6	MAN	O5-C5-C6-O6
3	F	2	NAG	C4-C5-C6-O6
3	F	4	MAN	O5-C5-C6-O6
3	E	6	MAN	O5-C5-C6-O6
3	F	6	MAN	C4-C5-C6-O6
3	F	4	MAN	C4-C5-C6-O6

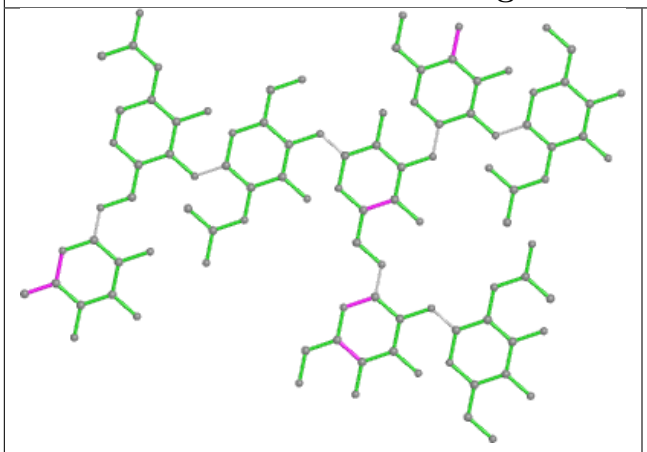
There are no ring outliers.

8 monomers are involved in 8 short contacts:

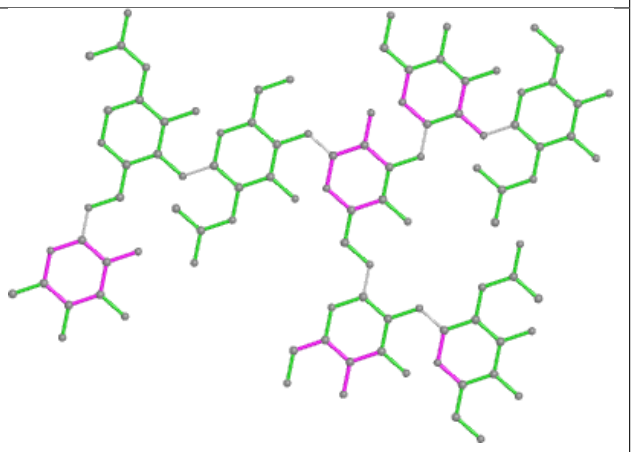
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	E	8	FUC	1	0
3	E	4	MAN	1	0
3	E	5	NAG	1	0
3	F	5	NAG	2	0
3	F	4	MAN	2	0
3	F	1	NAG	2	0
3	E	2	NAG	1	0
3	E	1	NAG	1	0

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

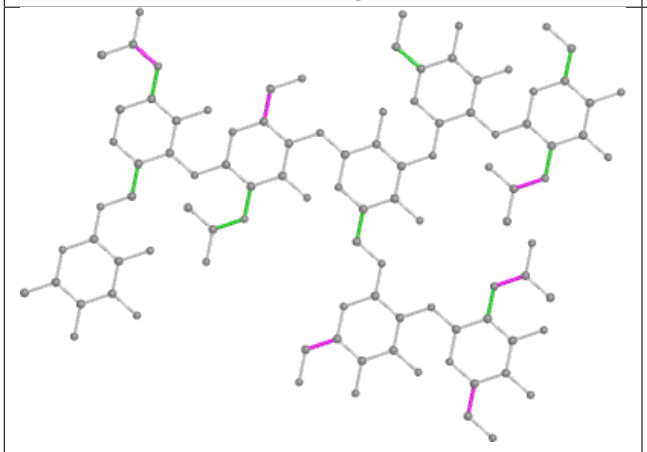
Oligosaccharide Chain E



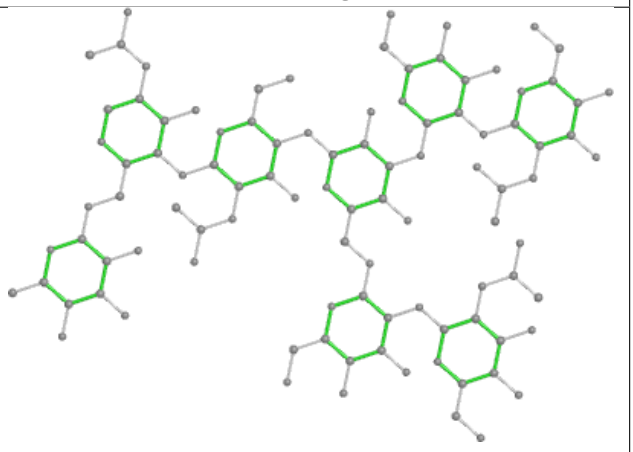
Bond lengths



Bond angles

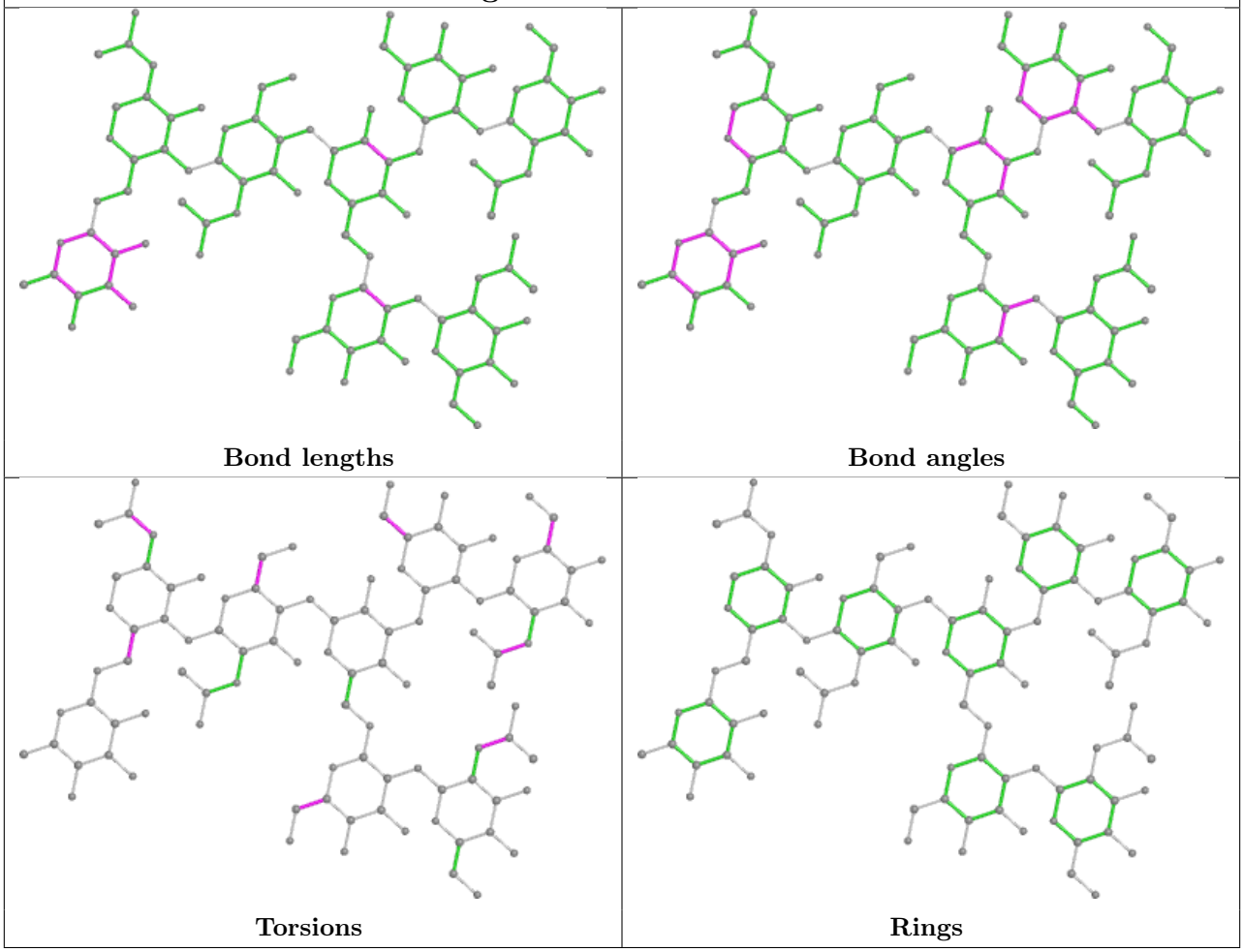


Torsions



Rings

Oligosaccharide Chain F



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	207/210 (98%)	0.06	2 (0%)	82	59	25, 41, 56, 76	0
1	B	207/210 (98%)	0.14	4 (1%)	66	37	25, 43, 59, 72	0
2	C	12/17 (70%)	0.19	0	100	100	38, 45, 58, 62	0
2	D	12/17 (70%)	0.08	0	100	100	38, 42, 60, 61	0
All	All	438/454 (96%)	0.10	6 (1%)	75	49	25, 42, 59, 76	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	389	ASN	3.2
1	B	329	PRO	2.7
1	B	265	ASP	2.6
1	B	345	GLU	2.2
1	B	402	GLY	2.1
1	A	324	SER	2.0

6.2 Non-standard residues in protein, DNA, RNA chains (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	A1LWV	C	606	16/17	0.85	0.35	51,56,64,64	0
2	A1LWV	D	606	16/17	0.87	0.40	34,51,75,77	0

6.3 Carbohydrates (i)

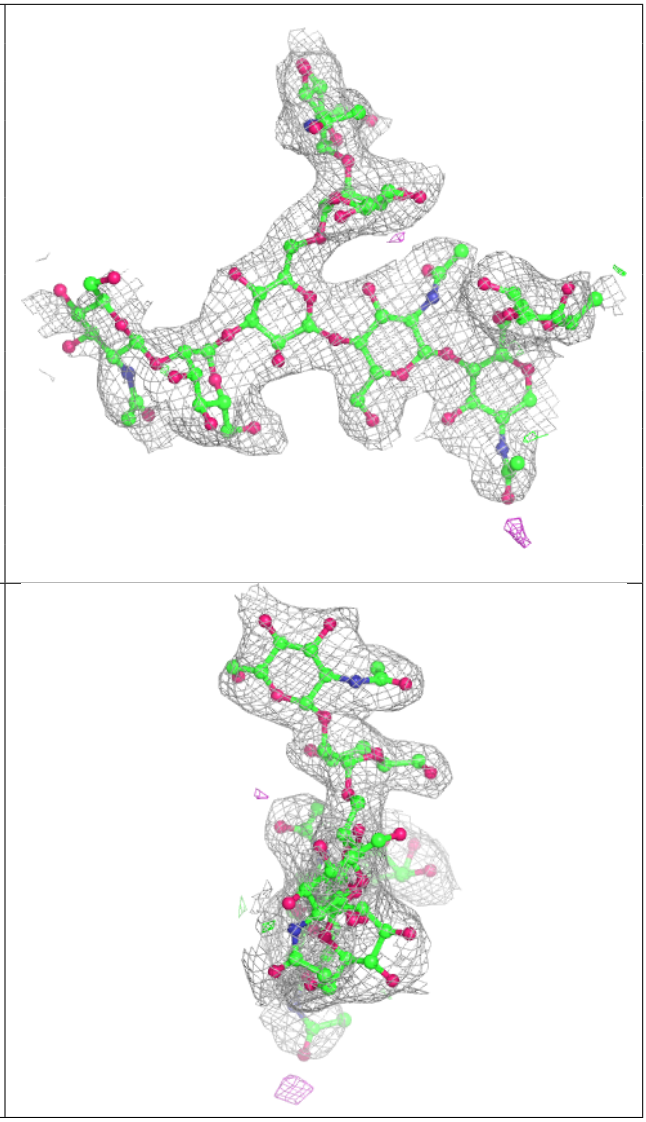
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

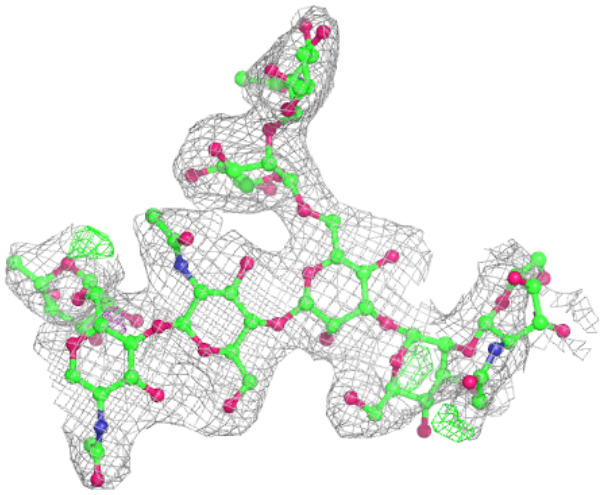
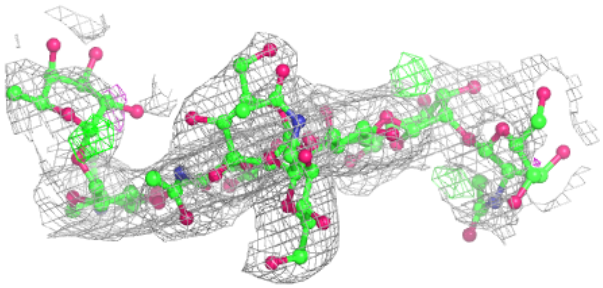
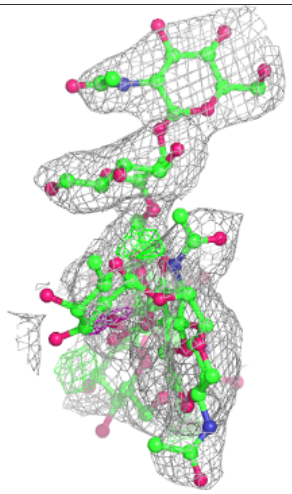
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	FUC	F	8	10/11	0.57	0.46	59,78,83,87	0
3	NAG	F	5	14/15	0.64	0.34	87,107,115,116	0
3	NAG	E	5	14/15	0.70	0.32	71,98,110,111	0
3	MAN	F	4	11/12	0.80	0.19	47,70,93,97	0
3	NAG	F	1	14/15	0.88	0.16	48,52,62,65	0
3	MAN	F	6	11/12	0.89	0.15	47,54,61,64	0
3	FUC	E	8	10/11	0.90	0.16	34,41,52,52	0
3	NAG	F	2	14/15	0.90	0.17	40,48,57,58	0
3	NAG	F	7	14/15	0.91	0.18	41,52,62,77	0
3	BMA	F	3	11/12	0.91	0.13	45,49,53,61	0
3	MAN	E	4	11/12	0.92	0.11	36,61,69,76	0
3	MAN	E	6	11/12	0.92	0.16	40,51,55,56	0
3	NAG	E	7	14/15	0.92	0.16	37,41,48,49	0
3	NAG	E	1	14/15	0.94	0.21	36,45,51,57	0
3	BMA	E	3	11/12	0.95	0.10	39,46,52,55	0
3	NAG	E	2	14/15	0.96	0.14	32,38,42,44	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 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 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.