



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 11, 2021 – 02:48 AM EDT

PDB ID : 3BPO
Title : Crystal structure of the IL13-IL4R-IL13Ra ternary complex
Authors : Garcia, K.C.
Deposited on : 2007-12-18
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 ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) 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 : 2.23.2
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) : 2.23.2

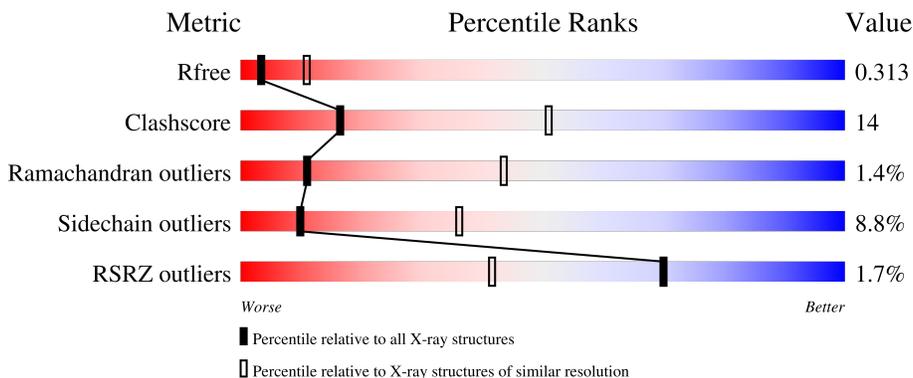
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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 $\leq 5\%$. 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	127	
2	B	205	
3	C	314	
4	D	2	

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	NAG	D	1	X	-	-	X
5	NAG	B	901	X	-	-	-

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 4574 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 Interleukin 13.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	98	725	464	122	132	7	0	0	0

- Molecule 2 is a protein called Interleukin-4 receptor alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	198	1532	981	255	287	9	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-2	ALA	-	expression tag	UNP P24394
B	-1	ASP	-	expression tag	UNP P24394
B	0	PRO	-	expression tag	UNP P24394
B	1	PHE	-	expression tag	UNP P24394
B	28	GLN	ASN	engineered mutation	UNP P24394
B	73	GLN	ASN	engineered mutation	UNP P24394
B	109	GLN	ASN	engineered mutation	UNP P24394
B	151	GLN	ASN	engineered mutation	UNP P24394

- Molecule 3 is a protein called Interleukin-13 receptor alpha-1 chain.

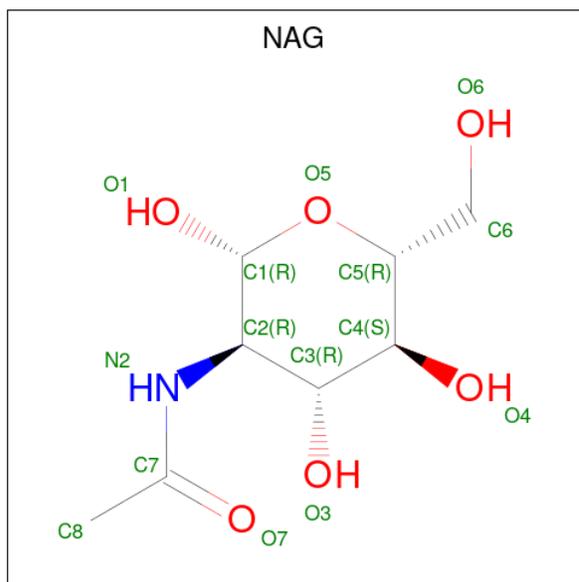
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	C	286	2221	1413	377	414	17	0	0	0

- Molecule 4 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	
			Total	C	N				O
4	D	2	28	16	2	10	0	0	0

- Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: C₈H₁₅NO₆).



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

- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	8	Total O 8 8	0	0
6	B	20	Total O 20 20	0	0
6	C	26	Total O 26 26	0	0

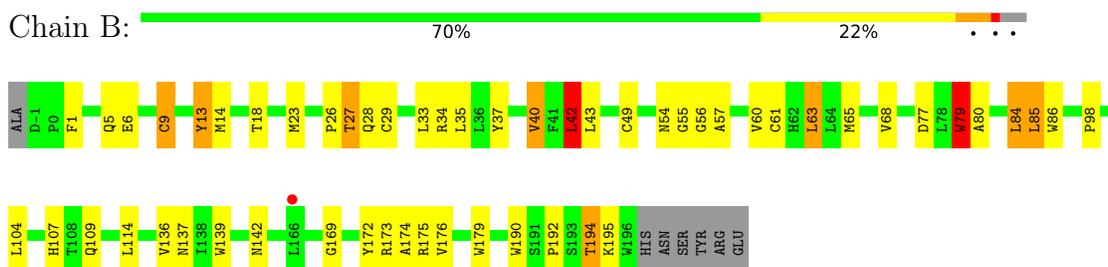
3 Residue-property plots

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

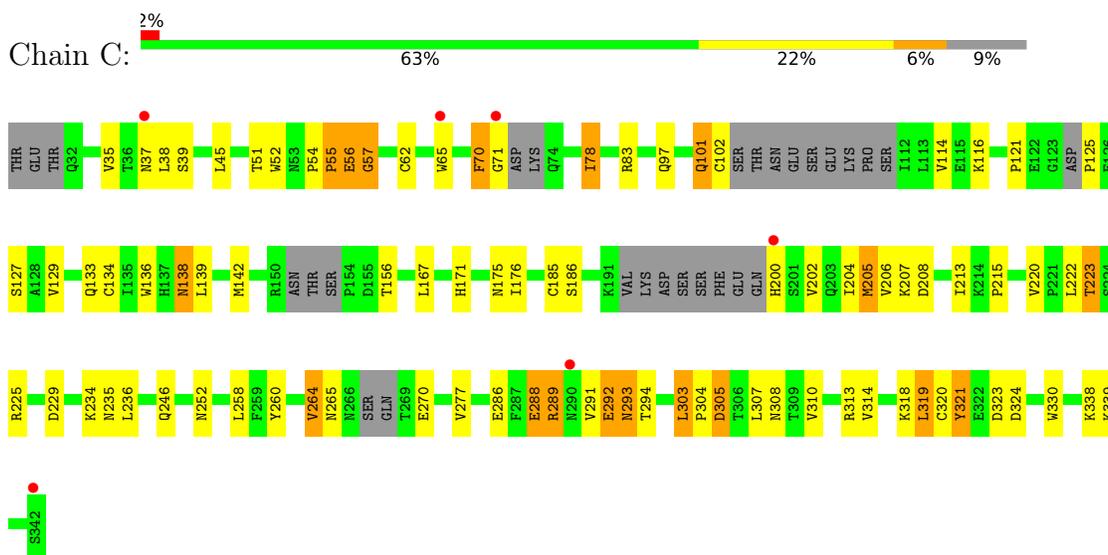
- Molecule 1: Interleukin 13



- Molecule 2: Interleukin-4 receptor alpha chain



- Molecule 3: Interleukin-13 receptor alpha-1 chain



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

Chain D: 

MAG1
MAG2

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	211.51Å 58.17Å 64.24Å 90.00° 100.51° 90.00°	Depositor
Resolution (Å)	30.00 – 3.00 34.66 – 2.90	Depositor EDS
% Data completeness (in resolution range)	99.1 (30.00-3.00) 98.7 (34.66-2.90)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.30 (at 2.90Å)	Xtrriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.253 , 0.312 0.252 , 0.313	Depositor DCC
R_{free} test set	868 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å ²)	64.5	Xtrriage
Anisotropy	0.473	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 63.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	4574	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.25% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	A	0.32	0/735	0.46	0/991
2	B	0.31	0/1582	0.48	0/2180
3	C	0.38	0/2278	0.55	2/3105 (0.1%)
All	All	0.35	0/4595	0.51	2/6276 (0.0%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	4
2	B	0	7
3	C	0	3
All	All	0	14

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
3	C	125	PRO	N-CA-CB	5.89	110.37	103.30
3	C	37	ASN	N-CA-C	5.00	124.51	111.00

There are no chirality outliers.

All (14) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	21	THR	Peptide
1	A	27	PRO	Peptide

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Mol	Chain	Res	Type	Group
1	A	43	MET	Peptide
1	A	54	VAL	Peptide
2	B	40	VAL	Peptide
2	B	42	LEU	Peptide
2	B	43	LEU	Peptide
2	B	79	TRP	Peptide
2	B	80	ALA	Peptide
2	B	84	LEU	Peptide
2	B	85	LEU	Peptide
3	C	288	GLU	Peptide
3	C	54	PRO	Peptide
3	C	57	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	725	0	711	33	1
2	B	1532	0	1400	30	0
3	C	2221	0	2045	60	0
4	D	28	0	25	4	0
5	B	14	0	13	2	0
6	A	8	0	0	0	0
6	B	20	0	0	3	0
6	C	26	0	0	3	0
All	All	4574	0	4194	125	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 14.

All (125) 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)
2:B:84:LEU:HB3	2:B:85:LEU:HA	1.18	1.15
3:C:57:GLY:HA3	6:C:364:HOH:O	1.47	1.13
3:C:35:VAL:HG12	3:C:55:PRO:HD3	1.51	0.92

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:3:PRO:HB2	4:D:1:NAG:O6	1.70	0.91
3:C:200:HIS:O	3:C:222:LEU:HD22	1.69	0.91
2:B:84:LEU:CB	2:B:85:LEU:HA	2.00	0.91
2:B:85:LEU:CB	6:B:919:HOH:O	2.23	0.86
3:C:127:SER:O	3:C:213:ILE:HB	1.73	0.86
2:B:6:GLU:HG3	2:B:86:TRP:CZ2	2.14	0.82
2:B:84:LEU:HB3	2:B:85:LEU:CA	2.05	0.82
1:A:43:MET:HB3	1:A:45:CYS:H	1.45	0.81
1:A:33:MET:HB3	1:A:89:LYS:HG2	1.64	0.78
3:C:38:LEU:HD12	3:C:52:TRP:HB3	1.67	0.77
3:C:200:HIS:O	3:C:222:LEU:CD2	2.32	0.77
2:B:34:ARG:N	2:B:79:TRP:O	2.18	0.75
1:A:43:MET:HB3	1:A:45:CYS:N	2.01	0.74
1:A:3:PRO:O	4:D:1:NAG:O6	2.05	0.70
2:B:35:LEU:HB3	2:B:49:CYS:HB3	1.74	0.68
3:C:142:MET:HE3	3:C:202:VAL:HG11	1.75	0.67
2:B:104:LEU:HD12	2:B:176:VAL:HG22	1.75	0.67
1:A:89:LYS:HB3	3:C:78:ILE:HG23	1.77	0.65
3:C:223:THR:HG22	3:C:321:TYR:HA	1.77	0.65
3:C:260:TYR:HB2	3:C:277:VAL:CG1	2.28	0.64
3:C:35:VAL:CG1	3:C:55:PRO:HD3	2.27	0.63
3:C:70:PHE:O	6:C:354:HOH:O	2.15	0.63
5:B:901:NAG:H83	6:B:918:HOH:O	1.99	0.61
3:C:292:GLU:N	3:C:294:THR:HG22	2.16	0.61
1:A:18:VAL:HG23	1:A:100:LEU:HD11	1.83	0.61
2:B:13:TYR:HA	2:B:68:VAL:HG11	1.83	0.61
2:B:63:LEU:H	2:B:63:LEU:HD12	1.66	0.61
1:A:17:LEU:HB2	1:A:100:LEU:HD13	1.83	0.60
5:B:901:NAG:C8	6:B:918:HOH:O	2.50	0.60
2:B:173:ARG:HA	2:B:194:THR:O	2.02	0.59
3:C:264:VAL:C	3:C:265:ASN:HD22	2.06	0.59
3:C:121:PRO:O	3:C:215:PRO:HD2	2.03	0.59
1:A:3:PRO:CB	4:D:1:NAG:O6	2.50	0.59
1:A:17:LEU:HD22	1:A:96:VAL:HG13	1.85	0.58
3:C:260:TYR:HB2	3:C:277:VAL:HG12	1.86	0.58
2:B:77:ASP:HA	2:B:86:TRP:O	2.03	0.57
3:C:305:ASP:O	3:C:338:LYS:HB3	2.04	0.57
3:C:246:GLN:HE22	3:C:289:ARG:H	1.51	0.56
1:A:17:LEU:O	1:A:20:ILE:HG22	2.06	0.56
3:C:291:VAL:C	3:C:293:ASN:H	2.09	0.56
3:C:318:LYS:CD	3:C:323:ASP:HB2	2.36	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:318:LYS:HD3	3:C:323:ASP:HB2	1.89	0.55
2:B:175:ARG:HD2	2:B:190:TRP:HB3	1.90	0.54
2:B:136:VAL:HA	2:B:176:VAL:HG12	1.89	0.54
3:C:223:THR:O	3:C:321:TYR:HB3	2.07	0.53
1:A:18:VAL:O	1:A:21:THR:O	2.27	0.53
3:C:142:MET:CE	3:C:202:VAL:HG11	2.37	0.53
3:C:127:SER:HB2	3:C:213:ILE:HG22	1.91	0.53
2:B:137:ASN:O	2:B:174:ALA:HA	2.09	0.53
2:B:23:MET:HG3	2:B:27:THR:HG21	1.90	0.52
3:C:246:GLN:OE1	3:C:289:ARG:O	2.26	0.52
2:B:34:ARG:HB2	2:B:79:TRP:HB3	1.91	0.52
2:B:98:PRO:HG3	2:B:179:TRP:HA	1.93	0.51
1:A:51:LEU:C	1:A:53:ASN:H	2.14	0.51
4:D:1:NAG:H61	4:D:2:NAG:N2	2.26	0.51
3:C:138:ASN:ND2	3:C:229:ASP:H	2.09	0.50
3:C:39:SER:HB2	3:C:51:THR:HG22	1.93	0.50
2:B:40:VAL:HG12	2:B:40:VAL:O	2.12	0.50
2:B:142:ASN:H	2:B:142:ASN:HD22	1.60	0.49
3:C:292:GLU:H	3:C:294:THR:HG22	1.76	0.49
1:A:20:ILE:HG12	1:A:28:LEU:HD11	1.95	0.49
3:C:234:LYS:O	3:C:235:ASN:HB3	2.12	0.48
1:A:37:ILE:HG12	1:A:46:ALA:HB1	1.95	0.48
3:C:56:GLU:OE1	3:C:56:GLU:O	2.31	0.48
3:C:134:CYS:SG	3:C:204:ILE:HD13	2.54	0.48
2:B:172:TYR:O	2:B:195:LYS:HA	2.13	0.48
3:C:97:GLN:NE2	6:C:354:HOH:O	2.47	0.48
2:B:26:PRO:HA	2:B:56:GLY:O	2.14	0.48
1:A:28:LEU:O	1:A:29:CYS:HB2	2.14	0.47
1:A:10:LEU:HD11	1:A:103:LEU:HB3	1.97	0.47
3:C:208:ASP:CB	3:C:213:ILE:HG12	2.44	0.47
2:B:63:LEU:HD12	2:B:63:LEU:N	2.29	0.47
1:A:51:LEU:O	1:A:54:VAL:HG23	2.15	0.47
1:A:32:SER:O	1:A:92:VAL:HG22	2.15	0.46
3:C:292:GLU:OE2	3:C:292:GLU:HA	2.15	0.46
1:A:16:GLU:HG3	1:A:66:MET:CE	2.45	0.46
3:C:136:TRP:N	3:C:225:ARG:O	2.46	0.46
1:A:54:VAL:HG12	1:A:55:SER:H	1.81	0.46
3:C:205:MET:HE1	3:C:207:LYS:HB3	1.96	0.46
3:C:265:ASN:O	3:C:308:ASN:HA	2.15	0.46
1:A:43:MET:HG2	1:A:45:CYS:SG	2.57	0.45
2:B:107:HIS:CE1	2:B:109:GLN:HE21	2.33	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:313:ARG:HD3	3:C:330:TRP:CE2	2.52	0.45
3:C:45:LEU:HD12	3:C:45:LEU:HA	1.91	0.45
3:C:291:VAL:O	3:C:293:ASN:N	2.50	0.45
3:C:97:GLN:HA	3:C:114:VAL:O	2.17	0.44
2:B:55:GLY:C	2:B:57:ALA:H	2.20	0.44
1:A:7:SER:O	1:A:11:ARG:HB2	2.17	0.44
3:C:303:LEU:H	3:C:303:LEU:HD12	1.81	0.44
1:A:28:LEU:HD23	1:A:29:CYS:SG	2.58	0.44
2:B:37:TYR:CG	2:B:63:LEU:HD23	2.53	0.43
1:A:51:LEU:C	1:A:53:ASN:N	2.72	0.43
3:C:258:LEU:HD22	3:C:314:VAL:HG22	2.00	0.43
3:C:291:VAL:C	3:C:293:ASN:N	2.72	0.43
1:A:41:ALA:HA	1:A:42:GLY:HA3	1.68	0.43
1:A:54:VAL:HG12	1:A:55:SER:N	2.34	0.43
2:B:54:ASN:HB2	2:B:60:VAL:HG12	2.00	0.43
3:C:175:ASN:HB2	3:C:186:SER:O	2.19	0.43
3:C:129:VAL:HG21	3:C:206:VAL:CG1	2.49	0.42
1:A:83:LEU:HG	1:A:84:HIS:H	1.84	0.42
3:C:264:VAL:HG13	3:C:310:VAL:HG22	1.99	0.42
3:C:70:PHE:HD2	3:C:70:PHE:HA	1.65	0.42
3:C:319:LEU:O	3:C:320:CYS:HB2	2.20	0.42
1:A:16:GLU:HG3	1:A:66:MET:HE1	2.02	0.42
3:C:56:GLU:HA	3:C:57:GLY:HA2	1.77	0.42
3:C:70:PHE:O	3:C:71:GLY:C	2.58	0.42
3:C:307:LEU:HD22	3:C:339:LYS:HG2	2.02	0.42
3:C:313:ARG:HB2	3:C:330:TRP:CZ3	2.55	0.41
2:B:139:TRP:HZ3	2:B:175:ARG:HG2	1.85	0.41
1:A:29:CYS:HB2	1:A:92:VAL:HG21	2.01	0.41
3:C:167:LEU:HD22	3:C:171:HIS:NE2	2.35	0.41
2:B:9:CYS:HA	2:B:18:THR:O	2.20	0.41
3:C:38:LEU:HD23	3:C:116:LYS:HB2	2.02	0.41
3:C:303:LEU:HA	3:C:304:PRO:HD3	1.93	0.41
1:A:65:ARG:HH11	1:A:65:ARG:HG3	1.86	0.41
3:C:288:GLU:HB3	3:C:289:ARG:O	2.20	0.41
1:A:54:VAL:HG11	1:A:57:CYS:HB3	2.02	0.41
3:C:101:GLN:O	3:C:102:CYS:SG	2.79	0.40
2:B:29:CYS:SG	2:B:33:LEU:HD12	2.61	0.40
1:A:60:ILE:O	1:A:64:GLN:HG3	2.22	0.40
3:C:136:TRP:CE2	3:C:139:LEU:HA	2.56	0.40
3:C:202:VAL:CG1	3:C:220:VAL:HB	2.51	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the sym-

metry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:55:SER:CB	1:A:55:SER:CB[2_554]	2.15	0.05

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	90/127 (71%)	73 (81%)	15 (17%)	2 (2%)	6	31
2	B	196/205 (96%)	174 (89%)	19 (10%)	3 (2%)	10	42
3	C	272/314 (87%)	241 (89%)	28 (10%)	3 (1%)	14	50
All	All	558/646 (86%)	488 (88%)	62 (11%)	8 (1%)	11	43

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	C	292	GLU
2	B	42	LEU
2	B	169	GLY
3	C	55	PRO
3	C	138	ASN
1	A	52	ILE
1	A	85	VAL
2	B	192	PRO

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	76/112 (68%)	73 (96%)	3 (4%)	32	69
2	B	161/184 (88%)	147 (91%)	14 (9%)	10	37
3	C	242/299 (81%)	217 (90%)	25 (10%)	7	28
All	All	479/595 (80%)	437 (91%)	42 (9%)	10	36

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	LEU
1	A	43	MET
1	A	92	VAL
2	B	1	PHE
2	B	5	GLN
2	B	9	CYS
2	B	13	TYR
2	B	14	MET
2	B	27	THR
2	B	28	GLN
2	B	42	LEU
2	B	61	CYS
2	B	63	LEU
2	B	65	MET
2	B	79	TRP
2	B	114	LEU
2	B	194	THR
3	C	56	GLU
3	C	62	CYS
3	C	65	TRP
3	C	70	PHE
3	C	78	ILE
3	C	83	ARG
3	C	101	GLN
3	C	133	GLN
3	C	156	THR
3	C	176	ILE
3	C	185	CYS
3	C	205	MET
3	C	223	THR
3	C	236	LEU
3	C	252	ASN
3	C	264	VAL
3	C	270	GLU

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Mol	Chain	Res	Type
3	C	286	GLU
3	C	289	ARG
3	C	293	ASN
3	C	303	LEU
3	C	305	ASP
3	C	319	LEU
3	C	321	TYR
3	C	324	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	19	ASN
2	B	24	ASN
2	B	38	GLN
2	B	95	HIS
2	B	109	GLN
2	B	120	ASN
2	B	142	ASN
2	B	151	GLN
3	C	74	GLN
3	C	138	ASN
3	C	209	ASN
3	C	246	GLN
3	C	265	ASN
3	C	273	ASN
3	C	293	ASN
3	C	308	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](#)

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
4	NAG	D	1	4,2	14,14,15	2.67	4 (28%)	17,19,21	2.63	6 (35%)
4	NAG	D	2	4	14,14,15	0.64	0	17,19,21	0.91	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	D	1	4,2	1/1/5/7	0/6/23/26	0/1/1/1
4	NAG	D	2	4	-	2/6/23/26	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	1	NAG	C2-N2	7.32	1.58	1.46
4	D	1	NAG	O5-C1	4.36	1.50	1.43
4	D	1	NAG	C8-C7	4.27	1.59	1.50
4	D	1	NAG	C7-N2	-2.02	1.27	1.34

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	1	NAG	C2-N2-C7	-6.43	113.75	122.90
4	D	1	NAG	C1-C2-N2	-4.47	102.85	110.49
4	D	1	NAG	O4-C4-C3	-4.38	100.22	110.35
4	D	1	NAG	O7-C7-N2	-3.26	115.96	121.95
4	D	1	NAG	C1-O5-C5	-2.76	108.45	112.19
4	D	1	NAG	O5-C1-C2	-2.37	107.55	111.29

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
4	D	1	NAG	C1

All (2) torsion outliers are listed below:

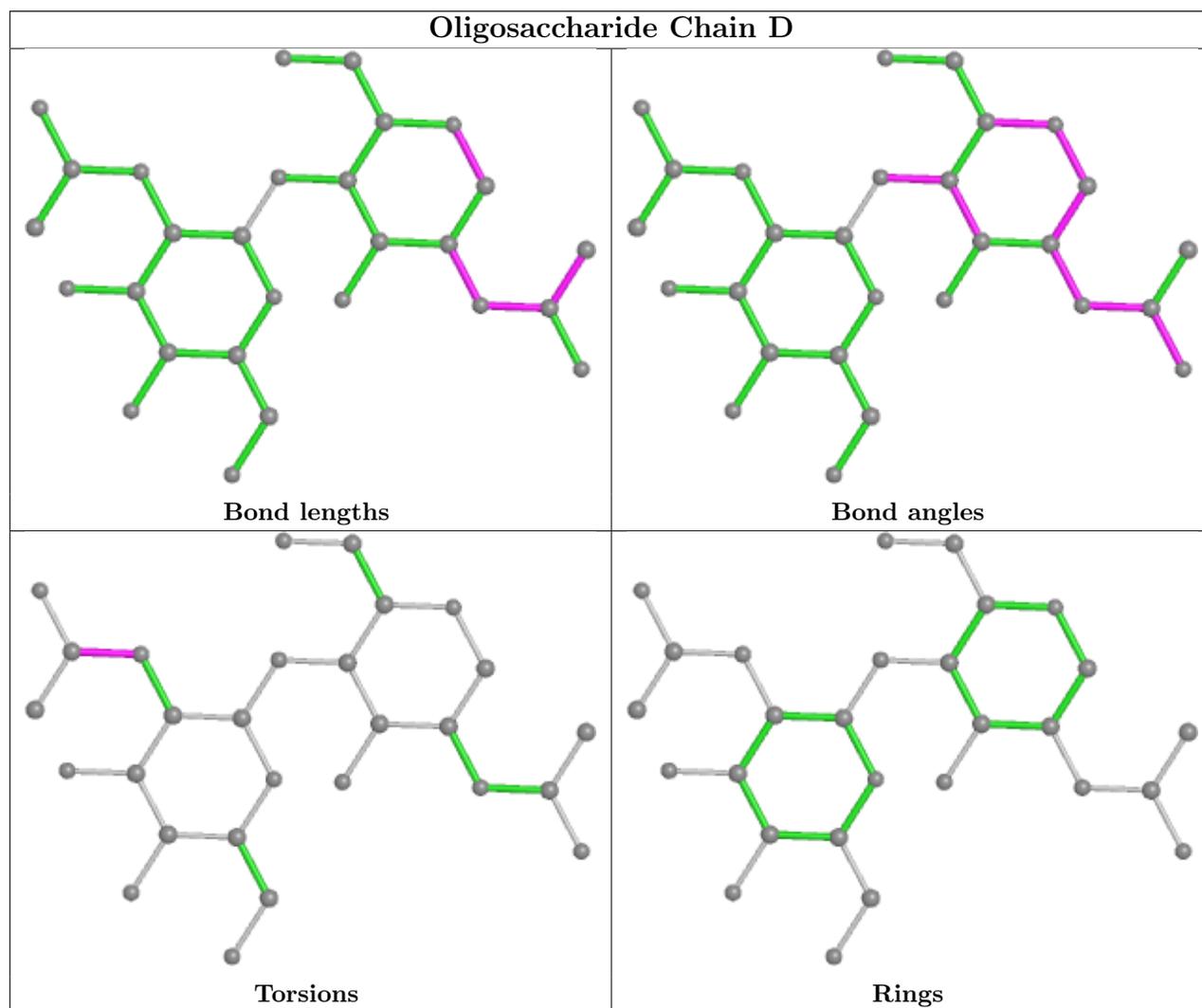
Mol	Chain	Res	Type	Atoms
4	D	2	NAG	C8-C7-N2-C2
4	D	2	NAG	O7-C7-N2-C2

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	D	1	NAG	4	0
4	D	2	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.



5.6 Ligand geometry

1 ligand is 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
5	NAG	B	901	2	14,14,15	2.66	4 (28%)	17,19,21	2.63	6 (35%)

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
5	NAG	B	901	2	1/1/5/7	0/6/23/26	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	901	NAG	C2-N2	7.33	1.58	1.46
5	B	901	NAG	O5-C1	4.34	1.50	1.43
5	B	901	NAG	C8-C7	4.19	1.59	1.50
5	B	901	NAG	C7-N2	-2.03	1.27	1.34

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	901	NAG	C2-N2-C7	-6.36	113.84	122.90
5	B	901	NAG	C1-C2-N2	-4.42	102.93	110.49
5	B	901	NAG	O4-C4-C3	-4.41	100.16	110.35
5	B	901	NAG	O7-C7-N2	-3.31	115.86	121.95
5	B	901	NAG	C1-O5-C5	-2.75	108.46	112.19
5	B	901	NAG	O5-C1-C2	-2.35	107.58	111.29

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
5	B	901	NAG	C1

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	901	NAG	2	0

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	98/127 (77%)	-0.00	3 (3%) 49 21	23, 51, 63, 64	0
2	B	198/205 (96%)	-0.10	1 (0%) 91 75	20, 48, 61, 64	0
3	C	286/314 (91%)	-0.15	6 (2%) 63 34	20, 31, 63, 66	0
All	All	582/646 (90%)	-0.11	10 (1%) 70 41	20, 45, 62, 66	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	C	290	ASN	3.5
3	C	71	GLY	3.3
3	C	342	SER	2.6
3	C	65	TRP	2.4
2	B	166	LEU	2.4
3	C	200	HIS	2.3
3	C	37	ASN	2.2
1	A	5	PRO	2.2
1	A	26	ALA	2.2
1	A	3	PRO	2.1

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

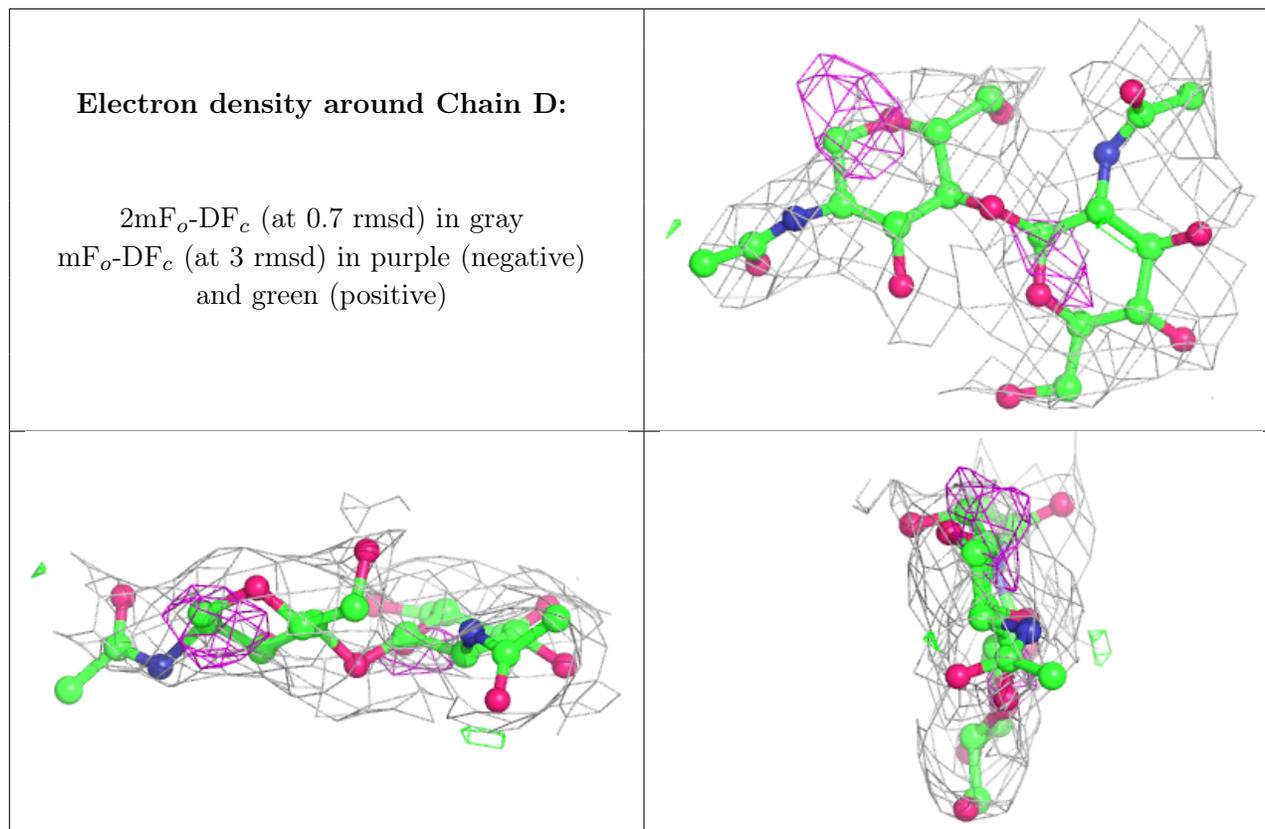
There are no non-standard protein/DNA/RNA residues in this entry.

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(\AA^2)	Q<0.9
4	NAG	D	2	14/15	0.75	0.33	81,83,84,84	0
4	NAG	D	1	14/15	0.79	0.42	70,74,75,77	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 [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(\AA^2)	Q<0.9
5	NAG	B	901	14/15	0.75	0.26	56,56,57,64	0

6.5 Other polymers [i](#)

There are no such residues in this entry.