



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 7, 2020 – 01:08 AM BST

PDB ID : 4XLW
Title : Complex of Notch1 (EGF11-13) bound to Delta-like 4 (N-EGF2)
Authors : Luca, V.C.; Jude, K.M.; Garcia, K.C.
Deposited on : 2015-01-13
Resolution : 3.39 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 : 2.13.1
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

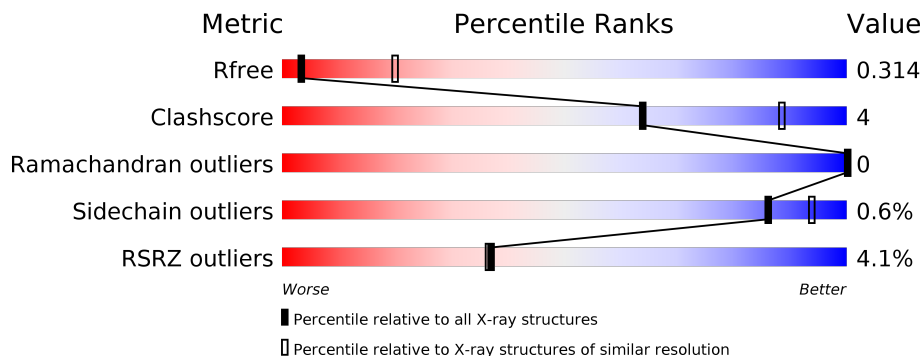
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.39 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	1026 (3.48-3.32)
Clashscore	141614	1055 (3.48-3.32)
Ramachandran outliers	138981	1038 (3.48-3.32)
Sidechain outliers	138945	1038 (3.48-3.32)
RSRZ outliers	127900	2173 (3.50-3.30)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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 $\leq 5\%$. 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	124	
1	C	124	
1	E	124	
1	G	124	
2	B	261	
2	D	261	

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Mol	Chain	Length	Quality of chain
2	F	261	 5% 86% 13%
2	H	261	 3% 87% 13%

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
2	MLY	B	189	-	-	-	X
2	MLY	B	190	-	-	-	X
2	MLY	H	189	-	-	-	X

2 Entry composition i

There are 6 unique types of molecules in this entry. The entry contains 11773 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 Neurogenic locus notch homolog protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	113	834	503	141	171	19	0	0	0
1	C	118	872	525	149	179	19	0	0	0
1	E	117	863	520	148	176	19	0	0	0
1	G	117	863	520	148	176	19	0	0	0

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	527	SER	-	expression tag	UNP Q07008
A	528	GLY	-	expression tag	UNP Q07008
A	529	ARG	-	expression tag	UNP Q07008
A	530	LEU	-	expression tag	UNP Q07008
A	531	GLU	-	expression tag	UNP Q07008
A	532	VAL	-	expression tag	UNP Q07008
A	533	LEU	-	expression tag	UNP Q07008
A	534	PHE	-	expression tag	UNP Q07008
A	535	GLN	-	expression tag	UNP Q07008
C	527	SER	-	expression tag	UNP Q07008
C	528	GLY	-	expression tag	UNP Q07008
C	529	ARG	-	expression tag	UNP Q07008
C	530	LEU	-	expression tag	UNP Q07008
C	531	GLU	-	expression tag	UNP Q07008
C	532	VAL	-	expression tag	UNP Q07008
C	533	LEU	-	expression tag	UNP Q07008
C	534	PHE	-	expression tag	UNP Q07008
C	535	GLN	-	expression tag	UNP Q07008
E	527	SER	-	expression tag	UNP Q07008
E	528	GLY	-	expression tag	UNP Q07008
E	529	ARG	-	expression tag	UNP Q07008

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Chain	Residue	Modelled	Actual	Comment	Reference
E	530	LEU	-	expression tag	UNP Q07008
E	531	GLU	-	expression tag	UNP Q07008
E	532	VAL	-	expression tag	UNP Q07008
E	533	LEU	-	expression tag	UNP Q07008
E	534	PHE	-	expression tag	UNP Q07008
E	535	GLN	-	expression tag	UNP Q07008
G	527	SER	-	expression tag	UNP Q07008
G	528	GLY	-	expression tag	UNP Q07008
G	529	ARG	-	expression tag	UNP Q07008
G	530	LEU	-	expression tag	UNP Q07008
G	531	GLU	-	expression tag	UNP Q07008
G	532	VAL	-	expression tag	UNP Q07008
G	533	LEU	-	expression tag	UNP Q07008
G	534	PHE	-	expression tag	UNP Q07008
G	535	GLN	-	expression tag	UNP Q07008

- Molecule 2 is a protein called Delta-like protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	257	2012	1251	362	376	23	0	0	0
2	D	258	2021	1256	363	379	23	0	0	0
2	F	257	2016	1253	362	378	23	0	0	0
2	H	259	2026	1259	364	380	23	0	0	0

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	26	SER	-	expression tag	UNP D3ZHH1
B	28	SER	GLY	engineered mutation	UNP D3ZHH1
B	107	LEU	PHE	engineered mutation	UNP D3ZHH1
B	206	PRO	LEU	engineered mutation	UNP D3ZHH1
B	257	LYS	ASN	engineered mutation	UNP D3ZHH1
B	284	ALA	-	expression tag	UNP D3ZHH1
B	285	ALA	-	expression tag	UNP D3ZHH1
B	286	ALA	-	expression tag	UNP D3ZHH1
D	26	SER	-	expression tag	UNP D3ZHH1
D	28	SER	GLY	engineered mutation	UNP D3ZHH1
D	107	LEU	PHE	engineered mutation	UNP D3ZHH1

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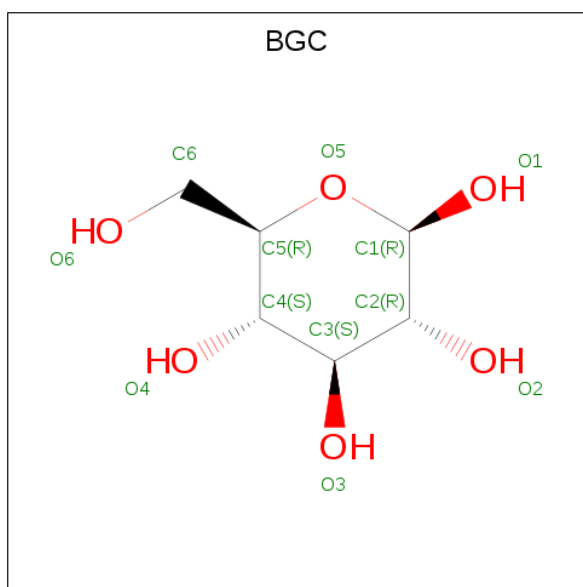
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Chain	Residue	Modelled	Actual	Comment	Reference
D	206	PRO	LEU	engineered mutation	UNP D3ZHH1
D	257	LYS	ASN	engineered mutation	UNP D3ZHH1
D	284	ALA	-	expression tag	UNP D3ZHH1
D	285	ALA	-	expression tag	UNP D3ZHH1
D	286	ALA	-	expression tag	UNP D3ZHH1
F	26	SER	-	expression tag	UNP D3ZHH1
F	28	SER	GLY	engineered mutation	UNP D3ZHH1
F	107	LEU	PHE	engineered mutation	UNP D3ZHH1
F	206	PRO	LEU	engineered mutation	UNP D3ZHH1
F	257	LYS	ASN	engineered mutation	UNP D3ZHH1
F	284	ALA	-	expression tag	UNP D3ZHH1
F	285	ALA	-	expression tag	UNP D3ZHH1
F	286	ALA	-	expression tag	UNP D3ZHH1
H	26	SER	-	expression tag	UNP D3ZHH1
H	28	SER	GLY	engineered mutation	UNP D3ZHH1
H	107	LEU	PHE	engineered mutation	UNP D3ZHH1
H	206	PRO	LEU	engineered mutation	UNP D3ZHH1
H	257	LYS	ASN	engineered mutation	UNP D3ZHH1
H	284	ALA	-	expression tag	UNP D3ZHH1
H	285	ALA	-	expression tag	UNP D3ZHH1
H	286	ALA	-	expression tag	UNP D3ZHH1

- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

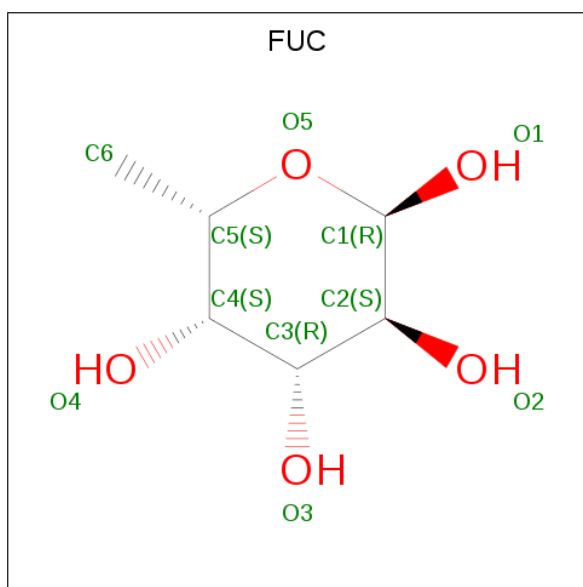
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	G	2	Total Ca 2 2	0	0
3	A	3	Total Ca 3 3	0	0
3	C	3	Total Ca 3 3	0	0
3	E	2	Total Ca 2 2	0	0

- Molecule 4 is beta-D-glucopyranose (three-letter code: BGC) (formula: C₆H₁₂O₆).



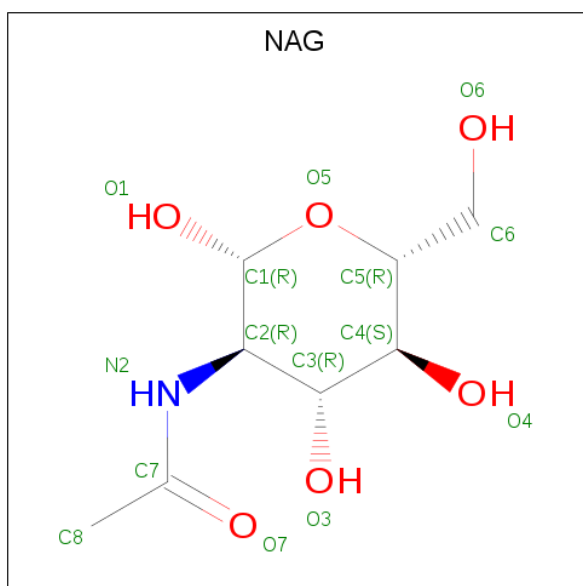
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O 11 6 5	0	0
4	A	1	Total C O 11 6 5	0	0
4	A	1	Total C O 11 6 5	0	0
4	C	1	Total C O 11 6 5	0	0
4	C	1	Total C O 11 6 5	0	0
4	C	1	Total C O 11 6 5	0	0
4	E	1	Total C O 11 6 5	0	0
4	E	1	Total C O 11 6 5	0	0
4	E	1	Total C O 11 6 5	0	0
4	G	1	Total C O 11 6 5	0	0
4	G	1	Total C O 11 6 5	0	0
4	G	1	Total C O 11 6 5	0	0

- Molecule 5 is alpha-L-fucopyranose (three-letter code: FUC) (formula: C₆H₁₂O₅).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			10	6	4		
5	C	1	Total	C	O	0	0
			10	6	4		
5	E	1	Total	C	O	0	0
			10	6	4		
5	G	1	Total	C	O	0	0
			10	6	4		

- Molecule 6 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).

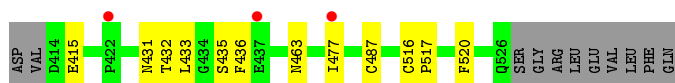
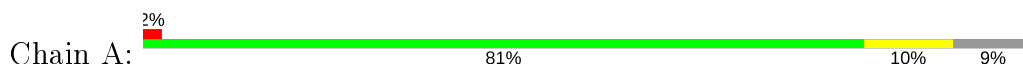


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	B	1	Total 14	C 8	N 1	O 5	0	0
6	D	1	Total 14	C 8	N 1	O 5	0	0
6	F	1	Total 14	C 8	N 1	O 5	0	0
6	F	1	Total 14	C 8	N 1	O 5	0	0
6	H	1	Total 14	C 8	N 1	O 5	0	0
6	H	1	Total 14	C 8	N 1	O 5	0	0

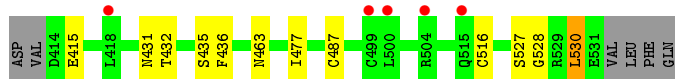
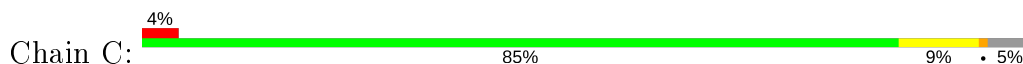
3 Residue-property plots [i](#)

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.

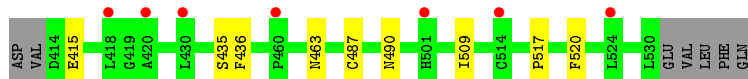
- Molecule 1: Neurogenic locus notch homolog protein 1



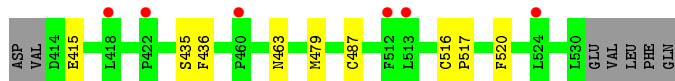
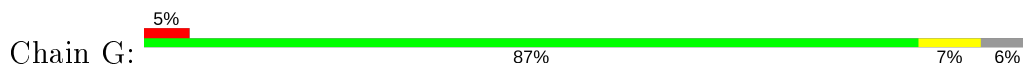
- Molecule 1: Neurogenic locus notch homolog protein 1



- Molecule 1: Neurogenic locus notch homolog protein 1



- Molecule 1: Neurogenic locus notch homolog protein 1

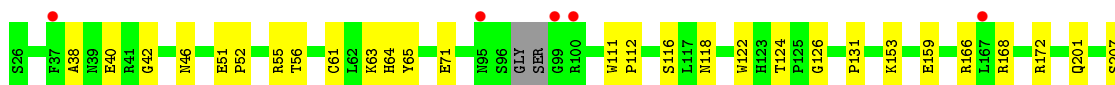
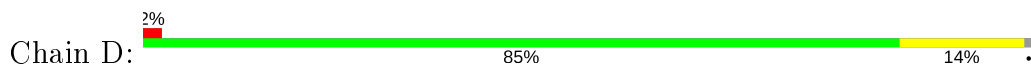


- Molecule 2: Delta-like protein

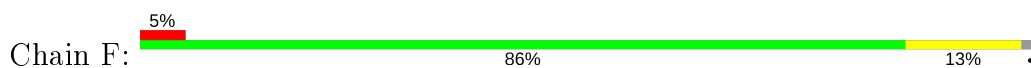




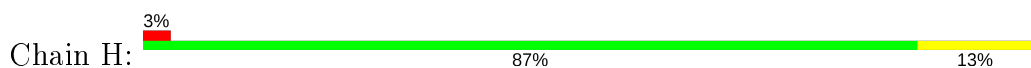
- Molecule 2: Delta-like protein



- Molecule 2: Delta-like protein



- Molecule 2: Delta-like protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	84.56Å 95.57Å 122.90Å 90.00° 96.21° 90.00°	Depositor
Resolution (Å)	84.06 – 3.39 84.06 – 3.39	Depositor EDS
% Data completeness (in resolution range)	98.8 (84.06-3.39) 99.2 (84.06-3.39)	Depositor EDS
R_{merge}	0.22	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.38 (at 3.41Å)	Xtrriage
Refinement program	PHENIX DEV-1839	Depositor
R, R_{free}	0.257 , 0.310 0.259 , 0.314	Depositor DCC
R_{free} test set	1339 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å ²)	87.0	Xtrriage
Anisotropy	0.664	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 74.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.42$, $\langle L^2 \rangle = 0.25$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	11773	wwPDB-VP
Average B, all atoms (Å ²)	144.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 29.37 % 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 1.5689e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: CA, BGC, NAG, MLY, FUC

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	0.23	0/849	0.46	0/1150
1	C	0.22	0/887	0.45	1/1200 (0.1%)
1	E	0.23	0/878	0.43	0/1188
1	G	0.22	0/878	0.42	0/1188
2	B	0.21	0/2014	0.39	0/2739
2	D	0.21	0/2023	0.39	0/2751
2	F	0.21	0/2018	0.39	0/2744
2	H	0.21	0/2028	0.39	0/2758
All	All	0.21	0/11575	0.40	1/15718 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	530	LEU	CA-CB-CG	5.05	126.92	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	834	0	729	8	0
1	C	872	0	768	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	863	0	762	4	0
1	G	863	0	762	4	0
2	B	2012	0	1864	19	0
2	D	2021	0	1873	24	0
2	F	2016	0	1867	19	0
2	H	2026	0	1877	20	0
3	A	3	0	0	0	0
3	C	3	0	0	0	0
3	E	2	0	0	0	0
3	G	2	0	0	0	0
4	A	33	0	30	1	0
4	C	33	0	30	0	0
4	E	33	0	30	0	0
4	G	33	0	30	0	0
5	A	10	0	10	0	0
5	C	10	0	10	1	0
5	E	10	0	10	1	0
5	G	10	0	10	1	0
6	B	14	0	13	0	0
6	D	14	0	13	0	0
6	F	28	0	25	0	0
6	H	28	0	25	0	0
All	All	11773	0	10738	94	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 4.

All (94) 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:431:ASN:OD1	1:A:432:THR:O	1.78	1.01
1:E:435:SER:OG	1:E:436:PHE:N	2.21	0.72
2:D:122:TRP:HH2	2:H:67:ALA:HB1	1.54	0.71
1:A:435:SER:OG	1:A:436:PHE:N	2.24	0.70
1:G:435:SER:OG	1:G:436:PHE:N	2.20	0.69
2:H:235:MLY:HB2	2:H:238:GLU:HG3	1.74	0.69
2:B:161:ASN:HB2	2:F:161:ASN:HB2	1.74	0.68
1:A:432:THR:HG22	1:A:433:LEU:H	1.59	0.68
1:C:435:SER:OG	1:C:436:PHE:N	2.27	0.67
2:F:124:THR:HG21	2:F:131:PRO:HB3	1.78	0.66
1:E:463:ASN:ND2	1:E:487:CYS:O	2.30	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:229:GLN:O	2:H:242:ARG:NH1	2.30	0.64
2:D:61:CYS:HB3	2:D:118:ASN:HB3	1.79	0.63
2:H:124:THR:HG21	2:H:131:PRO:HB3	1.80	0.63
1:C:463:ASN:ND2	1:C:487:CYS:O	2.32	0.62
1:E:490:ASN:HD21	1:E:509:ILE:HG12	1.65	0.62
2:B:124:THR:HG21	2:B:131:PRO:HB3	1.84	0.60
5:E:904:FUC:O4	2:F:65:TYR:O	2.19	0.60
2:F:235:MLY:HB2	2:F:238:GLU:HG3	1.83	0.60
2:D:124:THR:HG21	2:D:131:PRO:HB3	1.84	0.60
2:B:235:MLY:HB2	2:B:238:GLU:HG3	1.84	0.59
2:H:277:GLY:N	2:H:283:GLN:O	2.35	0.59
2:F:229:GLN:O	2:F:242:ARG:NH1	2.35	0.59
1:A:463:ASN:ND2	1:A:487:CYS:O	2.36	0.57
1:A:431:ASN:CG	1:A:432:THR:O	2.43	0.57
2:D:38:ALA:HB3	2:D:166:ARG:HB2	1.88	0.55
2:B:61:CYS:HB3	2:B:118:ASN:HB3	1.88	0.54
2:H:61:CYS:HB3	2:H:118:ASN:HB3	1.90	0.53
2:F:277:GLY:N	2:F:283:GLN:O	2.42	0.53
1:G:479:MET:HG2	5:G:904:FUC:H62	1.91	0.53
2:B:159:GLU:HG2	2:B:168:ARG:HG3	1.92	0.52
1:C:527:SER:OG	1:C:528:GLY:N	2.41	0.52
2:F:61:CYS:HB3	2:F:118:ASN:HB3	1.93	0.51
2:D:159:GLU:HG2	2:D:168:ARG:HG3	1.93	0.51
2:D:71:GLU:HG3	2:H:72:GLY:H	1.76	0.49
2:F:38:ALA:HB3	2:F:166:ARG:HB2	1.95	0.48
2:H:38:ALA:HB3	2:H:166:ARG:HB2	1.95	0.48
2:B:153:MLY:HH12	2:B:172:ARG:HH22	1.79	0.48
2:B:282:ASP:OD1	2:B:282:ASP:N	2.48	0.47
2:F:197:HIS:HB3	2:F:209:LEU:HD12	1.96	0.47
1:A:517:PRO:HG2	1:A:520:PHE:CD2	2.50	0.47
2:D:282:ASP:N	2:D:282:ASP:OD1	2.46	0.47
2:D:111:TRP:HA	2:D:112:PRO:HD3	1.75	0.47
2:H:159:GLU:HG2	2:H:168:ARG:HG3	1.98	0.46
2:D:55:ARG:NH2	2:D:126:GLY:O	2.48	0.46
2:B:197:HIS:HB3	2:B:209:LEU:HD12	1.96	0.46
2:H:63:LYS:HB3	2:H:74:CYS:HA	1.98	0.46
2:D:56:THR:HA	2:D:122:TRP:O	2.17	0.45
2:F:201:GLN:HG3	2:F:207:SER:HB2	1.99	0.45
2:H:180:TYR:CE1	2:H:188:CYS:HB3	2.52	0.44
2:F:93:ASP:HA	2:F:101:ASN:HD22	1.82	0.44
2:F:111:TRP:HA	2:F:112:PRO:HD3	1.71	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:201:GLN:HG3	2:D:207:SER:HB2	2.00	0.44
2:H:282:ASP:N	2:H:282:ASP:OD1	2.49	0.44
2:H:201:GLN:HG3	2:H:207:SER:HB2	1.98	0.44
2:D:229:GLN:O	2:D:242:ARG:NH1	2.50	0.43
2:D:227:HIS:CE1	2:D:229:GLN:HB3	2.54	0.43
2:H:197:HIS:HB3	2:H:209:LEU:HD12	2.01	0.43
2:H:260:ARG:HB2	2:H:260:ARG:HE	1.68	0.43
2:B:180:TYR:CE1	2:B:188:CYS:HB3	2.54	0.43
2:B:55:ARG:NH2	2:B:126:GLY:O	2.52	0.43
1:E:517:PRO:HG2	1:E:520:PHE:CD2	2.54	0.42
2:B:229:GLN:O	2:B:242:ARG:NH1	2.48	0.42
2:D:235:MLY:HB2	2:D:238:GLU:HG3	2.00	0.42
2:B:111:TRP:HA	2:B:112:PRO:HD3	1.76	0.42
2:B:40:GLU:C	2:B:42:GLY:H	2.23	0.42
2:B:67:ALA:HB1	2:F:122:TRP:CH2	2.54	0.42
2:F:159:GLU:HG2	2:F:168:ARG:HG3	2.02	0.42
1:A:435:SER:HB2	4:A:606:BGC:H2	1.86	0.42
2:F:32:LEU:HD12	2:F:103:LEU:HD12	2.02	0.42
1:G:463:ASN:ND2	1:G:487:CYS:O	2.53	0.42
2:D:153:MLY:HH12	2:D:172:ARG:HH22	1.85	0.42
2:F:55:ARG:NH2	2:F:126:GLY:O	2.53	0.42
2:H:111:TRP:HA	2:H:112:PRO:HD3	1.72	0.42
2:D:244:GLY:O	2:D:267:PRO:HG3	2.20	0.41
2:B:153:MLY:HB2	2:B:153:MLY:HE3	1.85	0.41
1:C:477:ILE:HG21	2:D:64:HIS:CE1	2.55	0.41
2:D:46:ASN:HA	2:H:148:SER:H	1.85	0.41
5:C:602:FUC:O4	2:D:65:TYR:O	2.26	0.41
2:H:36:GLU:HG3	2:H:90:VAL:HG22	2.02	0.41
2:D:63:LYS:O	2:D:116:SER:N	2.44	0.41
2:F:198:TYR:HB3	2:F:217:CYS:SG	2.60	0.41
1:A:477:ILE:HG21	2:B:64:HIS:CE1	2.55	0.41
2:D:40:GLU:C	2:D:42:GLY:H	2.24	0.41
2:H:51:GLU:HA	2:H:52:PRO:HA	1.90	0.41
2:B:122:TRP:CH2	2:F:67:ALA:HB1	2.56	0.41
2:D:51:GLU:HA	2:D:52:PRO:HA	1.90	0.41
2:B:178:ASN:O	2:B:187:LEU:HD12	2.21	0.40
2:D:260:ARG:HB2	2:D:260:ARG:HE	1.74	0.40
2:F:260:ARG:HB2	2:F:260:ARG:HE	1.68	0.40
1:G:517:PRO:HG2	1:G:520:PHE:CD2	2.56	0.40
2:D:122:TRP:CH2	2:H:67:ALA:HB1	2.43	0.40
2:B:198:TYR:HB3	2:B:217:CYS:SG	2.62	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:431:ASN:OD1	1:C:432:THR:O	2.39	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	111/124 (90%)	102 (92%)	9 (8%)	0	100	100
1	C	116/124 (94%)	103 (89%)	13 (11%)	0	100	100
1	E	115/124 (93%)	108 (94%)	7 (6%)	0	100	100
1	G	115/124 (93%)	108 (94%)	7 (6%)	0	100	100
2	B	248/261 (95%)	231 (93%)	17 (7%)	0	100	100
2	D	249/261 (95%)	229 (92%)	20 (8%)	0	100	100
2	F	248/261 (95%)	229 (92%)	19 (8%)	0	100	100
2	H	250/261 (96%)	232 (93%)	18 (7%)	0	100	100
All	All	1452/1540 (94%)	1342 (92%)	110 (8%)	0	100	100

There are no Ramachandran outliers to report.

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	95/107 (89%)	93 (98%)	2 (2%)	53	76
1	C	99/107 (92%)	96 (97%)	3 (3%)	41	68
1	E	98/107 (92%)	97 (99%)	1 (1%)	76	88
1	G	98/107 (92%)	96 (98%)	2 (2%)	55	77
2	B	218/220 (99%)	218 (100%)	0	100	100
2	D	219/220 (100%)	219 (100%)	0	100	100
2	F	219/220 (100%)	219 (100%)	0	100	100
2	H	219/220 (100%)	219 (100%)	0	100	100
All	All	1265/1308 (97%)	1257 (99%)	8 (1%)	86	94

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	415	GLU
1	A	516	CYS
1	C	415	GLU
1	C	516	CYS
1	C	530	LEU
1	E	415	GLU
1	G	415	GLU
1	G	516	CYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

20 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	MLY	B	190	2	9,10,11	0.54	0	6,11,13	0.77	0
2	MLY	D	189	2	9,10,11	0.50	0	6,11,13	0.85	0
2	MLY	B	189	2	9,10,11	0.55	0	6,11,13	0.83	0
2	MLY	F	190	2	9,10,11	0.55	0	6,11,13	0.77	0
2	MLY	F	153	2	9,10,11	0.50	0	6,11,13	0.91	0
2	MLY	H	190	2	9,10,11	0.53	0	6,11,13	0.79	0
2	MLY	H	189	2	9,10,11	0.52	0	6,11,13	0.85	0
2	MLY	D	190	2	9,10,11	0.54	0	6,11,13	0.78	0
2	MLY	D	235	2	9,10,11	0.53	0	6,11,13	0.87	0
2	MLY	F	235	2	9,10,11	0.51	0	6,11,13	0.89	0
2	MLY	B	235	2	9,10,11	0.55	0	6,11,13	0.88	0
2	MLY	F	215	2	9,10,11	0.51	0	6,11,13	0.79	0
2	MLY	B	153	2	9,10,11	0.49	0	6,11,13	0.92	0
2	MLY	D	215	2	9,10,11	0.53	0	6,11,13	0.78	0
2	MLY	H	215	2	9,10,11	0.53	0	6,11,13	0.77	0
2	MLY	B	215	2	9,10,11	0.52	0	6,11,13	0.79	0
2	MLY	H	153	2	9,10,11	0.51	0	6,11,13	0.90	0
2	MLY	H	235	2	9,10,11	0.54	0	6,11,13	0.83	0
2	MLY	F	189	2	9,10,11	0.53	0	6,11,13	0.85	0
2	MLY	D	153	2	9,10,11	0.48	0	6,11,13	0.89	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	MLY	B	190	2	-	1/8/9/11	-
2	MLY	D	189	2	-	0/8/9/11	-
2	MLY	B	189	2	-	0/8/9/11	-
2	MLY	F	190	2	-	0/8/9/11	-
2	MLY	F	153	2	-	0/8/9/11	-
2	MLY	H	190	2	-	1/8/9/11	-
2	MLY	H	189	2	-	0/8/9/11	-
2	MLY	D	190	2	-	1/8/9/11	-
2	MLY	D	235	2	-	4/8/9/11	-
2	MLY	F	235	2	-	2/8/9/11	-
2	MLY	B	235	2	-	3/8/9/11	-
2	MLY	F	215	2	-	2/8/9/11	-
2	MLY	B	153	2	-	2/8/9/11	-
2	MLY	D	215	2	-	3/8/9/11	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	MLY	H	215	2	-	1/8/9/11	-
2	MLY	B	215	2	-	1/8/9/11	-
2	MLY	H	153	2	-	0/8/9/11	-
2	MLY	H	235	2	-	1/8/9/11	-
2	MLY	F	189	2	-	0/8/9/11	-
2	MLY	D	153	2	-	0/8/9/11	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (22) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	H	190	MLY	O-C-CA-CB
2	D	235	MLY	N-CA-CB-CG
2	D	235	MLY	C-CA-CB-CG
2	F	235	MLY	O-C-CA-CB
2	D	215	MLY	O-C-CA-CB
2	B	235	MLY	N-CA-CB-CG
2	B	235	MLY	C-CA-CB-CG
2	H	235	MLY	O-C-CA-CB
2	D	235	MLY	CD-CE-NZ-CH1
2	D	235	MLY	CD-CE-NZ-CH2
2	B	235	MLY	CD-CE-NZ-CH2
2	H	215	MLY	CE-CD-CG-CB
2	F	235	MLY	CD-CE-NZ-CH2
2	B	215	MLY	CE-CD-CG-CB
2	F	215	MLY	CE-CD-CG-CB
2	D	190	MLY	CD-CE-NZ-CH2
2	B	153	MLY	CA-CB-CG-CD
2	D	215	MLY	CE-CD-CG-CB
2	D	215	MLY	CG-CD-CE-NZ
2	B	153	MLY	CD-CE-NZ-CH2
2	B	190	MLY	CD-CE-NZ-CH2
2	F	215	MLY	CG-CD-CE-NZ

There are no ring outliers.

6 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	235	MLY	1	0
2	F	235	MLY	1	0
2	B	235	MLY	1	0
2	B	153	MLY	2	0
2	H	235	MLY	1	0
2	D	153	MLY	1	0

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 32 ligands modelled in this entry, 10 are monoatomic - leaving 22 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	BGC	C	604	1	11,11,12	0.29	0	15,15,17	0.84	1 (6%)
4	BGC	E	903	1	11,11,12	0.30	0	15,15,17	0.69	0
5	FUC	G	904	1	10,10,11	0.70	0	14,14,16	0.71	0
5	FUC	A	605	1	10,10,11	0.64	0	14,14,16	0.73	0
6	NAG	H	302	2	14,14,15	0.27	0	17,19,21	0.53	0
4	BGC	E	906	1	11,11,12	0.26	0	15,15,17	0.71	0
4	BGC	C	603	1	11,11,12	0.34	0	15,15,17	0.87	0
5	FUC	C	602	1	10,10,11	0.72	0	14,14,16	0.68	0
4	BGC	A	606	1	11,11,12	0.34	0	15,15,17	0.78	0
4	BGC	C	601	1	11,11,12	0.30	0	15,15,17	0.64	0
6	NAG	H	301	2	14,14,15	0.23	0	17,19,21	0.39	0
5	FUC	E	904	1	10,10,11	0.67	0	14,14,16	0.73	0
4	BGC	A	604	1	11,11,12	0.28	0	15,15,17	0.70	0
6	NAG	F	301	2	14,14,15	0.25	0	17,19,21	0.40	0
6	NAG	D	301	2	14,14,15	0.34	0	17,19,21	0.50	0
4	BGC	G	905	1	11,11,12	0.34	0	15,15,17	0.97	0
6	NAG	F	302	2	14,14,15	0.21	0	17,19,21	0.51	0
6	NAG	B	301	2	14,14,15	0.25	0	17,19,21	0.52	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	BGC	A	607	1	11,11,12	0.27	0	15,15,17	0.70	0
4	BGC	E	905	1	11,11,12	0.35	0	15,15,17	0.81	0
4	BGC	G	906	1	11,11,12	0.26	0	15,15,17	0.68	0
4	BGC	G	903	1	11,11,12	0.30	0	15,15,17	0.81	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	BGC	C	604	1	-	2/2/19/22	0/1/1/1
4	BGC	E	903	1	-	2/2/19/22	0/1/1/1
5	FUC	G	904	1	-	-	0/1/1/1
5	FUC	A	605	1	-	-	0/1/1/1
6	NAG	H	302	2	-	2/6/23/26	0/1/1/1
4	BGC	E	906	1	-	2/2/19/22	0/1/1/1
4	BGC	C	603	1	-	2/2/19/22	0/1/1/1
5	FUC	C	602	1	-	-	0/1/1/1
4	BGC	A	606	1	-	0/2/19/22	0/1/1/1
4	BGC	C	601	1	-	2/2/19/22	0/1/1/1
6	NAG	H	301	2	-	1/6/23/26	0/1/1/1
5	FUC	E	904	1	-	-	0/1/1/1
4	BGC	A	604	1	-	2/2/19/22	0/1/1/1
6	NAG	F	301	2	-	0/6/23/26	0/1/1/1
6	NAG	D	301	2	-	2/6/23/26	0/1/1/1
4	BGC	G	905	1	-	0/2/19/22	0/1/1/1
6	NAG	F	302	2	-	4/6/23/26	0/1/1/1
6	NAG	B	301	2	-	4/6/23/26	0/1/1/1
4	BGC	A	607	1	-	2/2/19/22	0/1/1/1
4	BGC	E	905	1	-	2/2/19/22	0/1/1/1
4	BGC	G	906	1	-	2/2/19/22	0/1/1/1
4	BGC	G	903	1	-	2/2/19/22	0/1/1/1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	604	BGC	C1-C2-C3	2.24	112.42	109.67

There are no chirality outliers.

All (33) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	E	903	BGC	O5-C5-C6-O6
4	A	607	BGC	O5-C5-C6-O6
4	G	903	BGC	O5-C5-C6-O6
4	E	905	BGC	O5-C5-C6-O6
4	G	903	BGC	C4-C5-C6-O6
4	A	607	BGC	C4-C5-C6-O6
4	E	906	BGC	O5-C5-C6-O6
4	G	906	BGC	C4-C5-C6-O6
4	C	601	BGC	O5-C5-C6-O6
4	C	604	BGC	O5-C5-C6-O6
4	C	603	BGC	O5-C5-C6-O6
4	E	903	BGC	C4-C5-C6-O6
4	E	905	BGC	C4-C5-C6-O6
6	B	301	NAG	O5-C5-C6-O6
4	G	906	BGC	O5-C5-C6-O6
4	A	604	BGC	O5-C5-C6-O6
4	E	906	BGC	C4-C5-C6-O6
4	A	604	BGC	C4-C5-C6-O6
6	F	302	NAG	C8-C7-N2-C2
6	F	302	NAG	O7-C7-N2-C2
6	B	301	NAG	C8-C7-N2-C2
6	B	301	NAG	O7-C7-N2-C2
4	C	601	BGC	C4-C5-C6-O6
4	C	603	BGC	C4-C5-C6-O6
4	C	604	BGC	C4-C5-C6-O6
6	B	301	NAG	C4-C5-C6-O6
6	H	302	NAG	O5-C5-C6-O6
6	D	301	NAG	O5-C5-C6-O6
6	H	301	NAG	O5-C5-C6-O6
6	H	302	NAG	C4-C5-C6-O6
6	F	302	NAG	C4-C5-C6-O6
6	F	302	NAG	O5-C5-C6-O6
6	D	301	NAG	C4-C5-C6-O6

There are no ring outliers.

4 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	G	904	FUC	1	0
5	C	602	FUC	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	606	BGC	1	0
5	E	904	FUC	1	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

6.1 Protein, DNA and RNA chains

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	113/124 (91%)	0.19	3 (2%) 54 53	101, 160, 220, 298	0
1	C	118/124 (95%)	0.14	5 (4%) 36 35	76, 145, 207, 247	0
1	E	117/124 (94%)	0.28	7 (5%) 21 23	119, 174, 225, 243	0
1	G	117/124 (94%)	0.26	6 (5%) 28 28	94, 138, 203, 241	0
2	B	252/261 (96%)	0.24	11 (4%) 34 34	93, 139, 226, 310	0
2	D	253/261 (96%)	0.14	6 (2%) 59 57	60, 117, 175, 271	0
2	F	252/261 (96%)	0.37	13 (5%) 27 27	84, 162, 215, 291	0
2	H	254/261 (97%)	0.15	9 (3%) 44 43	67, 129, 181, 232	0
All	All	1476/1540 (95%)	0.22	60 (4%) 37 36	60, 142, 213, 310	0

All (60) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	F	95	ASN	5.6
2	H	99	GLY	5.4
2	B	36	GLU	5.2
2	B	37	PHE	4.8
2	F	96	SER	4.7
2	H	199	GLU	4.2
2	B	95	ASN	4.1
2	B	34	LEU	3.9
2	F	168	ARG	3.8
2	H	178	ASN	3.8
2	F	167	LEU	3.5
2	D	95	ASN	3.5
1	E	501	HIS	3.4
2	B	30	PHE	3.4
2	B	41	ARG	3.4
1	C	418	LEU	3.3

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Mol	Chain	Res	Type	RSRZ
2	B	174	VAL	3.3
2	D	99	GLY	3.1
2	H	130	ARG	3.1
1	C	500	LEU	3.0
2	F	92	ARG	3.0
2	B	92	ARG	3.0
1	C	515	GLN	3.0
2	H	146	GLN	2.9
2	B	91	ILE	2.9
1	C	499	CYS	2.9
1	G	513	LEU	2.9
1	A	422	PRO	2.9
2	H	96	SER	2.8
1	E	514	CYS	2.8
1	G	512	PHE	2.8
2	D	167	LEU	2.6
2	F	199	GLU	2.5
2	H	100	ARG	2.5
2	D	239	CYS	2.5
1	G	524	LEU	2.5
2	F	57	PHE	2.5
2	F	194	HIS	2.5
1	A	437	GLU	2.5
2	D	37	PHE	2.4
1	E	460	PRO	2.4
2	F	178	ASN	2.4
1	E	420	ALA	2.3
2	F	91	ILE	2.3
2	B	199	GLU	2.3
1	C	504	ARG	2.3
1	G	460	PRO	2.3
2	H	69	PHE	2.2
2	F	34	LEU	2.2
1	G	422	PRO	2.2
1	E	418	LEU	2.2
2	B	56	THR	2.2
1	E	430	LEU	2.1
2	H	167	LEU	2.1
2	F	90	VAL	2.1
2	D	100	ARG	2.1
1	E	524	LEU	2.1
2	F	104	GLN	2.1

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Mol	Chain	Res	Type	RSRZ
1	G	418	LEU	2.1
1	A	477	ILE	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	MLY	D	153	11/12	0.67	0.30	109,119,137,139	0
2	MLY	F	153	11/12	0.70	0.31	155,169,186,188	0
2	MLY	B	189	11/12	0.74	0.52	122,132,136,138	0
2	MLY	H	153	11/12	0.75	0.25	132,145,158,158	0
2	MLY	B	190	11/12	0.77	0.51	143,160,186,187	0
2	MLY	H	189	11/12	0.78	0.59	145,154,164,169	0
2	MLY	D	215	11/12	0.81	0.36	136,145,155,160	0
2	MLY	H	190	11/12	0.81	0.39	124,132,147,151	0
2	MLY	H	215	11/12	0.81	0.35	141,156,163,164	0
2	MLY	F	189	11/12	0.82	0.38	156,167,184,185	0
2	MLY	F	190	11/12	0.82	0.32	167,173,177,178	0
2	MLY	F	215	11/12	0.84	0.52	168,178,188,197	0
2	MLY	H	235	11/12	0.85	0.38	79,121,143,144	0
2	MLY	D	189	11/12	0.86	0.25	102,109,123,127	0
2	MLY	F	235	11/12	0.88	0.32	177,188,198,200	0
2	MLY	B	153	11/12	0.89	0.14	140,142,149,152	0
2	MLY	D	190	11/12	0.89	0.21	113,129,148,149	0
2	MLY	B	215	11/12	0.90	0.38	110,119,134,138	0
2	MLY	B	235	11/12	0.93	0.34	93,107,134,138	0
2	MLY	D	235	11/12	0.95	0.36	106,118,140,150	0

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

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(Å ²)	Q<0.9
3	CA	C	605	1/1	0.73	0.17	277,277,277,277	0
4	BGC	A	607	11/12	0.73	0.31	180,186,193,194	0
4	BGC	A	606	11/12	0.76	0.28	136,146,177,181	0
3	CA	A	603	1/1	0.77	0.09	261,261,261,261	0
3	CA	A	601	1/1	0.77	0.13	382,382,382,382	0
4	BGC	C	604	11/12	0.79	0.21	108,124,129,131	0
6	NAG	F	301	14/15	0.80	0.32	114,129,140,145	0
6	NAG	D	301	14/15	0.80	0.24	92,112,122,144	0
4	BGC	E	906	11/12	0.80	0.21	103,113,118,121	0
4	BGC	C	603	11/12	0.81	0.24	100,113,135,137	0
6	NAG	H	302	14/15	0.82	0.22	94,118,123,129	0
6	NAG	H	301	14/15	0.83	0.30	101,124,132,137	0
4	BGC	G	903	11/12	0.83	0.25	63,91,102,105	0
3	CA	A	602	1/1	0.84	0.16	116,116,116,116	0
6	NAG	B	301	14/15	0.85	0.23	115,134,143,151	0
4	BGC	C	601	11/12	0.85	0.22	81,91,110,115	0
4	BGC	G	906	11/12	0.85	0.27	99,116,130,142	0
6	NAG	F	302	14/15	0.85	0.17	118,130,140,141	0
4	BGC	G	905	11/12	0.86	0.25	100,114,155,156	0
4	BGC	A	604	11/12	0.86	0.22	108,113,118,120	0
4	BGC	E	903	11/12	0.89	0.17	95,100,108,108	0
5	FUC	E	904	10/11	0.91	0.14	87,110,121,121	0
4	BGC	E	905	11/12	0.92	0.20	99,116,122,122	0
5	FUC	A	605	10/11	0.92	0.18	107,118,135,143	0
3	CA	G	901	1/1	0.92	0.19	123,123,123,123	0
3	CA	E	902	1/1	0.94	0.07	152,152,152,152	0
3	CA	C	607	1/1	0.94	0.09	185,185,185,185	0
5	FUC	G	904	10/11	0.95	0.14	71,80,88,89	0
5	FUC	C	602	10/11	0.95	0.20	79,87,111,114	0
3	CA	C	606	1/1	0.96	0.14	102,102,102,102	0
3	CA	E	901	1/1	0.97	0.12	107,107,107,107	0
3	CA	G	902	1/1	0.98	0.10	188,188,188,188	0

6.5 Other polymers [i](#)

There are no such residues in this entry.