Full wwPDB X-ray Structure Validation Report (i)

Dec 12, 2023 – 05:16 pm GMT

PDB ID	:	2WUB
Title	:	Crystal structure of HGFA in complex with the allosteric non- inhibitory antibody Fab40.deltaTrp
Authors	:	Ganesan, R.; Eigenbrot, C.; Shia, S.
Deposited on	:	2009-10-01
Resolution	:	2.90 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:

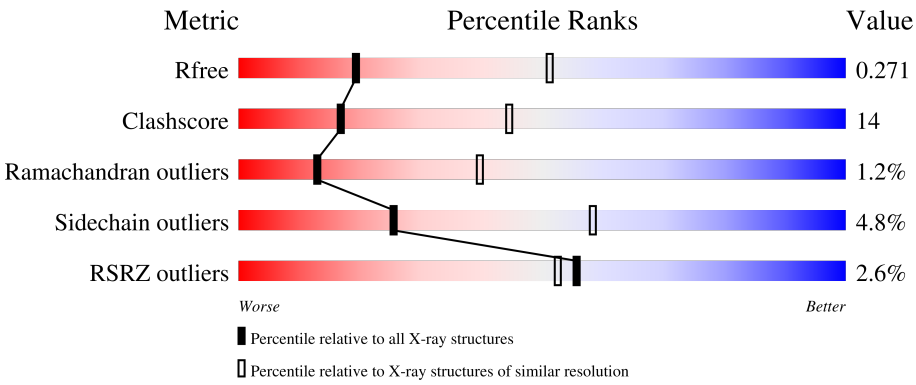
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	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 2.90 Å.

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



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

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	257	2% 65% 21% • • 10%
1	C	257	2% 59% 28% • 11%
2	B	35	17% • 80%
2	D	35	3% 14% • 83%
3	H	224	% 72% 23% • •

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Mol	Chain	Length	Quality of chain
3	R	224	6% 61% 30% 8%
4	L	214	68% 29%
4	Q	214	5% 68% 29%
5	E	3	67% 33%
5	F	3	100%

2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 10198 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 HEPATOCYTE GROWTH FACTOR ACTIVATOR LONG CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	232	1761	1119	303	326	13	0	0	1
1	C	229	1747	1110	305	319	13	0	0	1

- Molecule 2 is a protein called HEPATOCYTE GROWTH FACTOR ACTIVATOR SHORT CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	7	48	26	15	6	1	0	0	1
2	D	6	38	20	12	5	1	0	0	1

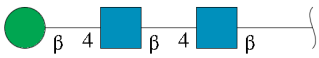
- Molecule 3 is a protein called FAB FRAGMENT FAB40.DELTATRP HEAVY CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	H	214	1584	1010	263	305	6	0	0	1
3	R	206	1522	971	252	294	5	0	0	1

- Molecule 4 is a protein called FAB FRAGMENT FAB40.DELTATRP LIGHT CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
4	L	214	1633	1018	275	334	6	0	0	0
4	Q	214	1633	1018	275	334	6	0	0	0

- Molecule 5 is an oligosaccharide called beta-D-mannopyranose-(1-4)-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			
5	E	3	39	22	2	15	0	0	0
5	F	3	39	22	2	15	0	0	0

- Molecule 6 is water.

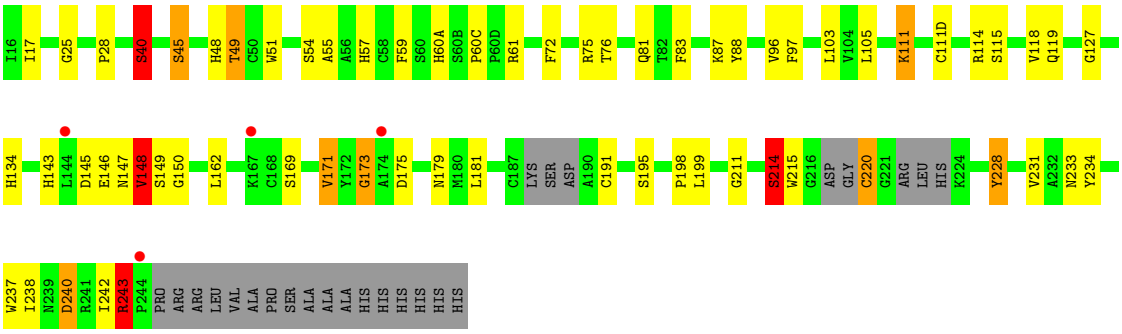
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
6	A	24	24	24	0	0
6	B	2	2	2	0	0
6	C	23	23	23	0	0
6	D	1	1	1	0	0
6	H	25	25	25	0	0
6	L	22	22	22	0	0
6	Q	27	27	27	0	0
6	R	30	30	30	0	0

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.

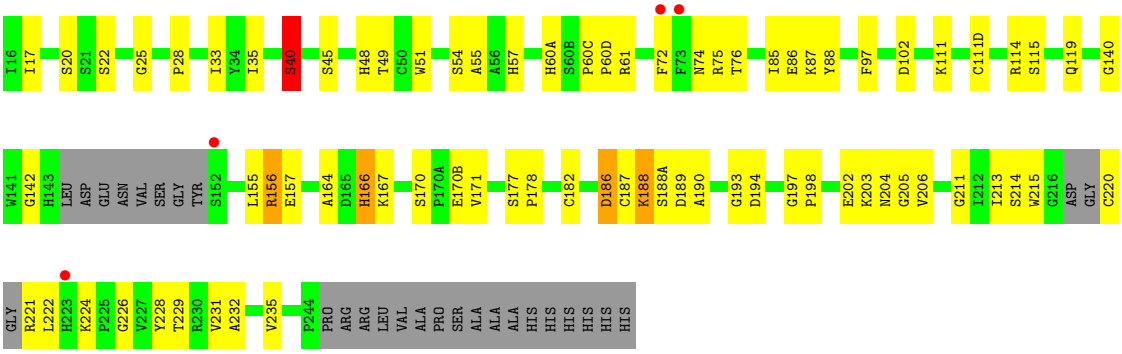
• Molecule 1: HEPATOCYTE GROWTH FACTOR ACTIVATOR LONG CHAIN

Chain A: 2% 65% 21% 10%



• Molecule 1: HEPATOCYTE GROWTH FACTOR ACTIVATOR LONG CHAIN

Chain C: 2% 59% 28% 11%

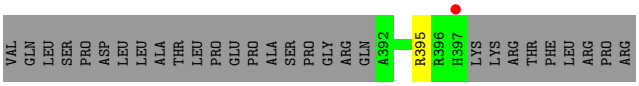


• Molecule 2: HEPATOCYTE GROWTH FACTOR ACTIVATOR SHORT CHAIN

Chain B: 17% 80%

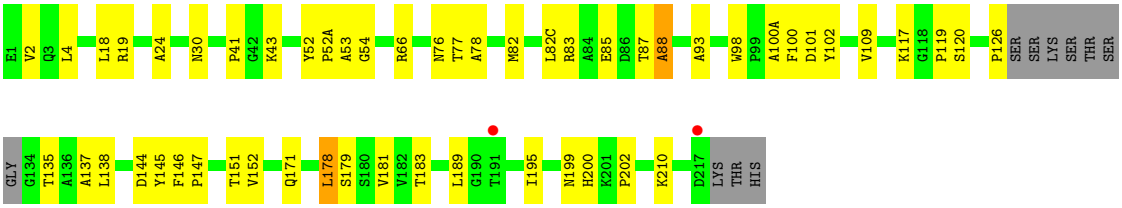
• Molecule 2: HEPATOCYTE GROWTH FACTOR ACTIVATOR SHORT CHAIN

Chain D: 3% 14% 83%



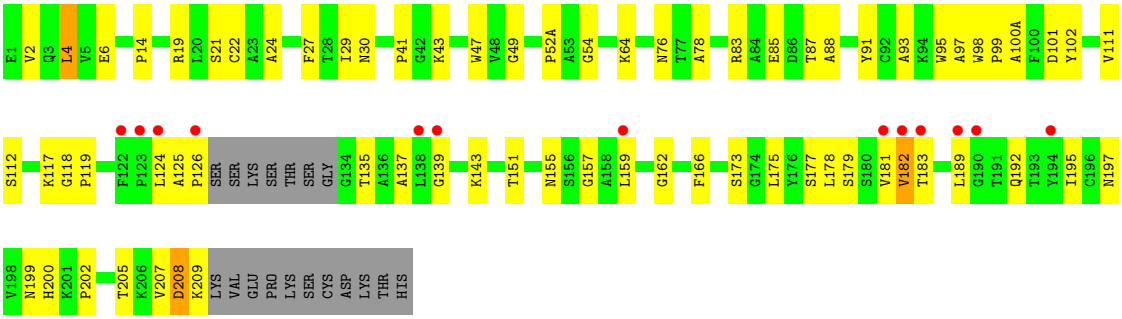
• Molecule 3: FAB FRAGMENT FAB40.DELTATRP HEAVY CHAIN

Chain H:



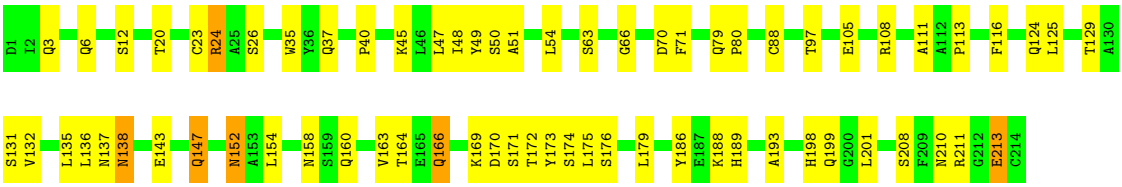
• Molecule 3: FAB FRAGMENT FAB40.DELTATRP HEAVY CHAIN

Chain R:



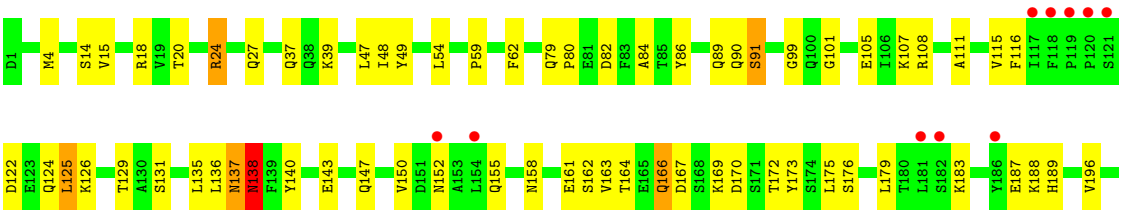
• Molecule 4: FAB FRAGMENT FAB40.DELTATRP LIGHT CHAIN

Chain L:



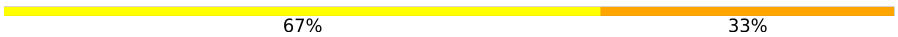
• Molecule 4: FAB FRAGMENT FAB40.DELTATRP LIGHT CHAIN

Chain Q:



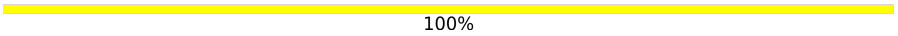
• Molecule 5: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain E:



• Molecule 5: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-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	72.36Å 89.53Å 118.47Å 90.00° 91.08° 90.00°	Depositor
Resolution (Å)	19.99 – 2.90 49.40 – 2.90	Depositor EDS
% Data completeness (in resolution range)	98.8 (19.99-2.90) 98.8 (49.40-2.90)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.49 (at 2.91Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.216 , 0.275 0.212 , 0.271	Depositor DCC
R_{free} test set	3318 reflections (9.95%)	wwPDB-VP
Wilson B-factor (Å^2)	60.7	Xtriage
Anisotropy	0.314	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.27 , 22.4	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	0.060 for h,-k,-l	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	10198	wwPDB-VP
Average B, all atoms (Å^2)	71.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.72% of the height of the origin peak. No significant pseudotranslation is detected.*

[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: BMA, 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	0.44	1/1811 (0.1%)	0.58	0/2465
1	C	0.43	0/1798	0.63	1/2447 (0.0%)
2	B	0.67	0/48	0.49	0/62
2	D	0.77	0/37	0.52	0/47
3	H	0.48	0/1627	0.65	0/2224
3	R	0.48	0/1564	0.64	0/2141
4	L	0.43	0/1667	0.64	0/2262
4	Q	0.41	0/1667	0.61	0/2262
All	All	0.45	1/10219 (0.0%)	0.62	1/13910 (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	1
1	C	0	2
All	All	0	3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	40	SER	C-N	-6.87	1.18	1.34

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	40	SER	O-C-N	-7.54	110.63	122.70

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	40	SER	Mainchain
1	C	40	SER	Mainchain
1	C	74	ASN	Mainchain

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	1761	0	1674	51	0
1	C	1747	0	1672	54	0
2	B	48	0	44	1	0
2	D	38	0	37	0	0
3	H	1584	0	1543	34	0
3	R	1522	0	1473	55	0
4	L	1633	0	1585	50	0
4	Q	1633	0	1586	50	0
5	E	39	0	34	1	0
5	F	39	0	34	0	0
6	A	24	0	0	4	0
6	B	2	0	0	1	0
6	C	23	0	0	0	0
6	D	1	0	0	0	0
6	H	25	0	0	2	0
6	L	22	0	0	3	0
6	Q	27	0	0	3	0
6	R	30	0	0	6	0
All	All	10198	0	9682	279	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 14.

All (279) 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)
4:L:24:ARG:HH11	4:L:24:ARG:HB2	1.32	0.93
4:L:210:ASN:HB2	4:L:213:GLU:HG3	1.53	0.89
1:C:164:ALA:HB3	1:C:167:LYS:HG2	1.56	0.86
3:H:82:MET:HB3	3:H:82(C):LEU:HD21	1.58	0.84
3:R:195:ILE:HG23	3:R:209:LYS:N	1.93	0.83
4:Q:18:ARG:HD2	6:Q:2005:HOH:O	1.78	0.82
6:A:2024:HOH:O	5:E:3:BMA:O2	2.00	0.78
1:C:170:SER:HB2	1:C:171:VAL:HG23	1.65	0.77
1:A:49:THR:O	1:A:111(D):CYS:HB2	1.88	0.74
3:H:200:HIS:CD2	3:H:202:PRO:HD2	2.23	0.74
3:R:159:LEU:HD21	3:R:182:VAL:HG21	1.68	0.72
1:C:72:PHE:HB2	1:C:75:ARG:HB2	1.70	0.72
1:A:48:HIS:HB3	1:A:51:TRP:HB2	1.73	0.71
1:C:186:ASP:HA	1:C:222:LEU:HD12	1.74	0.70
4:Q:189:HIS:O	4:Q:211:ARG:HD3	1.91	0.70
1:C:167:LYS:O	1:C:171:VAL:HG21	1.92	0.70
4:L:37:GLN:HB2	4:L:47:LEU:HD11	1.74	0.69
3:R:24:ALA:HB3	3:R:76:ASN:ND2	2.07	0.69
4:L:108:ARG:NH2	4:L:111:ALA:HB2	2.08	0.68
4:L:24:ARG:HH11	4:L:24:ARG:CB	2.03	0.68
4:L:79:GLN:HB3	4:L:80:PRO:HD2	1.76	0.68
1:A:72:PHE:HB2	1:A:75:ARG:HB2	1.76	0.67
3:H:195:ILE:HG12	3:H:210:LYS:HA	1.77	0.66
1:C:48:HIS:HB3	1:C:51:TRP:HB2	1.78	0.65
3:H:119:PRO:HB3	3:H:145:TYR:HB3	1.78	0.65
1:C:177:SER:HB2	1:C:178:PRO:HD2	1.77	0.65
4:Q:158:ASN:HD21	4:Q:179:LEU:HD11	1.61	0.65
4:Q:91:SER:HB3	3:R:95:TRP:HH2	1.63	0.63
1:C:22:SER:HB3	1:C:155:LEU:O	1.99	0.62
4:Q:135:LEU:HD21	4:Q:137:ASN:HD22	1.63	0.62
3:R:98:TRP:O	3:R:100(A):ALA:N	2.32	0.62
3:H:135:THR:HB	6:H:2023:HOH:O	2.00	0.61
3:R:195:ILE:CG2	3:R:209:LYS:N	2.63	0.61
1:A:87:LYS:HB2	3:H:54:GLY:O	2.01	0.61
1:A:134:HIS:HB3	1:A:162:LEU:HD12	1.82	0.61
3:R:22:CYS:HB3	3:R:78:ALA:HB3	1.81	0.61
3:H:24:ALA:HB3	3:H:76:ASN:ND2	2.15	0.61
3:H:85:GLU:H	3:H:85:GLU:CD	2.02	0.61
4:L:97:THR:HA	6:L:2013:HOH:O	2.01	0.61
3:R:200:HIS:CD2	3:R:202:PRO:HD2	2.36	0.61
3:H:41:PRO:O	3:H:43:LYS:HG2	2.01	0.61
4:Q:187:GLU:HA	4:Q:211:ARG:HD2	1.82	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:87:LYS:HB2	3:R:54:GLY:O	2.01	0.60
1:C:202:GLU:OE1	1:C:205:GLY:HA2	2.01	0.60
1:A:148:VAL:HG12	1:A:149:SER:H	1.66	0.60
1:C:28:PRO:HB2	1:C:119:GLN:H	1.67	0.60
3:H:52:TYR:CE1	3:H:53:ALA:HB3	2.36	0.60
1:A:25:GLY:O	1:A:28:PRO:HD3	2.03	0.59
1:C:28:PRO:HB2	1:C:119:GLN:N	2.18	0.59
1:C:232:ALA:HA	1:C:235:VAL:HG23	1.83	0.59
1:A:61:ARG:HD3	1:A:87:LYS:HA	1.85	0.59
3:R:98:TRP:C	3:R:100(A):ALA:H	2.07	0.58
1:C:22:SER:HB2	1:C:157:GLU:HG2	1.85	0.58
3:R:151:THR:HB	3:R:199:ASN:HB3	1.84	0.58
4:Q:79:GLN:HB3	4:Q:80:PRO:HD2	1.84	0.58
1:C:61:ARG:HD3	1:C:87:LYS:HA	1.85	0.57
4:Q:135:LEU:HD21	4:Q:137:ASN:ND2	2.19	0.57
1:C:197:GLY:O	1:C:213:ILE:HG13	2.04	0.57
4:Q:115:VAL:CG2	4:Q:196:VAL:HG21	2.35	0.57
1:A:111:LYS:HA	6:A:2010:HOH:O	2.03	0.57
3:H:19:ARG:HD2	6:H:2003:HOH:O	2.05	0.56
1:A:179:ASN:HA	1:A:233:ASN:HD22	1.71	0.56
1:C:61:ARG:HG2	1:C:85:ILE:HD11	1.87	0.56
4:L:158:ASN:HD21	4:L:179:LEU:HD11	1.71	0.56
3:R:47:TRP:CZ2	3:R:49:GLY:HA2	2.40	0.56
3:R:157:GLY:HA3	6:R:2024:HOH:O	2.05	0.56
4:L:188:LYS:HG3	4:L:189:HIS:ND1	2.21	0.55
4:Q:158:ASN:ND2	4:Q:179:LEU:HD11	2.21	0.55
3:R:192:GLN:HB2	6:R:2029:HOH:O	2.04	0.55
3:H:181:VAL:HG11	4:L:135:LEU:HD12	1.88	0.55
1:A:145:ASP:OD2	1:A:148:VAL:HG23	2.07	0.55
1:C:170:SER:HB2	1:C:171:VAL:CG2	2.36	0.55
4:Q:107:LYS:HA	4:Q:140:TYR:OH	2.07	0.55
4:L:166:GLN:NE2	4:L:173:TYR:OH	2.40	0.54
3:R:41:PRO:O	3:R:43:LYS:HG2	2.07	0.54
1:A:145:ASP:OD2	1:A:147:ASN:HB2	2.07	0.54
4:Q:116:PHE:HB3	6:Q:2022:HOH:O	2.07	0.54
1:C:45:SER:OG	1:C:198:PRO:HB3	2.08	0.54
4:Q:188:LYS:HG3	4:Q:189:HIS:ND1	2.22	0.54
1:A:17:ILE:HD11	1:A:220:CYS:HB2	1.90	0.54
1:C:17:ILE:CD1	1:C:220:CYS:HB2	2.38	0.54
1:C:49:THR:O	1:C:111(D):CYS:HB2	2.07	0.54
1:C:114:ARG:HB3	1:C:119:GLN:HE22	1.73	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:164:ALA:HB3	1:C:167:LYS:CG	2.32	0.54
1:A:28:PRO:HB2	1:A:119:GLN:N	2.23	0.53
4:Q:175:LEU:HD23	4:Q:176:SER:N	2.22	0.53
4:Q:135:LEU:HD23	4:Q:135:LEU:C	2.29	0.53
3:R:85:GLU:CD	3:R:85:GLU:H	2.11	0.53
3:H:181:VAL:HG21	4:L:135:LEU:HD11	1.91	0.53
2:B:397:HIS:HA	6:B:2002:HOH:O	2.08	0.53
4:Q:108:ARG:NH2	4:Q:111:ALA:HB2	2.23	0.53
4:Q:24:ARG:HB2	4:Q:24:ARG:HH11	1.74	0.53
3:R:2:VAL:HB	3:R:102:TYR:CZ	2.45	0.52
3:H:181:VAL:HG21	4:L:135:LEU:CD1	2.40	0.52
4:L:175:LEU:HD23	4:L:176:SER:N	2.25	0.52
3:R:2:VAL:HB	3:R:102:TYR:CE1	2.44	0.52
1:A:81:GLN:HE21	1:A:118:VAL:HG21	1.75	0.52
3:H:171:GLN:HA	4:L:160:GLN:HE22	1.75	0.52
4:Q:135:LEU:CD1	3:R:181:VAL:HG21	2.39	0.51
4:Q:125:LEU:CD1	4:Q:183:LYS:HG3	2.39	0.51
3:R:19:ARG:HD2	6:R:2003:HOH:O	2.08	0.51
1:C:202:GLU:HA	1:C:206:VAL:O	2.10	0.51
3:H:151:THR:HB	3:H:199:ASN:HB3	1.93	0.51
3:R:162:GLY:HA3	6:R:2025:HOH:O	2.09	0.51
3:H:87:THR:O	3:H:88:ALA:HB2	2.10	0.50
3:R:87:THR:O	3:R:88:ALA:HB2	2.11	0.50
1:C:17:ILE:HD11	1:C:220:CYS:HB2	1.93	0.50
1:A:105:LEU:HD12	1:A:237:TRP:CZ3	2.47	0.50
4:L:20:THR:CG2	4:Q:24:ARG:HH21	2.24	0.50
4:L:48:ILE:HG22	4:L:49:TYR:N	2.26	0.50
3:R:117:LYS:HG2	3:R:118:GLY:O	2.11	0.50
1:A:17:ILE:HD11	1:A:191:CYS:SG	2.52	0.50
4:L:50:SER:O	4:L:51:ALA:HB3	2.12	0.50
1:C:20:SER:C	1:C:156:ARG:NH2	2.65	0.50
3:H:2:VAL:HB	3:H:102:TYR:CZ	2.47	0.50
1:C:167:LYS:HA	1:C:170:SER:OG	2.12	0.50
3:H:126:PRO:HD3	3:H:138:LEU:HB3	1.93	0.50
3:R:93:ALA:HA	3:R:102:TYR:O	2.11	0.50
3:R:197:ASN:N	3:R:197:ASN:HD22	2.10	0.50
1:C:211:GLY:HA2	1:C:231:VAL:HG23	1.94	0.49
4:L:12:SER:HA	4:L:105:GLU:O	2.13	0.49
3:R:6:GLU:HA	3:R:21:SER:O	2.13	0.49
1:A:146:GLU:HA	1:A:191:CYS:SG	2.53	0.49
4:L:158:ASN:ND2	4:L:179:LEU:HD11	2.27	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:203:LYS:O	1:C:204:ASN:HB2	2.13	0.49
3:H:178:LEU:HD12	3:H:179:SER:N	2.28	0.49
4:L:154:LEU:H	4:L:154:LEU:HD22	1.77	0.49
1:A:114:ARG:HB3	1:A:119:GLN:HE22	1.76	0.49
4:Q:91:SER:HB3	3:R:95:TRP:CH2	2.45	0.49
1:A:75:ARG:HG3	1:A:75:ARG:HH11	1.77	0.48
1:A:211:GLY:HA2	1:A:231:VAL:HG23	1.95	0.48
1:A:60(A):HIS:HD2	6:A:2003:HOH:O	1.94	0.48
4:L:3:GLN:HB2	4:L:26:SER:HB3	1.95	0.48
1:C:17:ILE:HD11	1:C:189:ASP:HB3	1.95	0.48
1:C:33:ILE:HG22	1:C:35:ILE:HG13	1.94	0.48
4:L:186:TYR:CE2	4:L:211:ARG:HD3	2.49	0.48
1:A:146:GLU:HG2	1:A:191:CYS:SG	2.54	0.48
1:A:28:PRO:HB2	1:A:119:GLN:H	1.79	0.48
4:L:135:LEU:HD23	4:L:135:LEU:C	2.34	0.48
1:A:59:PHE:HB3	1:A:88:TYR:CE2	2.49	0.48
1:A:17:ILE:HD13	1:A:146:GLU:OE2	2.13	0.48
3:H:30:ASN:HA	3:H:52(A):PRO:HB2	1.95	0.47
4:Q:115:VAL:HG22	4:Q:196:VAL:HG21	1.95	0.47
3:R:155:ASN:HA	6:R:2030:HOH:O	2.13	0.47
1:A:169:SER:HA	1:A:173:GLY:HA2	1.96	0.47
1:A:242:ILE:O	1:A:243:ARG:C	2.52	0.47
1:A:240:ASP:HB2	6:A:2018:HOH:O	2.15	0.47
3:R:119:PRO:HD2	3:R:205:THR:CB	2.43	0.47
4:L:154:LEU:HD22	4:L:154:LEU:N	2.30	0.47
4:L:170:ASP:O	4:L:171:SER:HB2	2.15	0.47
4:Q:135:LEU:HD11	3:R:181:VAL:HG21	1.97	0.47
4:L:23:CYS:HB2	4:L:35:TRP:CH2	2.50	0.47
4:Q:143:GLU:CD	4:Q:143:GLU:H	2.18	0.47
4:Q:122:ASP:O	4:Q:126:LYS:HG2	2.14	0.47
4:L:132:VAL:HB	4:L:179:LEU:HB3	1.97	0.47
1:A:234:TYR:O	1:A:238:ILE:HG13	2.14	0.46
4:Q:135:LEU:HD23	4:Q:136:LEU:N	2.30	0.46
3:R:91:TYR:HE1	6:R:2019:HOH:O	1.98	0.46
4:Q:14:SER:O	4:Q:15:VAL:C	2.53	0.46
1:C:57:HIS:ND1	1:C:102:ASP:OD2	2.40	0.46
4:L:48:ILE:HD13	4:L:54:LEU:HA	1.98	0.46
1:A:45:SER:OG	1:A:198:PRO:HB3	2.16	0.46
1:A:75:ARG:HG3	1:A:75:ARG:NH1	2.30	0.46
1:A:115:SER:H	1:A:119:GLN:HE21	1.64	0.46
1:C:61:ARG:HG2	1:C:85:ILE:CD1	2.45	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:228:TYR:CD1	1:A:228:TYR:N	2.83	0.46
3:H:66:ARG:HD3	3:H:83:ARG:HH21	1.80	0.46
4:L:135:LEU:HD23	4:L:136:LEU:N	2.31	0.45
3:R:124:LEU:HD12	3:R:139:GLY:HA3	1.99	0.45
3:H:18:LEU:HD23	3:H:109:VAL:HG13	1.98	0.45
4:Q:48:ILE:HG22	4:Q:49:TYR:N	2.32	0.45
3:H:77:THR:HG22	3:H:78:ALA:N	2.32	0.45
1:C:221:ARG:HB3	1:C:222:LEU:H	1.50	0.45
1:C:20:SER:O	1:C:156:ARG:NH2	2.50	0.45
3:H:2:VAL:HB	3:H:102:TYR:CE1	2.51	0.45
4:L:45:LYS:HA	6:L:2005:HOH:O	2.17	0.45
4:L:124:GLN:HE22	4:L:131:SER:HB2	1.82	0.45
4:Q:79:GLN:HB2	4:Q:82:ASP:OD2	2.17	0.45
3:R:207:VAL:HG12	3:R:208:ASP:N	2.32	0.45
1:C:186:ASP:O	1:C:188:LYS:N	2.50	0.45
4:Q:161:GLU:OE1	4:Q:175:LEU:HD11	2.16	0.45
4:Q:24:ARG:HH11	4:Q:24:ARG:CB	2.29	0.45
4:Q:116:PHE:HD2	4:Q:135:LEU:HD22	1.82	0.45
4:Q:37:GLN:HB2	4:Q:47:LEU:HD11	1.98	0.44
1:A:57:HIS:NE2	1:A:195:SER:HB3	2.32	0.44
1:A:114:ARG:HG3	1:A:114:ARG:HH11	1.81	0.44
1:A:88:TYR:CE1	3:H:54:GLY:HA2	2.52	0.44
1:C:140:GLY:HA3	1:C:194:ASP:OD1	2.18	0.44
4:L:40:PRO:HG2	6:L:2003:HOH:O	2.18	0.44
4:L:143:GLU:CD	4:L:143:GLU:H	2.21	0.44
1:A:143:HIS:CD2	1:A:150:GLY:H	2.35	0.44
4:L:105:GLU:HB3	4:L:166:GLN:HE22	1.81	0.44
3:R:4:LEU:HD12	3:R:24:ALA:HA	1.99	0.44
3:R:98:TRP:C	3:R:100(A):ALA:N	2.71	0.44
1:C:166:HIS:HB2	1:C:167:LYS:HE3	1.99	0.44
3:R:178:LEU:HD12	3:R:179:SER:N	2.33	0.44
1:A:146:GLU:HB3	1:A:220:CYS:N	2.33	0.43
1:C:214:SER:HB3	1:C:215:TRP:HD1	1.82	0.43
3:R:30:ASN:HA	3:R:52(A):PRO:HB2	2.00	0.43
1:C:88:TYR:CE1	3:R:54:GLY:HA2	2.53	0.43
1:C:170:SER:HB3	1:C:170(B):GLU:OE1	2.18	0.43
3:H:117:LYS:HD2	3:H:144:ASP:O	2.18	0.43
1:A:51:TRP:CE3	1:A:105:LEU:HD22	2.53	0.43
4:L:135:LEU:C	4:L:136:LEU:HD12	2.39	0.43
4:Q:48:ILE:HD13	4:Q:54:LEU:HA	1.99	0.43
3:R:137:ALA:HB2	3:R:183:THR:HG22	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:54:SER:OG	1:A:55:ALA:N	2.50	0.43
1:C:25:GLY:O	1:C:28:PRO:HD3	2.18	0.43
4:Q:86:TYR:O	4:Q:101:GLY:HA2	2.19	0.43
4:Q:105:GLU:HB3	4:Q:166:GLN:HE22	1.82	0.43
3:R:6:GLU:OE2	3:R:91:TYR:HA	2.18	0.43
1:C:97:PHE:CE1	3:R:97:ALA:HB1	2.54	0.43
4:L:152:ASN:HD22	4:L:152:ASN:HA	1.59	0.43
4:L:198:HIS:CG	4:L:199:GLN:N	2.87	0.43
1:A:179:ASN:HA	1:A:233:ASN:ND2	2.33	0.43
1:C:171:VAL:HA	1:C:224:LYS:HE2	2.00	0.43
3:H:93:ALA:HB1	3:H:100:PHE:HB3	2.01	0.43
4:L:116:PHE:HD2	4:L:135:LEU:HD22	1.83	0.43
4:L:136:LEU:O	4:L:174:SER:HA	2.17	0.43
1:C:189:ASP:CG	1:C:190:ALA:H	2.22	0.43
1:C:213:ILE:HG12	1:C:228:TYR:CE2	2.53	0.43
4:L:70:ASP:OD2	4:Q:20:THR:HG21	2.18	0.43
3:R:2:VAL:HG13	3:R:27:PHE:CD1	2.53	0.43
3:R:197:ASN:N	3:R:197:ASN:ND2	2.66	0.43
4:L:113:PRO:HD2	4:L:201:LEU:HD13	2.01	0.43
4:Q:166:GLN:NE2	4:Q:173:TYR:OH	2.52	0.43
4:Q:4:MET:O	4:Q:99:GLY:HA2	2.19	0.42
3:R:14:PRO:HD3	3:R:112:SER:C	2.39	0.42
3:H:98:TRP:C	3:H:100(A):ALA:H	2.22	0.42
3:R:124:LEU:HB2	3:R:139:GLY:O	2.20	0.42
3:R:182:VAL:HG13	3:R:182:VAL:O	2.19	0.42
3:R:173:SER:O	3:R:175:LEU:HD12	2.19	0.42
4:L:35:TRP:CZ3	4:L:88:CYS:HB3	2.55	0.42
3:H:151:THR:O	3:H:152:VAL:HG23	2.19	0.42
4:Q:170:ASP:CG	4:Q:172:THR:HG23	2.40	0.42
1:A:96:VAL:HG13	1:A:97:PHE:CD2	2.54	0.42
1:A:97:PHE:HD1	3:H:98:TRP:CD2	2.38	0.42
1:C:60(A):HIS:CE1	1:C:60(C):PRO:HB3	2.55	0.42
1:C:188:LYS:HB3	1:C:188(A):SER:H	1.60	0.41
4:L:108:ARG:HD3	4:L:172:THR:HG22	2.02	0.41
4:Q:27:GLN:HB2	6:Q:2006:HOH:O	2.20	0.41
1:A:60(A):HIS:CE1	1:A:60(C):PRO:HB3	2.55	0.41
1:A:114:ARG:CB	1:A:119:GLN:HE22	2.33	0.41
1:A:199:LEU:HD22	1:A:228:TYR:CD2	2.56	0.41
1:A:214:SER:O	1:A:215:TRP:C	2.59	0.41
4:L:66:GLY:HA3	4:L:71:PHE:HA	2.02	0.41
3:H:137:ALA:HB2	3:H:183:THR:HG22	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:178:LEU:HD12	3:H:178:LEU:C	2.41	0.41
4:L:138:ASN:HA	4:L:173:TYR:O	2.21	0.41
4:Q:89:GLN:HE21	4:Q:90:GLN:C	2.23	0.41
4:Q:150:VAL:HG23	4:Q:155:GLN:HG3	2.02	0.41
3:R:119:PRO:HD3	3:R:200:HIS:ND1	2.35	0.41
1:C:86:GLU:O	1:C:87:LYS:HB3	2.21	0.41
1:C:182:CYS:HA	1:C:226:GLY:O	2.20	0.41
4:Q:167:ASP:HB3	4:Q:170:ASP:OD1	2.21	0.41
3:R:83:ARG:C	3:R:111:VAL:HG11	2.41	0.41
1:C:142:GLY:HA2	1:C:193:GLY:HA3	2.02	0.41
4:Q:138:ASN:ND2	4:Q:138:ASN:N	2.68	0.41
3:R:125:ALA:HA	3:R:126:PRO:HD2	1.96	0.41
4:L:6:GLN:HG2	4:L:88:CYS:SG	2.60	0.41
4:L:193:ALA:HB2	4:L:208:SER:HB3	2.02	0.41
3:R:143:LYS:HG3	3:R:177:SER:OG	2.21	0.41
1:A:214:SER:HB3	1:A:215:TRP:H	1.73	0.41
1:C:60(C):PRO:HA	1:C:60(D):PRO:HD3	1.93	0.41
1:C:211:GLY:HA2	1:C:229:THR:O	2.21	0.41
4:L:163:VAL:HG12	4:L:164:THR:N	2.36	0.41
4:Q:39:LYS:HG2	4:Q:84:ALA:HB2	2.03	0.41
4:Q:162:SER:OG	3:R:166:PHE:HB3	2.21	0.41
3:R:29:ILE:HG13	3:R:30:ASN:N	2.35	0.41
4:L:147:GLN:HE21	4:L:147:GLN:HB2	1.69	0.41
1:A:17:ILE:HD13	1:A:146:GLU:HG2	2.03	0.40
3:H:146:PHE:CE1	3:H:147:PRO:HB3	2.56	0.40
1:A:81:GLN:HB3	1:A:83:PHE:CE1	2.56	0.40
4:Q:163:VAL:HG12	4:Q:164:THR:N	2.37	0.40
3:R:119:PRO:HD2	3:R:205:THR:HB	2.02	0.40
4:Q:59:PRO:HG2	4:Q:62:PHE:HD2	1.86	0.40
4:Q:124:GLN:HE22	4:Q:131:SER:HB2	1.86	0.40
1:C:54:SER:OG	1:C:55:ALA:N	2.54	0.40
3:R:24:ALA:HB3	3:R:76:ASN:HD21	1.84	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	224/257 (87%)	186 (83%)	32 (14%)	6 (3%)	5	19
1	C	223/257 (87%)	186 (83%)	34 (15%)	3 (1%)	12	37
2	B	5/35 (14%)	2 (40%)	3 (60%)	0	100	100
2	D	4/35 (11%)	3 (75%)	1 (25%)	0	100	100
3	H	210/224 (94%)	195 (93%)	14 (7%)	1 (0%)	29	61
3	R	202/224 (90%)	186 (92%)	14 (7%)	2 (1%)	15	45
4	L	212/214 (99%)	190 (90%)	20 (9%)	2 (1%)	17	48
4	Q	212/214 (99%)	190 (90%)	21 (10%)	1 (0%)	29	61
All	All	1292/1460 (88%)	1138 (88%)	139 (11%)	15 (1%)	13	40

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
4	Q	138	ASN
1	A	148	VAL
1	C	187	CYS
4	L	138	ASN
1	A	171	VAL
1	A	173	GLY
1	C	186	ASP
4	L	213	GLU
1	A	214	SER
1	C	188	LYS
3	H	88	ALA
3	R	99	PRO
1	A	127	GLY
1	A	243	ARG
3	R	182	VAL

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	192/213 (90%)	177 (92%)	15 (8%)	12	34
1	C	191/213 (90%)	185 (97%)	6 (3%)	40	74
2	B	4/30 (13%)	4 (100%)	0	100	100
2	D	3/30 (10%)	2 (67%)	1 (33%)	0	0
3	H	171/181 (94%)	166 (97%)	5 (3%)	42	76
3	R	163/181 (90%)	157 (96%)	6 (4%)	34	68
4	L	186/186 (100%)	177 (95%)	9 (5%)	25	58
4	Q	186/186 (100%)	175 (94%)	11 (6%)	19	49
All	All	1096/1220 (90%)	1043 (95%)	53 (5%)	25	58

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	40	SER
1	A	45	SER
1	A	49	THR
1	A	76	THR
1	A	103	LEU
1	A	111	LYS
1	A	148	VAL
1	A	171	VAL
1	A	175	ASP
1	A	181	LEU
1	A	214	SER
1	A	220	CYS
1	A	228	TYR
1	A	240	ASP
1	A	243	ARG
1	C	40	SER
1	C	76	THR
1	C	111	LYS
1	C	115	SER
1	C	156	ARG
1	C	166	HIS
2	D	395	ARG
3	H	4	LEU
3	H	101	ASP
3	H	120	SER

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Mol	Chain	Res	Type
3	H	178	LEU
3	H	189	LEU
4	L	24	ARG
4	L	63	SER
4	L	125	LEU
4	L	129	THR
4	L	137	ASN
4	L	147	GLN
4	L	152	ASN
4	L	166	GLN
4	L	169	LYS
4	Q	24	ARG
4	Q	91	SER
4	Q	125	LEU
4	Q	129	THR
4	Q	137	ASN
4	Q	138	ASN
4	Q	147	GLN
4	Q	152	ASN
4	Q	166	GLN
4	Q	169	LYS
4	Q	201	LEU
3	R	4	LEU
3	R	64	LYS
3	R	101	ASP
3	R	135	THR
3	R	189	LEU
3	R	208	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (34) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	60(A)	HIS
1	A	71	HIS
1	A	116	GLN
1	A	119	GLN
1	A	233	ASN
1	A	239	ASN
2	B	397	HIS
1	C	60(A)	HIS
1	C	71	HIS
1	C	116	GLN

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Mol	Chain	Res	Type
1	C	119	GLN
3	H	3	GLN
3	H	30	ASN
3	H	192	GLN
3	H	204	ASN
4	L	27	GLN
4	L	124	GLN
4	L	137	ASN
4	L	138	ASN
4	L	147	GLN
4	L	152	ASN
4	L	160	GLN
4	Q	27	GLN
4	Q	124	GLN
4	Q	137	ASN
4	Q	138	ASN
4	Q	152	ASN
4	Q	158	ASN
3	R	3	GLN
3	R	30	ASN
3	R	164	HIS
3	R	192	GLN
3	R	197	ASN
3	R	204	ASN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

6 monosaccharides are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond

length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
5	NAG	E	1	5,1	14,14,15	0.58	0	17,19,21	1.05	1 (5%)
5	NAG	E	2	5	14,14,15	0.76	1 (7%)	17,19,21	1.95	4 (23%)
5	BMA	E	3	5	11,11,12	0.56	0	15,15,17	1.20	2 (13%)
5	NAG	F	1	5,1	14,14,15	0.42	0	17,19,21	1.96	3 (17%)
5	NAG	F	2	5	14,14,15	0.65	0	17,19,21	1.41	3 (17%)
5	BMA	F	3	5	11,11,12	0.72	0	15,15,17	1.49	1 (6%)

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

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	E	2	NAG	C1-C2	2.03	1.55	1.52

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
5	F	1	NAG	C1-O5-C5	5.97	120.27	112.19
5	F	3	BMA	C1-O5-C5	4.90	118.84	112.19
5	E	2	NAG	C1-O5-C5	4.75	118.63	112.19
5	E	2	NAG	O4-C4-C5	3.80	118.73	109.30
5	F	1	NAG	C4-C3-C2	-3.51	105.88	111.02
5	E	3	BMA	C1-O5-C5	3.21	116.54	112.19
5	F	2	NAG	C4-C3-C2	3.12	115.60	111.02
5	E	2	NAG	O5-C1-C2	2.92	115.91	111.29

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
5	F	1	NAG	O5-C5-C6	2.56	111.21	107.20
5	E	1	NAG	C1-O5-C5	2.42	115.47	112.19
5	F	2	NAG	O4-C4-C5	2.36	115.17	109.30
5	E	2	NAG	C2-N2-C7	2.26	126.12	122.90
5	E	3	BMA	C1-C2-C3	2.12	112.27	109.67
5	F	2	NAG	O5-C5-C4	-2.06	105.81	110.83

There are no chirality outliers.

All (15) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	E	1	NAG	C8-C7-N2-C2
5	E	1	NAG	O7-C7-N2-C2
5	E	2	NAG	C3-C2-N2-C7
5	E	2	NAG	C8-C7-N2-C2
5	E	2	NAG	O7-C7-N2-C2
5	F	1	NAG	C8-C7-N2-C2
5	F	1	NAG	O7-C7-N2-C2
5	F	2	NAG	C8-C7-N2-C2
5	F	2	NAG	O7-C7-N2-C2
5	F	2	NAG	O5-C5-C6-O6
5	F	2	NAG	C4-C5-C6-O6
5	F	1	NAG	O5-C5-C6-O6
5	F	1	NAG	C4-C5-C6-O6
5	E	2	NAG	C4-C5-C6-O6
5	E	2	NAG	O5-C5-C6-O6

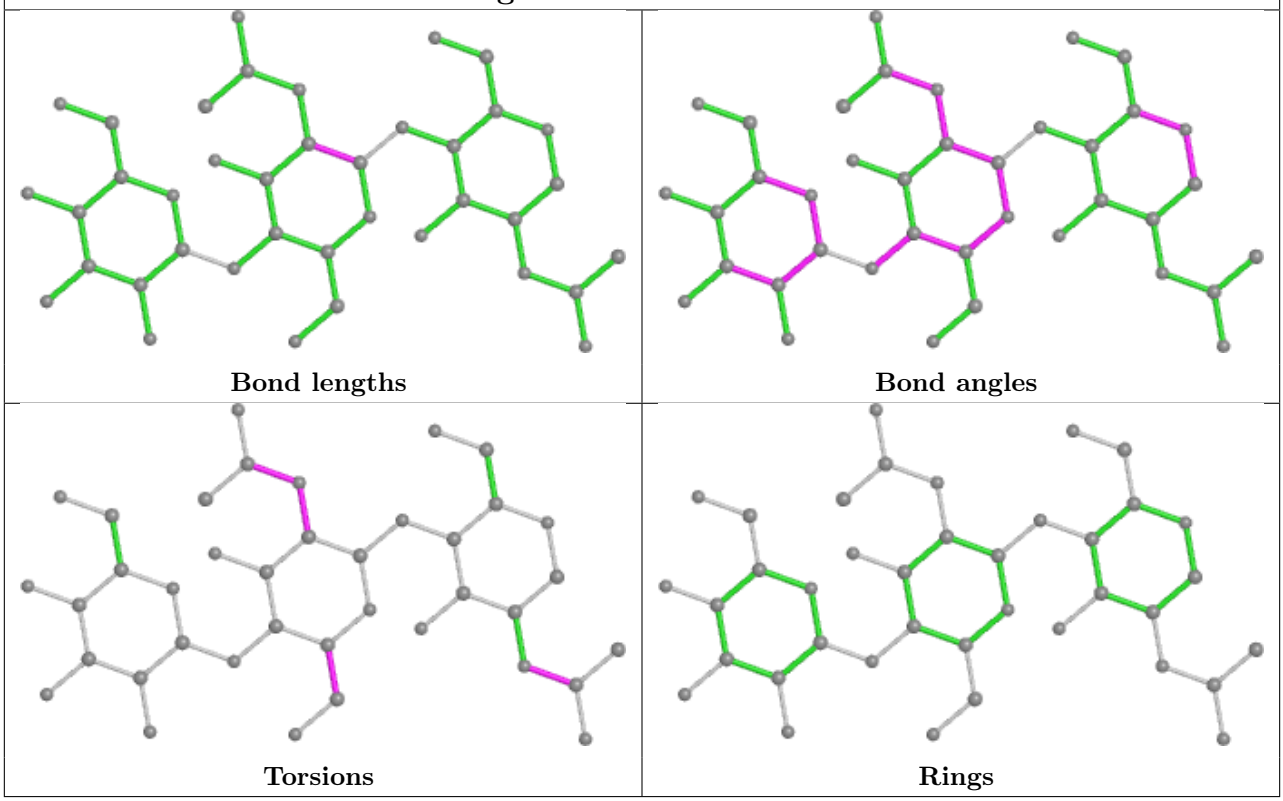
There are no ring outliers.

1 monomer is involved in 1 short contact:

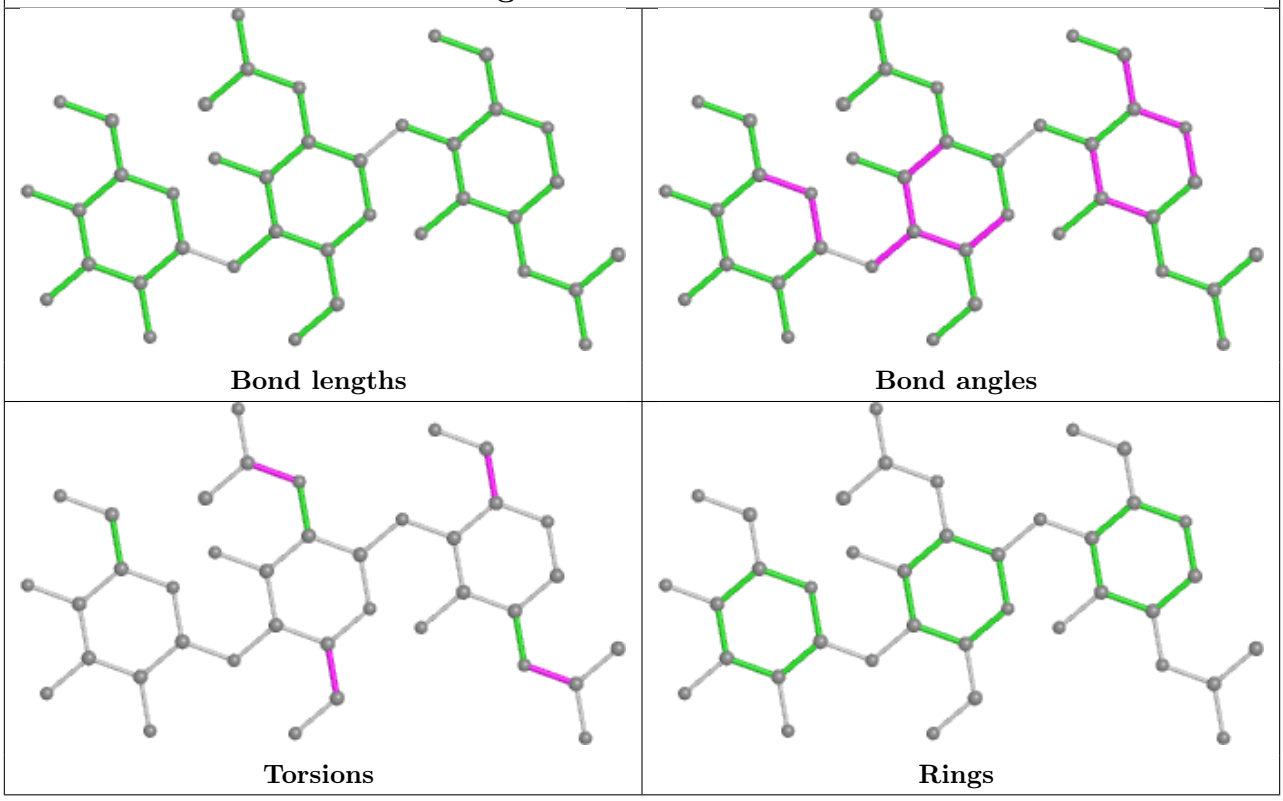
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	E	3	BMA	1	0

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



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)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	40:SER	C	41:PHE	N	1.18

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	232/257 (90%)	-0.16	4 (1%) 70 69	38, 71, 122, 133	0
1	C	229/257 (89%)	-0.19	4 (1%) 70 69	36, 74, 122, 139	0
2	B	7/35 (20%)	0.07	0 100 100	77, 84, 102, 104	0
2	D	6/35 (17%)	1.23	1 (16%) 1 1	83, 92, 107, 109	0
3	H	214/224 (95%)	-0.36	2 (0%) 84 84	33, 53, 99, 132	0
3	R	206/224 (91%)	-0.01	13 (6%) 20 16	34, 55, 123, 140	0
4	L	214/214 (100%)	-0.43	0 100 100	40, 63, 108, 124	0
4	Q	214/214 (100%)	0.01	10 (4%) 31 28	36, 76, 133, 142	0
All	All	1322/1460 (90%)	-0.18	34 (2%) 56 52	33, 64, 123, 142	0

All (34) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	R	189	LEU	6.9
3	R	190	GLY	5.5
3	R	138	LEU	5.0
3	R	124	LEU	5.0
3	R	123	PRO	4.3
3	R	122	PHE	4.1
4	Q	186	TYR	4.0
4	Q	152	ASN	3.4
4	Q	182	SER	3.2
4	Q	117	ILE	3.1
1	C	223	HIS	3.1
2	D	397	HIS	3.0
4	Q	119	PRO	3.0
3	H	217	ASP	3.0
3	R	183	THR	2.9
4	Q	154	LEU	2.9

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Mol	Chain	Res	Type	RSRZ
3	R	194	TYR	2.8
3	R	126	PRO	2.8
3	R	139	GLY	2.7
1	C	72	PHE	2.7
3	R	159	LEU	2.7
4	Q	120	PRO	2.5
4	Q	181	LEU	2.5
1	A	244	PRO	2.4
3	R	181	VAL	2.4
4	Q	121	SER	2.4
1	C	73	PHE	2.4
1	A	144	LEU	2.3
1	A	174	ALA	2.3
4	Q	118	PHE	2.2
3	H	191	THR	2.1
1	C	152	SER	2.1
3	R	182	VAL	2.1
1	A	167	LYS	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)

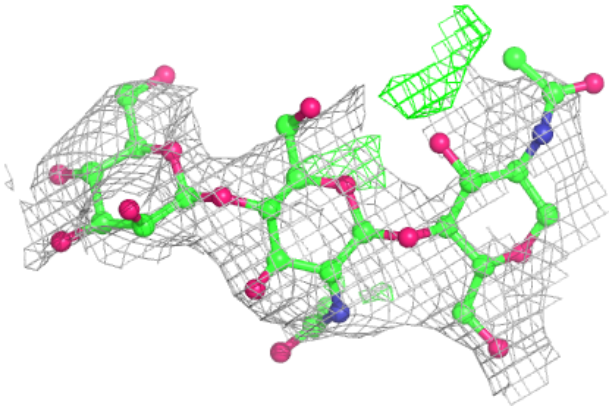
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

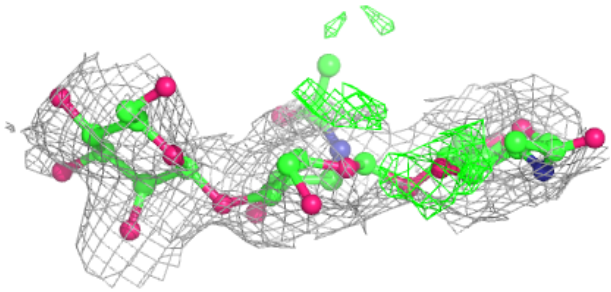
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
5	BMA	F	3	11/12	0.38	0.34	142,143,143,143	0
5	NAG	F	2	14/15	0.75	0.26	136,139,140,143	0
5	BMA	E	3	11/12	0.75	0.21	153,153,153,153	0
5	NAG	E	1	14/15	0.79	0.17	125,132,136,141	0
5	NAG	F	1	14/15	0.84	0.22	119,123,127,132	0
5	NAG	E	2	14/15	0.85	0.20	146,148,150,153	0

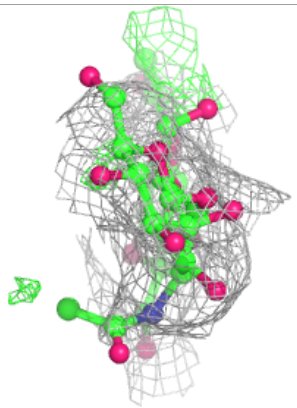
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

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)

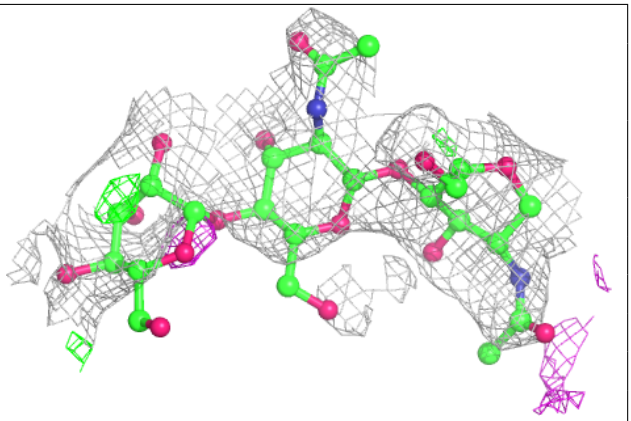


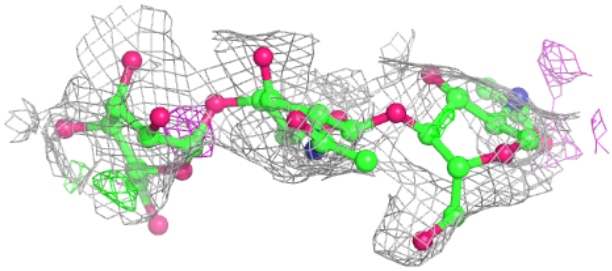


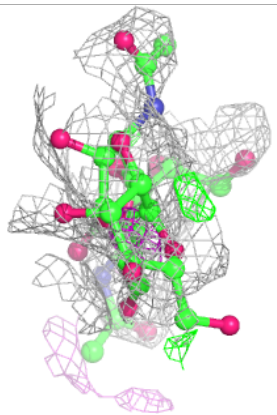


Electron density around Chain F:

$2mF_o$-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.