



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 08:17 PM GMT

PDB ID : 1JMJ
Title : Crystal Structure of Native Heparin Cofactor II
Authors : Baglin, T.P.; Carrell, R.W.; Church, F.C.; Huntington, J.A.
Deposited on : 2001-07-18
Resolution : 2.35 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
<http://wwpdb.org/validation/2016/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 : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

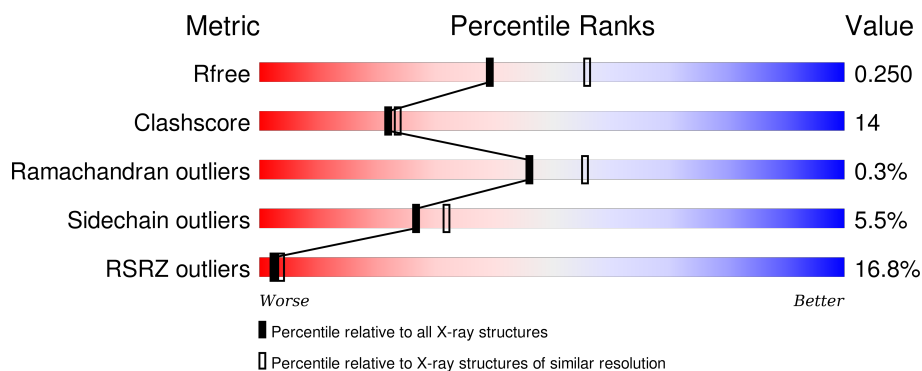
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.35 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}	91344	1352 (2.38-2.34)
Clashscore	102246	1456 (2.38-2.34)
Ramachandran outliers	100387	1435 (2.38-2.34)
Sidechain outliers	100360	1436 (2.38-2.34)
RSRZ outliers	91569	1358 (2.38-2.34)

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. 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	480	
1	B	480	

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 6723 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.

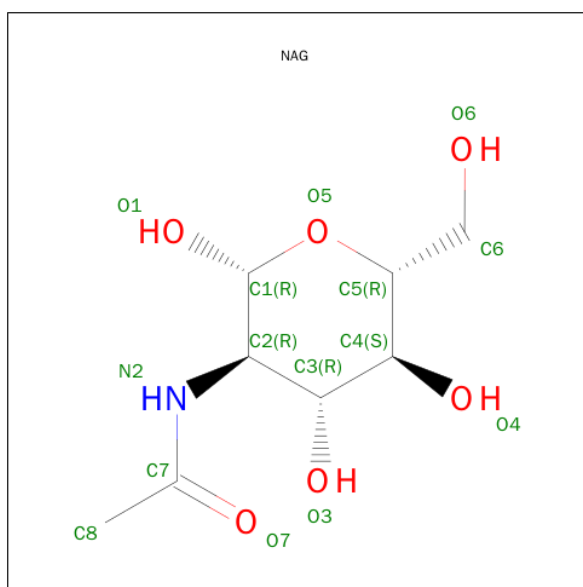
- Molecule 1 is a protein called HEPARIN COFACTOR II.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	397	Total	C	N	O	S	20	0	0
			3228	2069	560	579	20			
1	B	387	Total	C	N	O	S	14	0	0
			3144	2017	549	558	20			

- Molecule 2 is a polymer of unknown type called SUGAR (3-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	3	Total	C	N	O	0	0
			39	22	2	15		

- Molecule 3 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



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

- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total 1	Ca 1	0	0

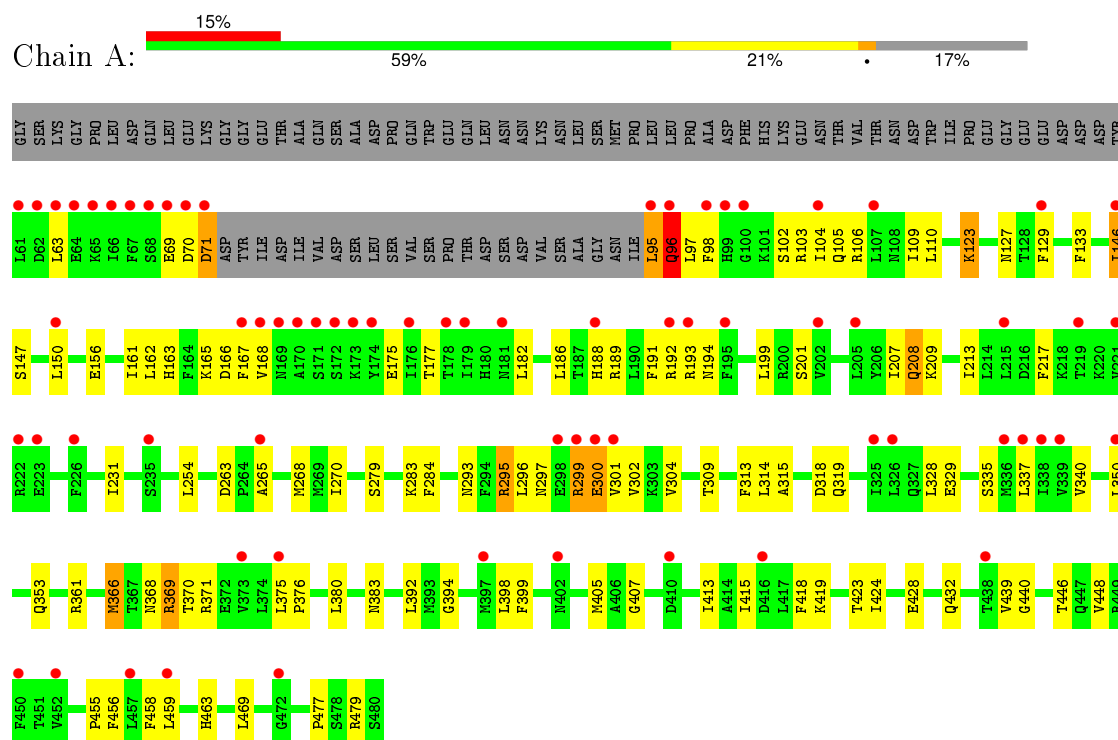
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	157	Total 157	O 157	0	0
5	B	140	Total 140	O 140	0	0

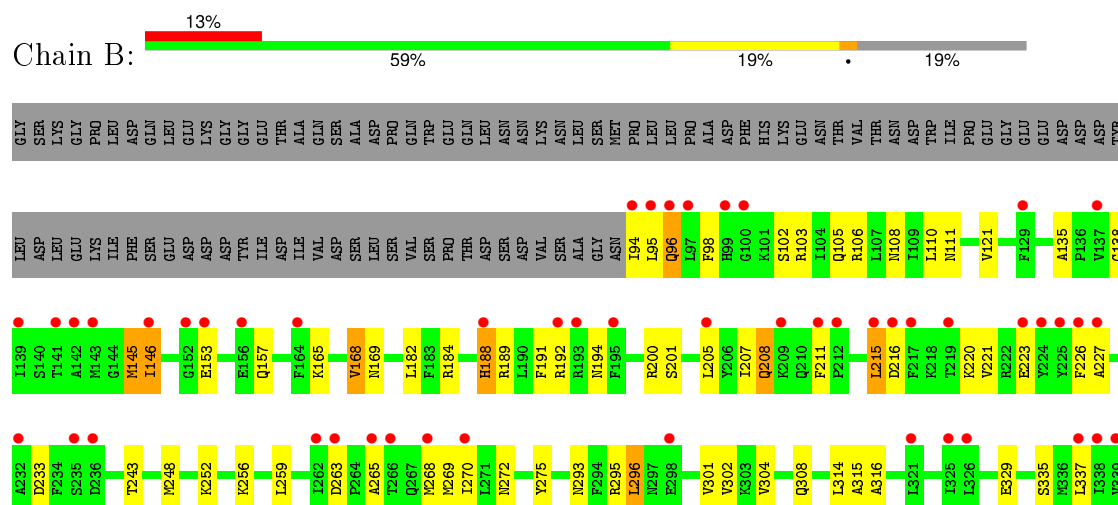
3 Residue-property plots

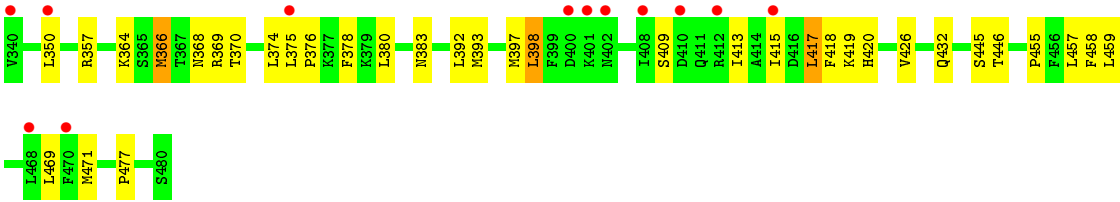
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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: HEPARIN COFACTOR II



• Molecule 1: HEPARIN COFACTOR II





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	74.73Å 80.00Å 92.24Å 90.00° 102.04° 90.00°	Depositor
Resolution (Å)	25.00 – 2.35 25.05 – 2.35	Depositor EDS
% Data completeness (in resolution range)	88.0 (25.00-2.35) 88.1 (25.05-2.35)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	0.14	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.57 (at 2.36Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.208 , 0.255 0.209 , 0.250	Depositor DCC
R_{free} test set	994 reflections (2.54%)	DCC
Wilson B-factor (Å ²)	31.6	Xtriage
Anisotropy	0.624	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 54.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 39171 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6723	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.12% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CA, 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	0.36	0/3291	0.62	0/4431
1	B	0.35	0/3207	0.63	0/4320
All	All	0.36	0/6498	0.62	0/8751

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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	3228	0	3279	95	0
1	B	3144	0	3207	84	0
2	A	39	0	34	4	0
3	B	14	0	13	0	0
4	A	1	0	0	0	0
5	A	157	0	0	2	0
5	B	140	0	0	1	0
All	All	6723	0	6533	179	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 (179) 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:314:LEU:HB3	1:A:366:MET:HG2	1.48	0.96
1:B:96:GLN:H	1:B:96:GLN:NE2	1.64	0.95
1:A:405:MET:HE1	1:A:413:ILE:HG13	1.52	0.92
1:A:110:LEU:HD12	1:A:167:PHE:HB2	1.59	0.84
2:A:482:NAG:H3	2:A:482:NAG:H82	1.61	0.83
1:A:329:GLU:OE1	1:A:335:SER:HB3	1.80	0.82
1:B:221:VAL:HG13	1:B:227:ALA:HB3	1.60	0.82
1:B:192:ARG:HG3	1:B:192:ARG:HH11	1.43	0.82
1:A:70:ASP:HB3	5:A:631:HOH:O	1.82	0.80
1:B:314:LEU:HB3	1:B:366:MET:HG2	1.65	0.78
1:B:135:ALA:H	1:B:420:HIS:CE1	2.03	0.77
1:A:95:LEU:O	1:A:96:GLN:HB3	1.87	0.73
1:B:398:LEU:HD11	1:B:415:ILE:HD12	1.70	0.73
1:A:370:THR:HG22	1:A:446:THR:HG22	1.69	0.72
1:B:329:GLU:OE2	1:B:335:SER:HB3	1.90	0.71
1:B:295:ARG:O	1:B:455:PRO:HD3	1.90	0.71
1:B:96:GLN:H	1:B:96:GLN:HE21	1.38	0.70
1:B:350:LEU:HD11	1:B:457:LEU:HD12	1.72	0.70
1:B:215:LEU:HD13	1:B:216:ASP:N	2.08	0.69
1:A:293:ASN:HB3	1:A:301:VAL:HG11	1.74	0.68
1:B:145:MET:HG2	1:B:221:VAL:HG22	1.75	0.67
1:B:304:VAL:HG11	1:B:477:PRO:HG2	1.77	0.67
1:A:63:LEU:HD22	1:A:63:LEU:H	1.63	0.64
1:A:375:LEU:HD12	1:A:376:PRO:HD2	1.79	0.64
1:A:95:LEU:HD22	1:A:98:PHE:CD1	2.32	0.64
1:B:216:ASP:OD1	1:B:220:LYS:HE2	1.98	0.63
1:B:207:ILE:HD13	1:B:268:MET:HG3	1.80	0.63
1:A:69:GLU:OE2	1:B:445:SER:HB2	1.98	0.63
1:A:95:LEU:O	1:A:96:GLN:CB	2.47	0.63
1:B:375:LEU:HD12	1:B:376:PRO:HD2	1.81	0.63
1:A:405:MET:HE1	1:A:413:ILE:CG1	2.26	0.62
1:B:146:ILE:HG13	1:B:270:ILE:HD11	1.80	0.62
1:B:138:GLY:C	1:B:272:ASN:HD21	2.03	0.62
1:B:98:PHE:CD2	1:B:106:ARG:HD3	2.34	0.62
1:B:252:LYS:HD2	1:B:275:TYR:OH	2.00	0.61
1:B:98:PHE:HZ	1:B:110:LEU:HD21	1.63	0.61
1:B:94:ILE:O	1:B:357:ARG:NH2	2.31	0.61
1:A:405:MET:CE	1:A:413:ILE:HG13	2.26	0.61
1:B:350:LEU:CD1	1:B:457:LEU:HD12	2.31	0.60
1:B:94:ILE:HG22	1:B:95:LEU:HG	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:153:GLU:O	1:B:157:GLN:HG3	2.03	0.59
1:B:135:ALA:H	1:B:420:HIS:HE1	1.46	0.58
1:A:96:GLN:HE21	1:A:96:GLN:HA	1.66	0.58
1:A:370:THR:CG2	1:A:446:THR:HG22	2.32	0.58
1:B:208:GLN:HA	1:B:208:GLN:OE1	2.03	0.58
1:B:192:ARG:CG	1:B:192:ARG:HH11	2.15	0.58
1:B:296:LEU:HD22	1:B:302:VAL:HG13	1.84	0.58
1:B:272:ASN:HD22	1:B:420:HIS:HD2	1.50	0.58
1:B:184:ARG:HG3	1:B:188:HIS:CE1	2.38	0.57
1:A:304:VAL:HG11	1:A:477:PRO:HG2	1.85	0.57
1:A:63:LEU:HD22	1:A:63:LEU:N	2.19	0.56
1:A:102:SER:OG	1:A:104:ILE:HG22	2.05	0.56
1:B:96:GLN:N	1:B:96:GLN:NE2	2.46	0.56
1:A:299:ARG:HB2	1:A:299:ARG:NH1	2.20	0.56
1:B:192:ARG:HG3	1:B:192:ARG:NH1	2.19	0.56
1:A:146:ILE:HG13	1:A:270:ILE:HD11	1.87	0.56
1:A:209:LYS:HB2	1:A:231:ILE:HD11	1.87	0.56
1:A:299:ARG:O	1:A:300:GLU:HB3	2.05	0.56
1:A:70:ASP:O	1:A:71:ASP:HB2	2.06	0.55
1:B:308:GLN:HG3	1:B:374:LEU:HG	1.88	0.55
1:A:405:MET:HE1	1:A:413:ILE:O	2.06	0.55
1:B:417:LEU:HD11	1:B:419:LYS:HB2	1.87	0.55
1:A:95:LEU:O	1:A:95:LEU:HG	2.05	0.55
1:B:95:LEU:HD22	1:B:98:PHE:CE1	2.41	0.55
1:A:192:ARG:HH11	1:A:192:ARG:HG3	1.73	0.54
1:A:188:HIS:O	1:A:192:ARG:HG2	2.08	0.54
1:A:168:VAL:HG13	1:A:175:GLU:HA	1.90	0.54
1:B:398:LEU:CD1	1:B:415:ILE:HD12	2.38	0.53
1:B:205:LEU:HD12	1:B:269:MET:O	2.09	0.53
1:B:370:THR:HG22	1:B:446:THR:HG22	1.91	0.53
1:A:147:SER:HA	1:A:150:LEU:HG	1.91	0.53
1:A:337:LEU:O	1:A:458:PHE:HA	2.09	0.53
1:A:398:LEU:HD13	1:A:398:LEU:O	2.09	0.53
1:A:368:ASN:HD22	2:A:481:NAG:H83	1.74	0.52
1:A:263:ASP:OD1	1:A:265:ALA:HB3	2.09	0.52
1:B:102:SER:O	1:B:106:ARG:HG3	2.10	0.52
1:A:110:LEU:CD1	1:A:167:PHE:HB2	2.37	0.51
1:A:405:MET:HE1	1:A:413:ILE:C	2.31	0.51
1:A:335:SER:OG	1:A:463:HIS:HE1	1.92	0.51
1:B:96:GLN:N	1:B:96:GLN:HE21	2.05	0.51
1:B:184:ARG:O	1:B:188:HIS:ND1	2.44	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:175:GLU:HG3	1:A:177:THR:CG2	2.41	0.50
1:B:94:ILE:HG22	1:B:95:LEU:N	2.26	0.50
1:A:295:ARG:O	1:A:455:PRO:HD3	2.11	0.50
1:A:208:GLN:HE21	1:A:208:GLN:CA	2.24	0.50
1:B:208:GLN:OE1	1:B:233:ASP:HA	2.11	0.50
1:A:69:GLU:CD	1:B:445:SER:HB2	2.32	0.50
1:A:209:LYS:CB	1:A:231:ILE:HD11	2.42	0.50
1:B:316:ALA:N	1:B:366:MET:HE2	2.27	0.49
1:B:184:ARG:HD3	1:B:223:GLU:O	2.11	0.49
1:A:299:ARG:HH11	1:A:299:ARG:CB	2.26	0.49
1:A:104:ILE:HG23	1:A:105:GLN:N	2.26	0.49
1:A:254:LEU:HD21	1:A:423:THR:HB	1.95	0.49
1:B:188:HIS:CE1	1:B:226:PHE:CZ	3.00	0.49
1:B:368:ASN:O	1:B:369:ARG:HD2	2.12	0.49
1:A:313:PHE:HB2	1:A:371:ARG:HD2	1.94	0.49
1:B:215:LEU:O	1:B:215:LEU:HD22	2.12	0.48
1:B:296:LEU:HD22	1:B:302:VAL:CG1	2.43	0.48
1:A:315:ALA:HB3	1:A:371:ARG:CZ	2.43	0.48
1:A:439:VAL:HG22	1:A:440:GLY:N	2.28	0.48
2:A:482:NAG:C3	2:A:482:NAG:H82	2.39	0.48
1:A:371:ARG:HG2	1:A:448:VAL:CG2	2.43	0.48
1:A:314:LEU:HD23	1:A:368:ASN:HA	1.96	0.48
2:A:482:NAG:H3	2:A:482:NAG:C8	2.39	0.48
1:A:279:SER:HB2	1:A:283:LYS:NZ	2.29	0.48
1:B:188:HIS:O	1:B:192:ARG:HG2	2.14	0.48
1:A:123:LYS:HD2	5:A:605:HOH:O	2.13	0.48
1:A:63:LEU:H	1:A:63:LEU:CD2	2.27	0.47
1:B:153:GLU:HB3	1:B:397:MET:CE	2.45	0.47
1:A:103:ARG:HG2	1:A:182:LEU:HD11	1.97	0.47
1:A:104:ILE:HD11	1:A:186:LEU:HB2	1.96	0.47
1:A:217:PHE:HB2	1:A:407:GLY:O	2.15	0.47
1:B:145:MET:HE1	1:B:146:ILE:HG12	1.96	0.46
1:B:383:ASN:HD21	1:B:419:LYS:HE3	1.80	0.46
1:A:318:ASP:OD2	1:A:361:ARG:NH2	2.40	0.46
1:A:192:ARG:NH1	1:A:194:ASN:OD1	2.48	0.46
1:A:127:ASN:HB3	1:A:129:PHE:CE1	2.50	0.46
1:A:405:MET:CE	1:A:413:ILE:O	2.64	0.46
1:A:268:MET:HG2	1:A:415:ILE:HD13	1.98	0.46
1:B:293:ASN:HB3	1:B:301:VAL:HG11	1.97	0.46
1:A:424:ILE:HA	1:A:432:GLN:HE22	1.80	0.46
1:A:368:ASN:O	1:A:369:ARG:HD3	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:162:LEU:O	1:A:163:HIS:HB2	2.16	0.45
1:A:297:ASN:ND2	1:A:300:GLU:HG2	2.31	0.45
1:B:248:MET:SD	1:B:256:LYS:HG2	2.56	0.45
1:A:207:ILE:HD12	1:A:213:ILE:HG12	1.97	0.45
1:A:192:ARG:CG	1:A:192:ARG:HH11	2.29	0.45
1:B:192:ARG:HH12	1:B:194:ASN:CG	2.19	0.45
1:B:357:ARG:HH11	1:B:357:ARG:HG2	1.82	0.45
1:A:268:MET:HE2	1:A:415:ILE:HD13	1.98	0.45
1:A:95:LEU:HD23	1:A:97:LEU:HB2	1.99	0.45
1:A:105:GLN:O	1:A:109:ILE:HG13	2.17	0.44
1:A:284:PHE:O	1:A:428:GLU:HB3	2.16	0.44
1:A:133:PHE:CZ	1:A:424:ILE:HB	2.52	0.44
1:A:156:GLU:OE1	1:A:165:LYS:NZ	2.40	0.44
1:A:370:THR:HG22	1:A:446:THR:HA	1.99	0.44
1:B:98:PHE:HZ	1:B:110:LEU:CD2	2.30	0.44
1:B:215:LEU:C	1:B:215:LEU:HD22	2.38	0.44
1:A:398:LEU:HD12	1:A:399:PHE:CE2	2.53	0.44
1:A:315:ALA:HB3	1:A:371:ARG:NH1	2.34	0.43
1:B:103:ARG:HG2	1:B:182:LEU:HD11	2.00	0.43
1:B:458:PHE:O	1:B:471:MET:HA	2.18	0.43
1:A:95:LEU:CD2	1:A:97:LEU:HB2	2.49	0.43
1:A:161:ILE:HD11	1:A:394:GLY:HA3	2.00	0.43
1:B:108:ASN:HA	1:B:111:ASN:HB2	2.00	0.43
1:B:102:SER:OG	1:B:105:GLN:HG3	2.19	0.43
1:A:301:VAL:HG12	1:A:302:VAL:N	2.32	0.43
1:B:378:PHE:CZ	1:B:426:VAL:HG21	2.53	0.43
1:B:200:ARG:NH2	1:B:252:LYS:HE3	2.34	0.42
1:A:340:VAL:HG22	1:A:456:PHE:HB3	2.01	0.42
1:A:299:ARG:HB2	1:A:299:ARG:HH11	1.82	0.42
1:B:191:PHE:CZ	1:B:201:SER:HB2	2.54	0.42
1:A:104:ILE:CG2	1:A:105:GLN:N	2.82	0.42
1:B:409:SER:HB3	1:B:413:ILE:HD11	2.02	0.42
1:B:315:ALA:C	1:B:366:MET:HG3	2.41	0.41
1:A:98:PHE:CD2	1:A:106:ARG:HD3	2.55	0.41
1:A:208:GLN:HA	1:A:208:GLN:NE2	2.35	0.41
1:A:129:PHE:HD2	1:A:479:ARG:NH1	2.18	0.41
1:A:191:PHE:CZ	1:A:201:SER:HB2	2.55	0.41
1:B:145:MET:HE1	1:B:146:ILE:HA	2.01	0.41
1:B:145:MET:CE	1:B:146:ILE:HA	2.51	0.41
1:A:284:PHE:HA	1:A:309:THR:HB	2.02	0.41
1:B:121:VAL:HG21	1:B:393:MET:HG2	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:263:ASP:OD1	1:B:265:ALA:HB3	2.20	0.41
1:B:337:LEU:O	1:B:458:PHE:HA	2.20	0.41
1:B:165:LYS:O	1:B:169:ASN:HB2	2.19	0.41
1:A:383:ASN:OD1	1:A:419:LYS:HE3	2.21	0.41
1:A:175:GLU:HG3	1:A:177:THR:HG23	2.02	0.41
1:B:243:THR:HG21	1:B:259:LEU:CD1	2.51	0.41
1:A:319:GLN:HG3	1:B:364:LYS:O	2.21	0.41
1:A:268:MET:CE	1:A:415:ILE:HD13	2.50	0.41
1:A:424:ILE:HG23	1:A:424:ILE:O	2.20	0.41
1:B:192:ARG:CG	1:B:192:ARG:NH1	2.74	0.40
1:A:192:ARG:O	1:A:193:ARG:C	2.59	0.40
1:B:168:VAL:CG2	1:B:169:ASN:N	2.84	0.40
1:A:370:THR:CG2	1:A:446:THR:HA	2.51	0.40
1:B:191:PHE:CE1	1:B:201:SER:HB2	2.56	0.40
1:B:211:PHE:CD1	1:B:211:PHE:N	2.89	0.40
1:B:189:ARG:HB2	5:B:576:HOH:O	2.22	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	393/480 (82%)	374 (95%)	17 (4%)	2 (0%)	34	39
1	B	385/480 (80%)	375 (97%)	10 (3%)	0	100	100
All	All	778/960 (81%)	749 (96%)	27 (4%)	2 (0%)	46	55

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	96	GLN
1	A	300	GLU

5.3.2 Protein sidechains ⓘ

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	360/433 (83%)	338 (94%)	22 (6%)	23	27
1	B	350/433 (81%)	333 (95%)	17 (5%)	31	39
All	All	710/866 (82%)	671 (94%)	39 (6%)	27	32

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	71	ASP
1	A	95	LEU
1	A	96	GLN
1	A	123	LYS
1	A	146	ILE
1	A	166	ASP
1	A	189	ARG
1	A	199	LEU
1	A	208	GLN
1	A	295	ARG
1	A	296	LEU
1	A	299	ARG
1	A	328	LEU
1	A	350	LEU
1	A	353	GLN
1	A	366	MET
1	A	369	ARG
1	A	380	LEU
1	A	392	LEU
1	A	418	PHE
1	A	459	LEU
1	A	469	LEU
1	B	96	GLN
1	B	145	MET
1	B	146	ILE
1	B	168	VAL
1	B	188	HIS

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Mol	Chain	Res	Type
1	B	208	GLN
1	B	215	LEU
1	B	296	LEU
1	B	366	MET
1	B	380	LEU
1	B	392	LEU
1	B	398	LEU
1	B	417	LEU
1	B	418	PHE
1	B	432	GLN
1	B	459	LEU
1	B	469	LEU

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

Mol	Chain	Res	Type
1	A	96	GLN
1	A	208	GLN
1	A	292	HIS
1	A	363	GLN
1	A	463	HIS
1	B	96	GLN
1	B	169	ASN
1	B	272	ASN
1	B	411	GLN
1	B	420	HIS

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](#)

3 carbohydrates are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
2	NAG	A	481	1,2	14,14,15	0.58	0	15,19,21	0.74	0
2	NAG	A	482	2	14,14,15	0.59	0	15,19,21	0.97	1 (6%)
2	BMA	A	483	2	11,11,12	0.44	0	14,15,17	0.42	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	A	481	1,2	-	0/6/23/26	0/1/1/1
2	NAG	A	482	2	-	2/6/23/26	0/1/1/1
2	BMA	A	483	2	-	0/2/19/22	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($^{\circ}$)	Ideal($^{\circ}$)
2	A	482	NAG	C3-C4-C5	2.31	114.23	110.20

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	482	NAG	C8-C7-N2-C2
2	A	482	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
2	A	481	NAG	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	482	NAG	3	0

5.6 Ligand geometry

Of 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 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	$\# Z > 2$
3	NAG	B	481	1	14,14,15	0.55	0	15,19,21	0.63	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	B	481	1	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	481	NAG	C8-C7-N2-C2
3	B	481	NAG	O7-C7-N2-C2

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	397/480 (82%)	1.15	70 (17%) 2 3	19, 34, 69, 95	7 (1%)
1	B	387/480 (80%)	1.02	62 (16%) 3 4	17, 34, 60, 85	4 (1%)
All	All	784/960 (81%)	1.08	132 (16%) 2 4	17, 34, 62, 95	11 (1%)

All (132) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	215	LEU	7.4
1	B	94	ILE	7.4
1	A	95	LEU	7.3
1	A	61	LEU	7.3
1	A	215	LEU	7.0
1	A	172	SER	6.7
1	B	402	ASN	6.1
1	A	99	HIS	6.1
1	A	62	ASP	6.0
1	A	193	ARG	5.4
1	A	96	GLN	5.4
1	B	96	GLN	5.2
1	B	219	THR	5.2
1	A	65	LYS	5.0
1	B	195	PHE	5.0
1	A	170	ALA	4.9
1	B	99	HIS	4.8
1	A	71	ASP	4.8
1	A	402	ASN	4.7
1	A	171	SER	4.7
1	B	226	PHE	4.6
1	A	70	ASP	4.5
1	A	173	LYS	4.4
1	B	216	ASP	4.4

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Mol	Chain	Res	Type	RSRZ
1	B	188	HIS	4.3
1	A	195	PHE	4.3
1	B	193	ARG	4.2
1	A	168	VAL	4.1
1	A	174	TYR	4.0
1	A	66	ILE	4.0
1	A	169	ASN	4.0
1	A	98	PHE	4.0
1	A	64	GLU	3.9
1	A	100	GLY	3.9
1	B	265	ALA	3.9
1	B	146	ILE	3.9
1	B	298	GLU	3.8
1	B	410	ASP	3.8
1	A	223	GLU	3.5
1	A	298	GLU	3.4
1	B	235	SER	3.4
1	A	339	VAL	3.3
1	A	188	HIS	3.3
1	B	156	GLU	3.3
1	A	192	ARG	3.2
1	A	235	SER	3.2
1	B	401	LYS	3.2
1	B	129	PHE	3.1
1	A	265	ALA	3.0
1	B	263	ASP	3.0
1	A	452	VAL	2.9
1	A	326	LEU	2.9
1	A	375	LEU	2.9
1	A	221	VAL	2.9
1	B	412	ARG	2.9
1	B	212	PRO	2.9
1	B	270	ILE	2.9
1	A	129	PHE	2.8
1	A	338	ILE	2.8
1	A	68	SER	2.8
1	A	63	LEU	2.8
1	B	339	VAL	2.8
1	B	338	ILE	2.7
1	A	205	LEU	2.7
1	B	139	ILE	2.7
1	B	350	LEU	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	222	ARG	2.7
1	A	337	LEU	2.6
1	B	325	ILE	2.6
1	A	459	LEU	2.6
1	A	336	MET	2.6
1	A	410	ASP	2.6
1	A	179	ILE	2.6
1	A	146	ILE	2.6
1	A	176	ILE	2.6
1	B	408	ILE	2.6
1	B	192	ARG	2.6
1	A	104	ILE	2.5
1	B	266	THR	2.5
1	A	69	GLU	2.5
1	B	223	GLU	2.5
1	B	375	LEU	2.5
1	B	153	GLU	2.5
1	A	325	ILE	2.5
1	B	217	PHE	2.5
1	B	340	VAL	2.4
1	A	350	LEU	2.4
1	B	95	LEU	2.4
1	B	209	LYS	2.4
1	A	472	GLY	2.4
1	A	301	VAL	2.4
1	B	211	PHE	2.4
1	B	141	THR	2.3
1	A	167	PHE	2.3
1	B	143	MET	2.3
1	B	262	ILE	2.3
1	A	67	PHE	2.3
1	A	373	VAL	2.3
1	B	225	TYR	2.3
1	B	227	ALA	2.3
1	B	400	ASP	2.3
1	B	321	LEU	2.3
1	A	299	ARG	2.3
1	A	178	THR	2.2
1	B	137	VAL	2.2
1	A	202	VAL	2.2
1	A	450	PHE	2.2
1	A	219	THR	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	224	TYR	2.2
1	B	326	LEU	2.2
1	B	415	ILE	2.2
1	B	100	GLY	2.2
1	B	152	GLY	2.2
1	A	107	LEU	2.2
1	A	457	LEU	2.2
1	B	205	LEU	2.2
1	B	268	MET	2.2
1	B	142	ALA	2.2
1	B	337	LEU	2.1
1	B	468	LEU	2.1
1	B	470	PHE	2.1
1	A	300	GLU	2.1
1	B	236	ASP	2.1
1	B	164	PHE	2.1
1	B	232	ALA	2.1
1	A	416	ASP	2.1
1	B	97	LEU	2.0
1	A	181	ASN	2.0
1	A	438	THR	2.0
1	A	150	LEU	2.0
1	A	226	PHE	2.0
1	A	397	MET	2.0

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

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

6.3 Carbohydrates [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
2	NAG	A	482	14/15	0.52	0.75	-	86,90,93,95	0
2	BMA	A	483	11/12	0.48	0.69	-	96,97,97,97	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	NAG	A	481	14/15	0.86	0.42	-	61,65,71,80	0

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(\AA^2)	Q<0.9
3	NAG	B	481	14/15	0.79	0.45	-	51,56,58,59	0
4	CA	A	501	1/1	0.98	0.14	-	44,44,44,44	0

6.5 Other polymers [i](#)

There are no such residues in this entry.