



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:10 PM GMT

PDB ID : 4O02
Title : AlphaVBeta3 integrin in complex with monoclonal antibody FAB fragment.
Authors : Mahalingam, B.; van Agthoven, J.; Xiong, J.; Arnaout, M.A.
Deposited on : 2013-12-13
Resolution : 3.60 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

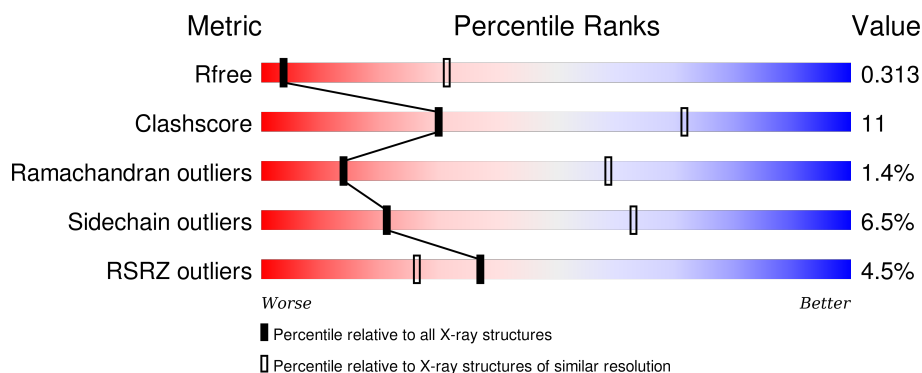
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.60 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	1408 (3.80-3.40)
Clashscore	102246	1010 (3.74-3.46)
Ramachandran outliers	100387	1007 (3.76-3.44)
Sidechain outliers	100360	1007 (3.76-3.44)
RSRZ outliers	91569	1003 (3.78-3.42)

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	962	<div> <div>5%</div> <div>77%</div> <div>17%</div> <div>• 5%</div> </div>
2	B	692	<div> <div>5%</div> <div>75%</div> <div>22%</div> <div>••</div> </div>
3	L	214	<div> <div>%</div> <div>73%</div> <div>24%</div> <div>••</div> </div>
4	H	218	<div> <div>3%</div> <div>72%</div> <div>25%</div> <div>••</div> </div>

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
11	NO3	A	1029	-	X	-	-
11	NO3	A	1030	-	X	-	-

2 Entry composition

There are 11 unique types of molecules in this entry. The entry contains 13526 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 Integrin alpha-V.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	915	Total	C	N	O	S	0	0	0
			5875	3627	1033	1192	23			

- Molecule 2 is a protein called Integrin beta-3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	687	Total	C	N	O	S	0	0	0
			4364	2657	771	871	65			

- Molecule 3 is a protein called 17E6 light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	L	214	Total	C	N	O	S	0	0	0
			1453	887	253	307	6			

- Molecule 4 is a protein called 17E6 heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	H	218	Total	C	N	O	S	0	0	0
			1440	891	247	293	9			

- Molecule 5 is a polymer of unknown type called SUGAR (6-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	6	Total	C	N	O	0	0
			72	40	2	30		

- Molecule 6 is a polymer of unknown type called SUGAR (3-MER).

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

- Molecule 7 is a polymer of unknown type called SUGAR (6-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	A	6	Total	C	N	O	0	0
			72	40	2	30		
7	A	6	Total	C	N	O	0	0
			72	40	2	30		

- Molecule 8 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



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

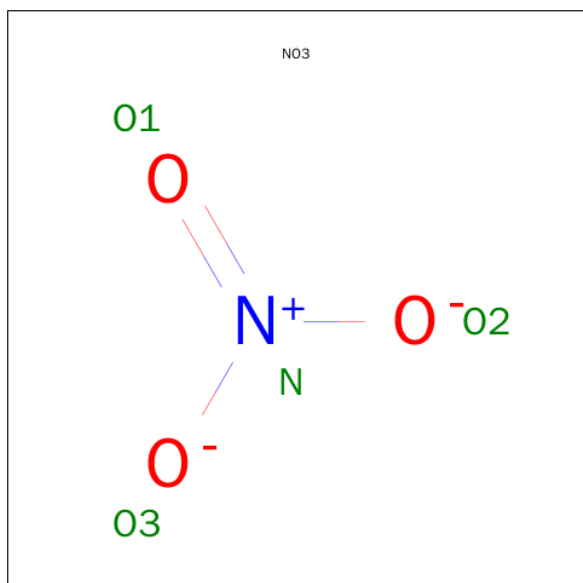
- Molecule 9 is a polymer of unknown type called SUGAR (2-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
9	A	2	Total	C	N	O	0	0
			28	16	2	10		
9	B	2	Total	C	N	O	0	0
			28	16	2	10		
9	B	2	Total	C	N	O	0	0
			28	16	2	10		

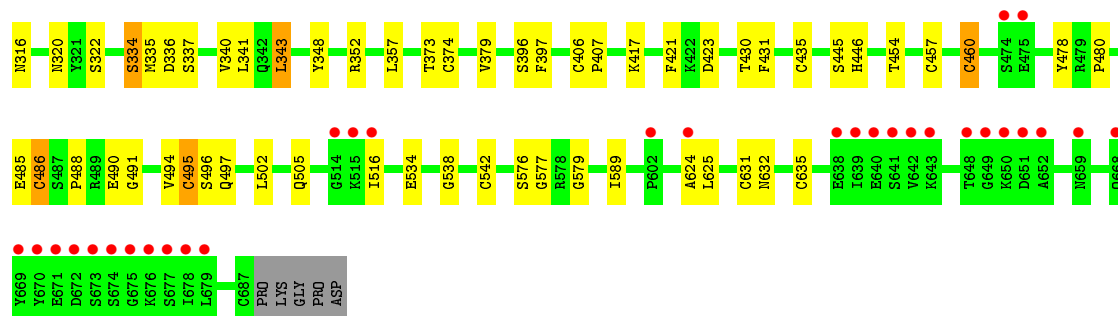
- Molecule 10 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	B	1	Total	Mn	0	0
			1	1		
10	A	4	Total	Mn	0	0
			4	4		

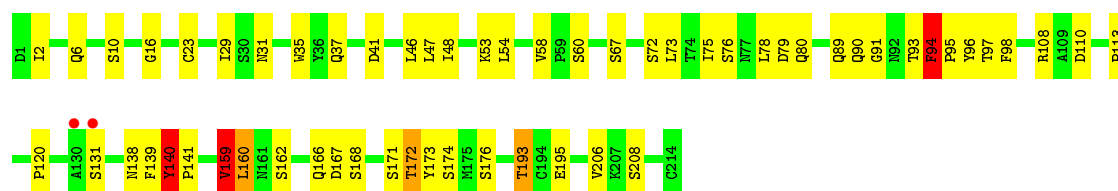
- Molecule 11 is NITRATE ION (three-letter code: NO3) (formula: NO₃).



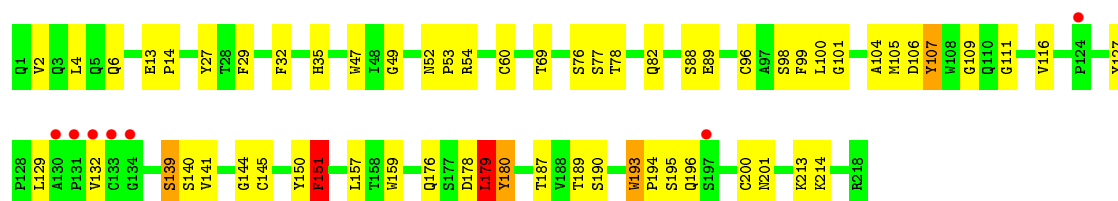
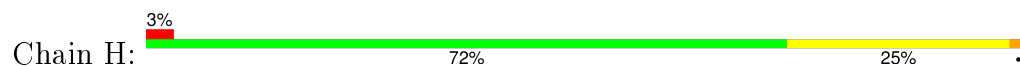
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
11	A	1	Total	N	O	0	0
			4	1	3		
11	A	1	Total	N	O	0	0
			4	1	3		



- Molecule 3: 17E6 light chain



- Molecule 4: 17E6 heavy chain



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	110.46 Å 266.99 Å 102.16 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.86 – 3.60 49.86 – 3.61	Depositor EDS
% Data completeness (in resolution range)	87.6 (49.86-3.60) 87.7 (49.86-3.61)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.13	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.06 (at 3.57 Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE: 1.8.2_1309)	Depositor
R, R_{free}	0.246 , 0.312 0.247 , 0.313	Depositor DCC
R_{free} test set	1561 reflections (4.99%)	DCC
Wilson B-factor (Å ²)	112.3	Xtriage
Anisotropy	0.556	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.24 , 123.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.27$	Xtriage
Outliers	1 of 31303 reflections (0.003%)	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	13526	wwPDB-VP
Average B, all atoms (Å ²)	155.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.42% 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: BMA, NO3, NAG, MN, MAN

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.28	0/5981	0.62	1/8211 (0.0%)
2	B	0.32	1/4438 (0.0%)	0.65	0/6091
3	L	0.29	0/1483	0.62	0/2037
4	H	0.29	0/1475	0.69	1/2027 (0.0%)
All	All	0.30	1/13377 (0.0%)	0.64	2/18366 (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	2
2	B	0	4
3	L	0	3
4	H	0	4
All	All	0	13

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	335	MET	C-O	5.78	1.34	1.23

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	173	GLY	N-CA-C	-5.39	99.62	113.10
4	H	179	LEU	N-CA-C	5.25	125.16	111.00

There are no chirality outliers.

All (13) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	134	THR	Peptide
1	A	768	THR	Peptide
2	B	50	ALA	Peptide
2	B	53	SER	Peptide
2	B	624	ALA	Peptide
2	B	8	ARG	Peptide
4	H	139	SER	Peptide
4	H	151	PHE	Peptide
4	H	178	ASP	Peptide
4	H	179	LEU	Peptide
3	L	140	TYR	Peptide
3	L	168	SER	Peptide
3	L	94	PHE	Peptide

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5875	0	4346	96	0
2	B	4364	0	3235	104	0
3	L	1453	0	1128	33	0
4	H	1440	0	1084	36	0
5	A	72	0	61	0	0
6	A	39	0	34	0	0
7	A	144	0	122	3	0
8	A	14	0	13	0	0
8	B	28	0	26	0	0
9	A	28	0	25	0	0
9	B	56	0	50	1	0
10	A	4	0	0	0	0
10	B	1	0	0	0	0
11	A	8	0	0	0	0
All	All	13526	0	10124	263	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 11.

All (263) 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:253:VAL:HG23	1:A:267:PHE:HB2	1.50	0.94
1:A:560:GLU:HG3	1:A:585:ASN:HB3	1.53	0.89
1:A:121:GLU:HA	2:B:169:PRO:HB3	1.57	0.85
1:A:100:SER:HB3	1:A:105:ILE:HG22	1.59	0.83
2:B:31:LEU:HG	2:B:34:GLY:HA3	1.67	0.77
2:B:579:GLY:HA2	2:B:589:ILE:HG13	1.67	0.76
1:A:232:ASN:HB3	1:A:264:LEU:HD13	1.67	0.76
2:B:488:PRO:HG2	2:B:491:GLY:HA3	1.73	0.70
1:A:195:TYR:HA	1:A:202:ILE:HD13	1.73	0.70
3:L:47:LEU:HA	3:L:58:VAL:HG11	1.74	0.70
4:H:193:TRP:CG	4:H:194:PRO:HD3	2.28	0.68
2:B:94:PRO:HB3	2:B:406:CYS:HB2	1.75	0.67
4:H:140:SER:O	4:H:190:SER:HB2	1.95	0.67
2:B:373:THR:HG22	2:B:379:VAL:HB	1.77	0.66
4:H:213:LYS:NZ	4:H:214:LYS:O	2.25	0.66
1:A:124:PRO:HG2	1:A:154:PHE:HA	1.76	0.66
2:B:126:ASP:OD2	2:B:127:ASP:N	2.28	0.66
3:L:46:LEU:HD11	4:H:104:ALA:HB1	1.77	0.65
1:A:657:VAL:HG22	1:A:659:ASN:H	1.60	0.65
1:A:131:GLN:HG3	1:A:136:THR:HG22	1.78	0.64
2:B:168:SER:HB2	2:B:173:LEU:HB3	1.79	0.64
3:L:162:SER:HB3	3:L:176:SER:HB3	1.80	0.64
4:H:140:SER:HA	4:H:189:THR:HA	1.80	0.64
2:B:127:ASP:HA	2:B:336:ASP:O	1.98	0.64
1:A:346:PRO:HA	1:A:358:ILE:HG12	1.80	0.63
1:A:88:PHE:HB2	1:A:113:HIS:HB2	1.80	0.63
3:L:29:ILE:HG21	3:L:90:GLN:HG3	1.81	0.63
1:A:14:PRO:HG2	1:A:17:SER:HB3	1.80	0.62
2:B:83:VAL:O	2:B:86:GLN:NE2	2.33	0.62
2:B:30:ALA:O	2:B:48:ASN:ND2	2.33	0.61
1:A:290:TYR:HB3	1:A:320:GLN:HB3	1.82	0.61
4:H:2:VAL:HG21	4:H:107:TYR:CD2	2.36	0.60
3:L:138:ASN:HA	3:L:172:THR:HB	1.82	0.60
3:L:110:ASP:HA	3:L:140:TYR:HB3	1.84	0.60
2:B:534:GLU:O	2:B:538:GLY:N	2.35	0.59
1:A:791:SER:OG	1:A:930:GLU:O	2.20	0.59
4:H:150:TYR:HB3	4:H:180:TYR:HB2	1.83	0.59
4:H:35:HIS:CD2	4:H:99:PHE:HB2	2.39	0.58
1:A:335:GLU:OE2	1:A:366:GLY:N	2.30	0.58
1:A:175:GLY:HA2	1:A:179:TRP:CD1	2.39	0.57
1:A:472:CYS:HA	1:A:541:TYR:HA	1.86	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:H:69:THR:OG1	4:H:82:GLN:HB2	2.05	0.57
2:B:158:ASP:OD1	2:B:289:TYR:OH	2.22	0.57
3:L:96:TYR:HB2	4:H:47:TRP:CG	2.40	0.57
2:B:486:CYS:HB3	2:B:494:VAL:O	2.03	0.57
1:A:395:TRP:CE2	1:A:431:ARG:HG2	2.40	0.57
2:B:64:LEU:HD11	2:B:89:ALA:HB2	1.86	0.57
4:H:35:HIS:HD2	4:H:99:PHE:HB2	1.69	0.57
4:H:6:GLN:HG3	4:H:111:GLY:H	1.70	0.56
1:A:350:LEU:HA	1:A:420:PRO:HD2	1.87	0.56
2:B:245:LEU:HD11	2:B:348:TYR:HD1	1.70	0.56
2:B:178:TYR:H	2:B:178:TYR:HD1	1.53	0.56
4:H:141:VAL:O	4:H:187:THR:HA	2.06	0.56
1:A:924:ALA:O	1:A:945:THR:N	2.32	0.55
1:A:925:SER:HA	1:A:944:SER:HA	1.88	0.55
2:B:27:SER:HA	2:B:457:CYS:HB3	1.88	0.55
1:A:570:ASP:HB3	1:A:573:GLY:HA2	1.87	0.55
1:A:100:SER:CB	1:A:105:ILE:HG22	2.35	0.55
3:L:166:GLN:HB2	3:L:173:TYR:CZ	2.41	0.55
2:B:249:THR:HG22	2:B:309:ALA:HB3	1.87	0.55
2:B:49:CYS:HA	2:B:51:PRO:HD2	1.88	0.54
1:A:399:SER:OG	1:A:400:MET:N	2.40	0.54
2:B:194:LEU:HD22	2:B:203:PHE:HA	1.89	0.54
2:B:221:GLY:HA2	2:B:289:TYR:HE2	1.72	0.54
3:L:16:GLY:H	3:L:78:LEU:HB3	1.73	0.54
1:A:141:PRO:HB2	1:A:184:ILE:HD13	1.89	0.54
7:A:1019:NAG:H61	7:A:1020:NAG:H2	1.90	0.54
2:B:106:GLN:HG2	2:B:352:ARG:HG3	1.90	0.54
2:B:340:VAL:HA	2:B:343:LEU:HD23	1.88	0.54
1:A:633:ASN:ND2	1:A:636:GLU:O	2.40	0.54
2:B:334:SER:OG	2:B:336:ASP:OD2	2.25	0.53
2:B:445:SER:OG	2:B:446:HIS:N	2.41	0.53
1:A:256:TYR:CD1	1:A:263:SER:HB3	2.42	0.53
1:A:255:ILE:HD12	1:A:265:TYR:HB2	1.90	0.53
2:B:454:THR:O	2:B:460:CYS:HB3	2.07	0.53
1:A:2:ASN:ND2	1:A:438:ARG:O	2.40	0.53
1:A:17:SER:HB2	1:A:43:ALA:HB2	1.91	0.53
3:L:167:ASP:O	3:L:171:SER:HA	2.09	0.53
1:A:301:MET:HA	1:A:311:GLU:HA	1.91	0.52
2:B:495:CYS:O	2:B:497:GLN:N	2.43	0.52
1:A:38:VAL:HG23	1:A:57:LEU:HB2	1.91	0.52
2:B:281:TYR:CE1	2:B:283:ALA:HB3	2.45	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:197:THR:HG22	2:B:199:GLN:H	1.74	0.52
1:A:253:VAL:HG22	1:A:276:PHE:HE2	1.75	0.51
4:H:13:GLU:HG3	4:H:14:PRO:HD2	1.91	0.51
1:A:338:ALA:HB1	1:A:362:ALA:HB1	1.93	0.51
2:B:122:TYR:N	2:B:213:SER:O	2.44	0.51
3:L:93:THR:OG1	3:L:94:PHE:N	2.44	0.51
2:B:169:PRO:HG2	2:B:170:PRO:HD3	1.92	0.51
4:H:151:PHE:HB2	4:H:179:LEU:CB	2.41	0.51
2:B:190:TYR:HB3	2:B:281:TYR:HB2	1.92	0.51
2:B:113:ASP:OD1	2:B:239:ARG:NE	2.42	0.50
1:A:114:TRP:HZ3	1:A:147:ILE:HG21	1.77	0.50
2:B:70:SER:HB2	2:B:80:VAL:O	2.11	0.50
1:A:86:LEU:HD11	1:A:200:TYR:CE1	2.47	0.50
1:A:76:GLY:O	1:A:89:LYS:NZ	2.44	0.50
3:L:48:ILE:HA	3:L:53:LYS:O	2.12	0.50
2:B:407:PRO:HD2	2:B:431:PHE:CD1	2.47	0.50
1:A:283:THR:CG2	1:A:358:ILE:HD11	2.42	0.49
1:A:339:ARG:HH12	2:B:268:PRO:HB3	1.77	0.49
1:A:159:PHE:HB3	1:A:174:PRO:HD3	1.92	0.49
3:L:54:LEU:HD11	3:L:60:SER:HA	1.94	0.49
1:A:640:GLU:HA	1:A:685:ASN:HA	1.95	0.49
1:A:314:GLN:OE1	1:A:332:ASN:ND2	2.46	0.49
1:A:299:LEU:HD11	2:B:258:LEU:HG	1.94	0.49
1:A:272:MET:HE1	2:B:320:ASN:HD22	1.78	0.49
1:A:487:LEU:O	1:A:489:ARG:N	2.40	0.49
2:B:283:ALA:HB1	2:B:287:MET:HB2	1.94	0.48
2:B:165:MET:HG2	2:B:166:TYR:H	1.77	0.48
4:H:29:PHE:CD2	4:H:77:SER:HA	2.48	0.48
3:L:159:VAL:O	3:L:160:LEU:HG	2.13	0.48
3:L:37:GLN:HB2	3:L:47:LEU:HD11	1.94	0.48
1:A:765:ASN:N	1:A:766:PRO:HD3	2.29	0.48
2:B:88:ILE:O	2:B:88:ILE:HG13	2.12	0.48
3:L:31:ASN:OD1	3:L:67:SER:HA	2.13	0.48
4:H:76:SER:OG	4:H:78:THR:OG1	2.30	0.48
1:A:404:PHE:HA	1:A:426:ALA:HB2	1.95	0.47
4:H:88:SER:HA	4:H:116:VAL:HB	1.96	0.47
2:B:64:LEU:HB2	2:B:87:ARG:HG3	1.95	0.47
4:H:144:GLY:HA2	4:H:159:TRP:CH2	2.49	0.47
2:B:134:LEU:HD11	2:B:341:LEU:HB2	1.95	0.47
2:B:143:ARG:HA	2:B:143:ARG:HD3	1.40	0.47
2:B:495:CYS:C	2:B:497:GLN:H	2.18	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:339:ARG:NH1	2:B:268:PRO:HB3	2.30	0.47
1:A:566:ARG:HA	1:A:566:ARG:HD3	1.73	0.47
4:H:193:TRP:O	4:H:195:SER:N	2.47	0.47
2:B:49:CYS:C	2:B:51:PRO:HD2	2.35	0.47
1:A:395:TRP:CZ2	1:A:431:ARG:HG2	2.50	0.47
2:B:154:GLY:HA2	2:B:193:VAL:HG23	1.97	0.47
1:A:186:ASP:OD1	1:A:207:GLN:HB2	2.15	0.47
2:B:82:GLN:N	2:B:423:ASP:OD2	2.48	0.47
2:B:110:TYR:HB3	2:B:352:ARG:HH22	1.79	0.47
1:A:408:MET:HA	1:A:423:ILE:O	2.13	0.47
1:A:230:ASP:HA	1:A:238:ASP:OD2	2.15	0.47
1:A:18:TYR:CD1	1:A:427:PHE:HD1	2.32	0.47
3:L:120:PRO:HG3	3:L:131:SER:O	2.15	0.46
1:A:577:ILE:HG22	1:A:578:LEU:O	2.14	0.46
2:B:139:ALA:O	2:B:143:ARG:HG2	2.16	0.46
3:L:2:ILE:O	3:L:97:THR:HG21	2.15	0.46
3:L:75:ILE:HG22	3:L:76:SER:O	2.16	0.46
2:B:200:VAL:O	2:B:203:PHE:HB3	2.15	0.46
1:A:181:GLY:HA3	1:A:222:LEU:HB3	1.96	0.46
2:B:153:PHE:HE2	2:B:210:GLN:HG3	1.79	0.46
1:A:20:GLY:O	1:A:407:SER:OG	2.31	0.46
1:A:657:VAL:HG13	1:A:658:ARG:H	1.79	0.46
2:B:158:ASP:HA	2:B:221:GLY:N	2.30	0.46
1:A:517:ARG:CG	1:A:517:ARG:HH11	2.29	0.46
2:B:139:ALA:HB2	2:B:200:VAL:HG21	1.96	0.46
4:H:100:LEU:HG	4:H:101:GLY:H	1.79	0.46
4:H:4:LEU:HB2	4:H:109:GLY:HA2	1.97	0.46
1:A:419:TYR:CE1	1:A:439:PRO:HA	2.50	0.46
3:L:108:ARG:HD2	3:L:171:SER:O	2.16	0.46
2:B:66:ASP:OD1	2:B:66:ASP:N	2.49	0.46
3:L:91:GLY:HA2	3:L:96:TYR:CE1	2.50	0.46
1:A:106:LEU:HD11	1:A:128:CYS:HB3	1.98	0.46
1:A:146:ASP:OD2	4:H:54:ARG:NE	2.41	0.46
1:A:492:ASN:ND2	1:A:564:ASP:OD2	2.49	0.45
4:H:127:TYR:O	4:H:129:LEU:N	2.49	0.45
1:A:216:ILE:HG13	1:A:217:PHE:N	2.31	0.45
1:A:257:ASP:HB2	1:A:264:LEU:HD21	1.98	0.45
4:H:100:LEU:HB2	4:H:106:ASP:CG	2.37	0.45
3:L:110:ASP:OD2	3:L:141:PRO:HD3	2.17	0.45
3:L:35:TRP:CE2	3:L:73:LEU:HB2	2.52	0.45
1:A:170:LEU:HD22	1:A:239:PHE:CD1	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:278:PHE:CD2	1:A:298:PRO:HG3	2.51	0.45
3:L:89:GLN:HB2	3:L:98:PHE:CD2	2.52	0.45
1:A:375:ILE:HD12	1:A:389:GLN:HB3	1.98	0.45
2:B:21:PRO:HG2	2:B:96:ASP:HB2	1.98	0.45
2:B:204:ASN:HA	2:B:207:VAL:HG12	1.99	0.45
2:B:221:GLY:HA2	2:B:289:TYR:CE2	2.51	0.44
4:H:49:GLY:HA3	4:H:60:CYS:HA	1.99	0.44
2:B:245:LEU:HD11	2:B:348:TYR:CD1	2.51	0.44
2:B:20:SER:HB3	2:B:22:MET:HG2	1.98	0.44
1:A:663:LEU:HA	1:A:694:LEU:O	2.16	0.44
2:B:169:PRO:CG	2:B:170:PRO:HD3	2.48	0.44
2:B:120:LEU:HD12	2:B:120:LEU:HA	1.68	0.44
1:A:605:LYS:H	1:A:635:GLY:HA3	1.82	0.44
2:B:30:ALA:HB1	2:B:49:CYS:HB3	1.99	0.44
2:B:176:PRO:CG	2:B:186:PRO:HA	2.47	0.44
1:A:719:SER:OG	1:A:720:SER:N	2.51	0.44
7:A:1013:MAN:H3	7:A:1014:MAN:H2	1.52	0.44
1:A:642:GLU:HG2	1:A:682:ASP:HA	1.99	0.44
2:B:27:SER:HB2	2:B:53:SER:OG	2.18	0.44
3:L:48:ILE:HG23	3:L:53:LYS:H	1.83	0.44
4:H:52:ASN:HD22	4:H:53:PRO:HD2	1.83	0.44
1:A:81:ALA:HB3	1:A:84:ASP:HB3	1.99	0.44
2:B:297:GLU:OE1	2:B:301:GLN:NE2	2.51	0.44
2:B:158:ASP:HA	2:B:221:GLY:H	1.83	0.44
4:H:98:SER:O	4:H:106:ASP:N	2.50	0.44
3:L:79:ASP:OD1	3:L:80:GLN:N	2.39	0.44
1:A:8:PRO:HB3	1:A:435:TYR:HD2	1.83	0.44
1:A:70:ILE:HB	1:A:72:PHE:CE1	2.53	0.44
3:L:46:LEU:HD12	4:H:105:MET:O	2.17	0.43
2:B:67:ARG:NH1	2:B:86:GLN:HG2	2.33	0.43
3:L:195:GLU:HG3	3:L:206:VAL:HG22	2.00	0.43
1:A:283:THR:HG21	1:A:358:ILE:HD11	2.00	0.43
2:B:220:GLU:H	2:B:254:THR:HA	1.84	0.43
4:H:159:TRP:CZ3	4:H:200:CYS:HB2	2.53	0.43
2:B:30:ALA:CB	2:B:49:CYS:HB3	2.49	0.43
2:B:136:THR:HA	2:B:200:VAL:HG11	2.00	0.43
2:B:136:THR:O	2:B:140:THR:HG23	2.18	0.43
1:A:554:PRO:HB2	1:A:589:GLN:HB3	1.99	0.43
2:B:283:ALA:O	2:B:287:MET:N	2.48	0.43
1:A:239:PHE:O	1:A:255:ILE:HA	2.18	0.43
3:L:113:PRO:HB3	3:L:139:PHE:CD2	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:357:LEU:HD11	2:B:397:PHE:CD2	2.53	0.43
2:B:28:ASP:N	2:B:457:CYS:O	2.52	0.43
1:A:763:LYS:O	1:A:766:PRO:HG3	2.19	0.43
2:B:265:ILE:HG23	2:B:286:THR:O	2.18	0.43
2:B:91:ARG:HA	2:B:430:THR:O	2.19	0.43
2:B:168:SER:HB2	2:B:173:LEU:CB	2.48	0.42
2:B:38:CYS:HB3	2:B:39:ASP:H	1.53	0.42
1:A:347:LEU:HD11	1:A:359:ALA:HB2	2.00	0.42
1:A:134:THR:OG1	1:A:135:LYS:N	2.53	0.42
1:A:829:ASN:HA	1:A:830:PRO:HD3	1.69	0.42
1:A:152:GLN:HB3	1:A:155:CYS:HB2	2.02	0.42
2:B:138:LEU:HA	2:B:341:LEU:HD13	2.01	0.42
2:B:336:ASP:O	2:B:337:SER:CB	2.68	0.42
1:A:525:MET:HE3	1:A:536:GLU:HG2	2.01	0.42
1:A:596:CYS:HB3	1:A:602:CYS:HB3	1.92	0.42
2:B:219:PRO:HG2	2:B:255:HIS:CE1	2.55	0.42
1:A:339:ARG:HB3	1:A:342:SER:OG	2.20	0.42
4:H:100:LEU:HG	4:H:101:GLY:N	2.34	0.42
2:B:316:ASN:HB3	9:B:702:NAG:HN2	1.85	0.42
2:B:170:PRO:HG2	2:B:172:ALA:HB3	2.01	0.42
4:H:47:TRP:CH2	4:H:49:GLY:HA2	2.55	0.42
2:B:115:TYR:OH	2:B:192:HIS:ND1	2.49	0.42
1:A:419:TYR:HA	1:A:420:PRO:HD3	1.77	0.41
2:B:152:GLY:HA2	2:B:194:LEU:HD23	2.02	0.41
3:L:139:PHE:CE2	3:L:174:SER:HA	2.54	0.41
3:L:96:TYR:HD1	3:L:96:TYR:HA	1.75	0.41
2:B:65:GLU:HG3	2:B:87:ARG:HG2	2.02	0.41
2:B:49:CYS:HB2	2:B:50:ALA:H	1.74	0.41
4:H:144:GLY:HA2	4:H:159:TRP:HH2	1.86	0.41
1:A:629:VAL:O	1:A:694:LEU:HA	2.20	0.41
1:A:248:ARG:O	1:A:249:THR:OG1	2.32	0.41
3:L:54:LEU:HD22	3:L:58:VAL:HG23	2.03	0.41
2:B:165:MET:HG3	2:B:187:MET:HE3	2.01	0.41
2:B:124:MET:O	2:B:127:ASP:HB2	2.21	0.41
1:A:8:PRO:HB3	1:A:435:TYR:CD2	2.56	0.41
1:A:748:SER:HA	1:A:778:HIS:HA	2.02	0.41
2:B:478:TYR:C	2:B:480:PRO:HD3	2.41	0.41
1:A:291:ALA:O	1:A:320:GLN:HB2	2.21	0.41
2:B:165:MET:HG2	2:B:166:TYR:N	2.35	0.41
2:B:269:ASN:HA	2:B:287:MET:SD	2.60	0.41
3:L:193:THR:HG22	3:L:208:SER:HA	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:H:27:TYR:HE2	4:H:32:PHE:HB2	1.86	0.41
1:A:794:MET:O	1:A:926:PHE:HA	2.21	0.41
2:B:76:ASP:O	2:B:78:SER:N	2.50	0.41
4:H:157:LEU:HA	4:H:201:ASN:O	2.21	0.41
2:B:502:LEU:O	2:B:505:GLN:O	2.39	0.41
2:B:10:VAL:C	2:B:12:SER:H	2.24	0.41
1:A:768:THR:OG1	1:A:769:GLU:N	2.54	0.40
2:B:407:PRO:HD2	2:B:431:PHE:CE1	2.56	0.40
1:A:298:PRO:O	1:A:339:ARG:HA	2.22	0.40
2:B:49:CYS:CA	2:B:51:PRO:HD2	2.50	0.40
2:B:249:THR:HA	2:B:309:ALA:O	2.21	0.40
7:A:1010:NAG:H61	7:A:1011:NAG:H82	2.02	0.40
2:B:485:GLU:CB	2:B:490:GLU:H	2.34	0.40
4:H:139:SER:HA	4:H:190:SER:HB3	2.02	0.40
1:A:764:GLU:C	1:A:766:PRO:HD3	2.42	0.40
2:B:4:ILE:C	2:B:6:THR:H	2.24	0.40
2:B:238:TRP:CE3	2:B:244:HIS:HD2	2.39	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	911/962 (95%)	815 (90%)	88 (10%)	8 (1%)	21	67
2	B	685/692 (99%)	595 (87%)	79 (12%)	11 (2%)	12	56
3	L	212/214 (99%)	196 (92%)	12 (6%)	4 (2%)	10	53
4	H	216/218 (99%)	190 (88%)	20 (9%)	6 (3%)	6	45
All	All	2024/2086 (97%)	1796 (89%)	199 (10%)	29 (1%)	14	59

All (29) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	135	LYS
1	A	619	ILE
1	A	710	SER
4	H	180	TYR
1	A	453	ILE
1	A	738	VAL
1	A	876	VAL
2	B	516	ILE
2	B	632	ASN
3	L	159	VAL
2	B	36	PRO
2	B	80	VAL
2	B	496	SER
3	L	94	PHE
3	L	140	TYR
4	H	151	PHE
2	B	10	VAL
2	B	76	ASP
2	B	625	LEU
4	H	107	TYR
4	H	193	TRP
1	A	555	ILE
2	B	577	GLY
4	H	196	GLN
2	B	77	SER
1	A	766	PRO
3	L	95	PRO
2	B	162	SER
4	H	132	VAL

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	400/816 (49%)	378 (94%)	22 (6%)	27	68
2	B	334/614 (54%)	308 (92%)	26 (8%)	16	55
3	L	125/192 (65%)	115 (92%)	10 (8%)	15	54

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
4	H	107/183 (58%)	102 (95%)	5 (5%)	32	72
All	All	966/1805 (54%)	903 (94%)	63 (6%)	21	63

All (63) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	29	SER
1	A	38	VAL
1	A	103	ASP
1	A	142	CYS
1	A	147	ILE
1	A	152	GLN
1	A	188	VAL
1	A	207	GLN
1	A	237	ASP
1	A	253	VAL
1	A	255	ILE
1	A	264	LEU
1	A	275	TYR
1	A	303	ARG
1	A	351	ASP
1	A	423	ILE
1	A	517	ARG
1	A	558	PHE
1	A	560	GLU
1	A	565	TYR
1	A	681	CYS
1	A	943	ASN
2	B	38	CYS
2	B	88	ILE
2	B	143	ARG
2	B	162	SER
2	B	177	CYS
2	B	178	TYR
2	B	232	CYS
2	B	284	SER
2	B	285	THR
2	B	300	SER
2	B	308	PHE
2	B	322	SER
2	B	334	SER
2	B	343	LEU

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Mol	Chain	Res	Type
2	B	374	CYS
2	B	396	SER
2	B	417	LYS
2	B	421	PHE
2	B	435	CYS
2	B	460	CYS
2	B	486	CYS
2	B	495	CYS
2	B	542	CYS
2	B	576	SER
2	B	631	CYS
2	B	635	CYS
3	L	6	GLN
3	L	10	SER
3	L	23	CYS
3	L	41	ASP
3	L	72	SER
3	L	94	PHE
3	L	159	VAL
3	L	160	LEU
3	L	172	THR
3	L	193	THR
4	H	89	GLU
4	H	96	CYS
4	H	145	CYS
4	H	151	PHE
4	H	176	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	214	GLN
1	A	310	GLN
1	A	314	GLN
1	A	332	ASN
2	B	313	ASN
3	L	138	ASN
4	H	52	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

27 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$
5	NAG	A	1001	1,5	14,14,15	0.30	0	15,19,21	0.43	0
5	NAG	A	1002	5	14,14,15	0.41	0	15,19,21	0.75	1 (6%)
5	BMA	A	1003	5	11,11,12	1.02	1 (9%)	14,15,17	1.26	1 (7%)
5	MAN	A	1004	5	11,11,12	0.97	1 (9%)	14,15,17	0.94	1 (7%)
5	BMA	A	1005	5	11,11,12	0.59	0	14,15,17	1.06	1 (7%)
5	MAN	A	1006	5	11,11,12	0.66	0	14,15,17	1.17	2 (14%)
6	NAG	A	1007	1,6	14,14,15	0.40	0	15,19,21	0.32	0
6	NAG	A	1008	6	14,14,15	0.41	0	15,19,21	0.38	0
6	BMA	A	1009	6	11,11,12	0.72	0	14,15,17	0.85	0
7	NAG	A	1010	1,7	14,14,15	0.51	0	15,19,21	0.92	1 (6%)
7	NAG	A	1011	7	14,14,15	0.42	0	15,19,21	0.32	0
7	BMA	A	1012	7	11,11,12	0.64	0	14,15,17	1.03	1 (7%)
7	MAN	A	1013	7	11,11,12	0.85	0	14,15,17	1.64	2 (14%)
7	MAN	A	1014	7	11,11,12	0.81	0	14,15,17	1.49	2 (14%)
7	MAN	A	1015	7	11,11,12	0.52	0	14,15,17	1.14	2 (14%)
9	NAG	A	1017	1,9	14,14,15	0.40	0	15,19,21	0.39	0
9	NAG	A	1018	9	14,14,15	0.31	0	15,19,21	0.17	0
7	NAG	A	1019	1,7	14,14,15	0.23	0	15,19,21	0.47	0
7	NAG	A	1020	7	14,14,15	0.62	1 (7%)	15,19,21	0.85	1 (6%)
7	BMA	A	1021	7	11,11,12	1.17	1 (9%)	14,15,17	1.00	1 (7%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
7	MAN	A	1022	7	11,11,12	1.21	1 (9%)	14,15,17	1.87	2 (14%)
7	MAN	A	1023	7	11,11,12	1.34	2 (18%)	14,15,17	1.86	4 (28%)
7	MAN	A	1024	7	11,11,12	0.78	1 (9%)	14,15,17	1.64	4 (28%)
9	NAG	B	702	9,2	14,14,15	0.74	1 (7%)	15,19,21	1.11	1 (6%)
9	NAG	B	703	9	14,14,15	0.37	0	15,19,21	0.98	1 (6%)
9	NAG	B	704	9,2	14,14,15	1.27	1 (7%)	15,19,21	1.76	2 (13%)
9	NAG	B	705	9	14,14,15	0.41	0	15,19,21	0.60	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	A	1001	1,5	-	0/6/23/26	0/1/1/1
5	NAG	A	1002	5	-	0/6/23/26	0/1/1/1
5	BMA	A	1003	5	-	0/2/19/22	0/1/1/1
5	MAN	A	1004	5	-	0/2/19/22	0/1/1/1
5	BMA	A	1005	5	-	0/2/19/22	0/1/1/1
5	MAN	A	1006	5	-	0/2/19/22	0/1/1/1
6	NAG	A	1007	1,6	-	0/6/23/26	0/1/1/1
6	NAG	A	1008	6	-	0/6/23/26	0/1/1/1
6	BMA	A	1009	6	-	0/2/19/22	0/1/1/1
7	NAG	A	1010	1,7	-	0/6/23/26	0/1/1/1
7	NAG	A	1011	7	-	0/6/23/26	0/1/1/1
7	BMA	A	1012	7	-	0/2/19/22	0/1/1/1
7	MAN	A	1013	7	-	0/2/19/22	1/1/1/1
7	MAN	A	1014	7	-	0/2/19/22	1/1/1/1
7	MAN	A	1015	7	-	0/2/19/22	0/1/1/1
9	NAG	A	1017	1,9	-	0/6/23/26	0/1/1/1
9	NAG	A	1018	9	-	0/6/23/26	0/1/1/1
7	NAG	A	1019	1,7	-	0/6/23/26	0/1/1/1
7	NAG	A	1020	7	-	0/6/23/26	0/1/1/1
7	BMA	A	1021	7	-	0/2/19/22	0/1/1/1
7	MAN	A	1022	7	-	0/2/19/22	1/1/1/1
7	MAN	A	1023	7	-	0/2/19/22	0/1/1/1
7	MAN	A	1024	7	-	0/2/19/22	0/1/1/1
9	NAG	B	702	9,2	-	0/6/23/26	0/1/1/1
9	NAG	B	703	9	-	0/6/23/26	0/1/1/1
9	NAG	B	704	9,2	-	0/6/23/26	0/1/1/1
9	NAG	B	705	9	-	0/6/23/26	0/1/1/1

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	1004	MAN	O5-C1	-2.53	1.39	1.43
5	A	1003	BMA	C2-C3	-2.06	1.49	1.52
7	A	1020	NAG	O5-C1	2.07	1.47	1.43
7	A	1024	MAN	C1-C2	2.15	1.57	1.52
7	A	1022	MAN	O3-C3	2.44	1.48	1.43
7	A	1021	BMA	C2-C3	2.60	1.56	1.52
9	B	702	NAG	O5-C1	2.61	1.48	1.43
7	A	1023	MAN	C2-C3	2.72	1.56	1.52
7	A	1023	MAN	C1-C2	2.75	1.58	1.52
9	B	704	NAG	O5-C1	4.64	1.51	1.43

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	1003	BMA	C1-C2-C3	-2.87	106.14	109.54
7	A	1013	MAN	O2-C2-C3	-2.51	105.08	110.12
7	A	1024	MAN	O2-C2-C3	-2.29	105.52	110.12
7	A	1023	MAN	O2-C2-C3	-2.28	105.53	110.12
7	A	1015	MAN	O2-C2-C3	-2.22	105.66	110.12
5	A	1004	MAN	O2-C2-C3	-2.21	105.67	110.12
5	A	1006	MAN	O2-C2-C3	-2.20	105.69	110.12
7	A	1022	MAN	O2-C2-C3	-2.15	105.80	110.12
7	A	1014	MAN	O2-C2-C3	-2.11	105.88	110.12
9	B	704	NAG	C4-C3-C2	-2.01	108.10	111.23
5	A	1002	NAG	C1-O5-C5	2.06	114.86	112.25
7	A	1024	MAN	C1-C2-C3	2.08	112.00	109.54
7	A	1021	BMA	C1-C2-C3	2.12	112.05	109.54
9	B	705	NAG	C1-O5-C5	2.14	114.97	112.25
5	A	1005	BMA	C1-O5-C5	2.24	115.09	112.25
7	A	1010	NAG	C1-O5-C5	2.42	115.32	112.25
7	A	1024	MAN	O5-C1-C2	2.47	114.87	110.86
7	A	1015	MAN	C1-O5-C5	2.89	115.92	112.25
7	A	1012	BMA	C1-O5-C5	3.12	116.21	112.25
7	A	1023	MAN	O5-C1-C2	3.17	116.00	110.86
7	A	1020	NAG	C1-O5-C5	3.18	116.28	112.25
5	A	1006	MAN	C1-O5-C5	3.23	116.34	112.25
9	B	703	NAG	C1-O5-C5	3.42	116.59	112.25
7	A	1023	MAN	C1-C2-C3	3.45	113.63	109.54
7	A	1023	MAN	C1-O5-C5	3.46	116.64	112.25
9	B	702	NAG	C1-O5-C5	3.92	117.22	112.25
7	A	1024	MAN	C1-O5-C5	4.02	117.35	112.25
7	A	1014	MAN	C1-O5-C5	4.46	117.91	112.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	A	1013	MAN	C1-O5-C5	4.86	118.41	112.25
7	A	1022	MAN	C1-O5-C5	5.40	119.10	112.25
9	B	704	NAG	C1-O5-C5	6.22	120.14	112.25

There are no chirality outliers.

There are no torsion outliers.

All (3) ring outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	A	1022	MAN	C1-C2-C3-C4-C5-O5
7	A	1013	MAN	C1-C2-C3-C4-C5-O5
7	A	1014	MAN	C1-C2-C3-C4-C5-O5

7 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	A	1010	NAG	1	0
7	A	1011	NAG	1	0
7	A	1013	MAN	1	0
7	A	1014	MAN	1	0
7	A	1019	NAG	1	0
7	A	1020	NAG	1	0
9	B	702	NAG	1	0

5.6 Ligand geometry [i](#)

Of 10 ligands modelled in this entry, 5 are monoatomic - leaving 5 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
8	NAG	A	1016	1	14,14,15	0.40	0	15,19,21	0.34	0
11	NO3	A	1029	-	3,3,3	3.29	3 (100%)	3,3,3	0.12	0
11	NO3	A	1030	-	3,3,3	3.28	3 (100%)	3,3,3	0.20	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
8	NAG	B	701	2	14,14,15	0.41	0	15,19,21	0.78	1 (6%)
8	NAG	B	706	2	14,14,15	0.31	0	15,19,21	0.50	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
8	NAG	A	1016	1	-	0/6/23/26	0/1/1/1
11	NO3	A	1029	-	-	0/0/0/0	0/0/0/0
11	NO3	A	1030	-	-	0/0/0/0	0/0/0/0
8	NAG	B	701	2	-	0/6/23/26	0/1/1/1
8	NAG	B	706	2	-	0/6/23/26	0/1/1/1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
11	A	1029	NO3	O3-N	2.84	1.40	1.25
11	A	1029	NO3	O2-N	2.89	1.40	1.25
11	A	1030	NO3	O3-N	2.90	1.40	1.25
11	A	1030	NO3	O2-N	3.03	1.41	1.25
11	A	1030	NO3	O1-N	3.84	1.40	1.24
11	A	1029	NO3	O1-N	4.01	1.40	1.24

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	B	701	NAG	C1-O5-C5	2.84	115.85	112.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

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	915/962 (95%)	0.01	47 (5%) 32 22	52, 165, 260, 296	0
2	B	687/692 (99%)	-0.07	35 (5%) 32 22	55, 145, 303, 428	0
3	L	214/214 (100%)	-0.33	2 (0%) 85 75	71, 127, 234, 297	0
4	H	218/218 (100%)	-0.05	7 (3%) 51 37	81, 154, 242, 329	0
All	All	2034/2086 (97%)	-0.06	91 (4%) 37 26	52, 149, 274, 428	0

All (91) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	515	LYS	10.3
4	H	133	CYS	9.6
2	B	678	ILE	9.4
2	B	670	TYR	9.1
2	B	650	LYS	9.1
1	A	683	LEU	8.9
1	A	454	LEU	7.6
2	B	672	ASP	7.4
2	B	651	ASP	7.4
4	H	131	PRO	7.4
1	A	822	CYS	6.5
1	A	452	SER	6.5
1	A	455	ASN	6.3
2	B	673	SER	6.3
2	B	668	GLN	6.1
2	B	638	GLU	5.9
4	H	130	ALA	5.9
2	B	669	TYR	5.7
2	B	659	ASN	5.4
4	H	132	VAL	5.4
2	B	639	ILE	5.2

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Mol	Chain	Res	Type	RSRZ
2	B	516	ILE	5.0
1	A	824	SER	4.9
2	B	677	SER	4.8
1	A	739	LEU	4.6
1	A	453	ILE	4.6
1	A	936	LEU	4.5
1	A	511	ALA	4.5
2	B	652	ALA	4.5
1	A	684	GLY	4.5
3	L	131	SER	4.4
2	B	675	GLY	4.3
1	A	917	SER	4.2
1	A	937	PRO	4.1
1	A	915	ASN	4.1
2	B	649	GLY	4.1
2	B	679	LEU	3.9
2	B	641	SER	3.8
2	B	36	PRO	3.7
2	B	648	THR	3.7
3	L	130	ALA	3.7
1	A	682	ASP	3.6
2	B	674	SER	3.6
1	A	667	SER	3.6
1	A	446	GLY	3.5
4	H	124	PRO	3.4
2	B	474	SER	3.4
1	A	451	PRO	3.3
1	A	483	GLY	3.3
4	H	197	SER	3.3
1	A	935	ASN	3.3
1	A	939	GLU	3.3
1	A	945	THR	3.2
2	B	602	PRO	3.2
1	A	543	ARG	3.1
1	A	681	CYS	3.1
1	A	916	HIS	3.1
1	A	445	ALA	3.0
1	A	507	ALA	3.0
1	A	774	PRO	2.9
1	A	601	VAL	2.8
2	B	642	VAL	2.8
4	H	134	GLY	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	737	ALA	2.8
1	A	884	CYS	2.7
1	A	447	LEU	2.7
2	B	643	LYS	2.7
2	B	640	GLU	2.6
1	A	592	ILE	2.6
2	B	475	GLU	2.6
1	A	590	ALA	2.6
2	B	676	LYS	2.5
1	A	900	LYS	2.5
1	A	662	ALA	2.5
1	A	740	ALA	2.5
1	A	465	GLY	2.4
2	B	39	ASP	2.4
1	A	591	HIS	2.4
1	A	738	VAL	2.3
2	B	514	GLY	2.3
1	A	493	PHE	2.3
1	A	731	SER	2.3
2	B	671	GLU	2.2
1	A	918	TYR	2.2
1	A	788	SER	2.2
1	A	669	ALA	2.2
2	B	215	ASN	2.1
2	B	624	ALA	2.1
1	A	762	HIS	2.1
2	B	33	LEU	2.1
1	A	450	TYR	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(\AA^2)	Q<0.9
7	NAG	A	1010	14/15	0.93	0.19	-0.80	73,83,91,95	0
5	NAG	A	1001	14/15	0.94	0.20	-0.91	57,77,88,91	0
9	NAG	B	702	14/15	0.93	0.23	-1.10	101,128,141,146	0
9	NAG	B	704	14/15	0.92	0.11	-1.14	104,130,146,163	0
7	NAG	A	1019	14/15	0.81	0.16	-	170,191,197,202	0
7	MAN	A	1015	11/12	0.89	0.13	-	150,152,158,164	0
5	NAG	A	1002	14/15	0.89	0.21	-	110,131,145,149	0
9	NAG	A	1018	14/15	0.91	0.20	-	178,183,190,192	0
9	NAG	A	1017	14/15	0.87	0.18	-	146,165,175,177	0
7	MAN	A	1024	11/12	0.75	0.18	-	229,231,239,245	0
5	BMA	A	1005	11/12	0.90	0.10	-	212,215,218,219	0
7	NAG	A	1020	14/15	0.86	0.14	-	207,215,221,224	0
9	NAG	B	703	14/15	0.89	0.20	-	139,147,159,160	0
5	MAN	A	1004	11/12	0.84	0.14	-	172,188,221,223	0
5	BMA	A	1003	11/12	0.80	0.16	-	161,174,187,190	0
9	NAG	B	705	14/15	0.80	0.17	-	175,184,186,188	0
6	NAG	A	1008	14/15	0.82	0.22	-	198,212,220,228	0
6	NAG	A	1007	14/15	0.89	0.13	-	132,150,165,181	0
5	MAN	A	1006	11/12	0.67	0.29	-	184,196,205,209	0
7	MAN	A	1022	11/12	0.80	0.18	-	253,259,264,264	0
7	NAG	A	1011	14/15	0.95	0.17	-	66,95,109,116	0
7	BMA	A	1021	11/12	0.75	0.20	-	225,229,266,267	0
7	MAN	A	1013	11/12	0.86	0.16	-	147,158,173,174	0
6	BMA	A	1009	11/12	0.61	0.23	-	223,233,237,238	0
7	MAN	A	1023	11/12	0.82	0.24	-	242,256,261,263	0
7	MAN	A	1014	11/12	0.86	0.18	-	170,178,183,185	0
7	BMA	A	1012	11/12	0.93	0.14	-	112,127,139,141	0

6.4 Ligands ⓘ

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
10	MN	A	1028	1/1	0.98	0.25	1.31	169,169,169,169	0
11	NO3	A	1029	4/4	0.94	0.29	0.82	95,98,100,103	0
11	NO3	A	1030	4/4	0.96	0.27	0.29	81,83,84,87	0
10	MN	B	707	1/1	0.77	0.16	0.11	271,271,271,271	0
8	NAG	B	706	14/15	0.85	0.18	-0.24	145,154,161,162	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
10	MN	A	1027	1/1	0.75	0.16	-0.91	187,187,187,187	0
10	MN	A	1025	1/1	0.84	0.06	-0.93	241,241,241,241	0
10	MN	A	1026	1/1	0.96	0.07	-1.39	181,181,181,181	0
8	NAG	A	1016	14/15	0.51	0.38	-	190,203,212,212	0
8	NAG	B	701	14/15	0.89	0.17	-	120,143,162,168	0

6.5 Other polymers [i](#)

There are no such residues in this entry.