



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 07:27 PM GMT

PDB ID : 1FQ8
Title : X-RAY STRUCTURE OF DIFLUOROSTATINE INHIBITOR CP81,198
BOUND TO SACCHAROPEPSIN
Authors : Cronin, N.B.; Badasso, M.O.; Tickle, I.J.; Dreyer, T.; Hoover, D.J.; Rosati,
R.L.; Humblet, C.C.; Lunney, E.A.; Cooper, J.B.
Deposited on : 2000-09-04
Resolution : 2.80 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

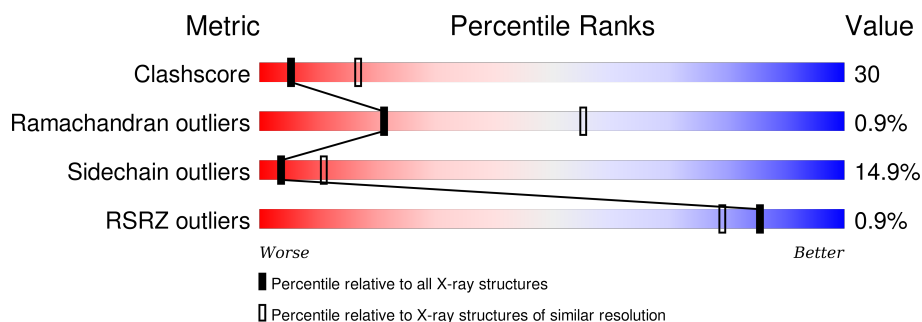
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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))
Clashscore	102246	2827 (2.80-2.80)
Ramachandran outliers	100387	2782 (2.80-2.80)
Sidechain outliers	100360	2784 (2.80-2.80)
RSRZ outliers	91569	2404 (2.80-2.80)

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	329	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	NAG	A	333	X	-	-	-
2	MAN	A	334	X	-	-	-
2	MAN	A	335	X	-	-	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 2812 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 SACCHAROPEPSIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	329	Total	C	N	O	S	0	0	0
			2528	1618	396	508	6			

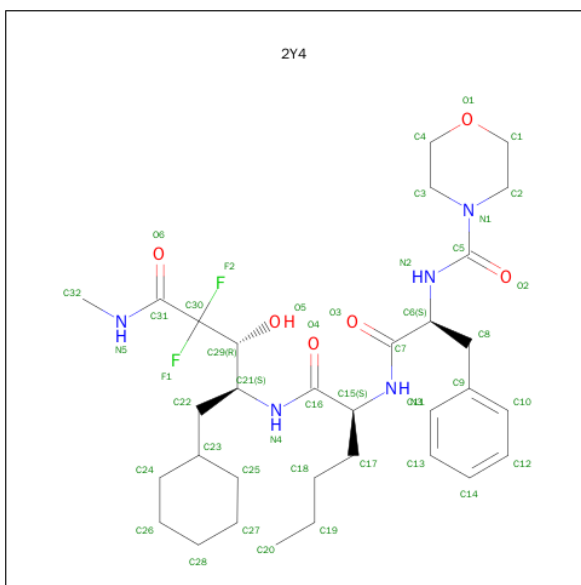
There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	315	ILE	LEU	CONFLICT	UNP P07267

- Molecule 2 is a polymer of unknown type called SUGAR (5-MER).

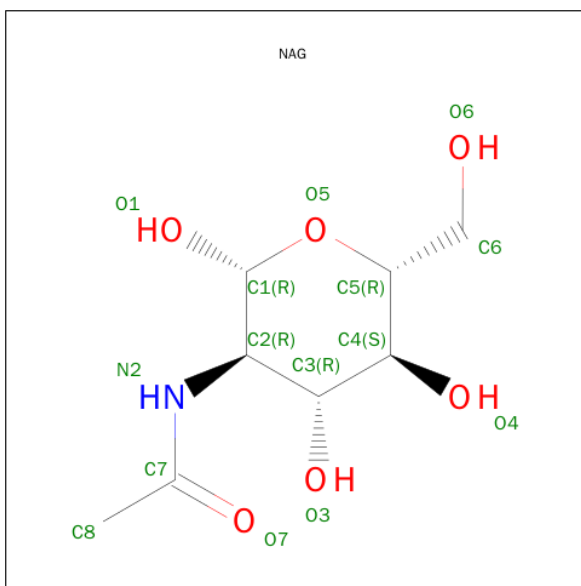
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	5	Total	C	N	O	0	0
			61	34	2	25		

- Molecule 3 is N-[(2S)-1-[[[(2S)-1-[[[(2S,3R)-1-CYCLOHEXYL-4,4-DIFLUORO-3-HYDROXY-5-(METHYLAMINO)-5-OXO-PENTAN-2-YL]AMINO]-1-OXO-HEXAN-2-YL]AMINO]-1-OXO-3-PHENYL-PROPAN-2-YL]MORPHOLINE-4-CARBOXAMIDE (three-letter code: 2Y4) (formula: C₃₂H₄₉F₂N₅O₆).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	F	N	O	0	0
			45	32	2	5	6		

- Molecule 4 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



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

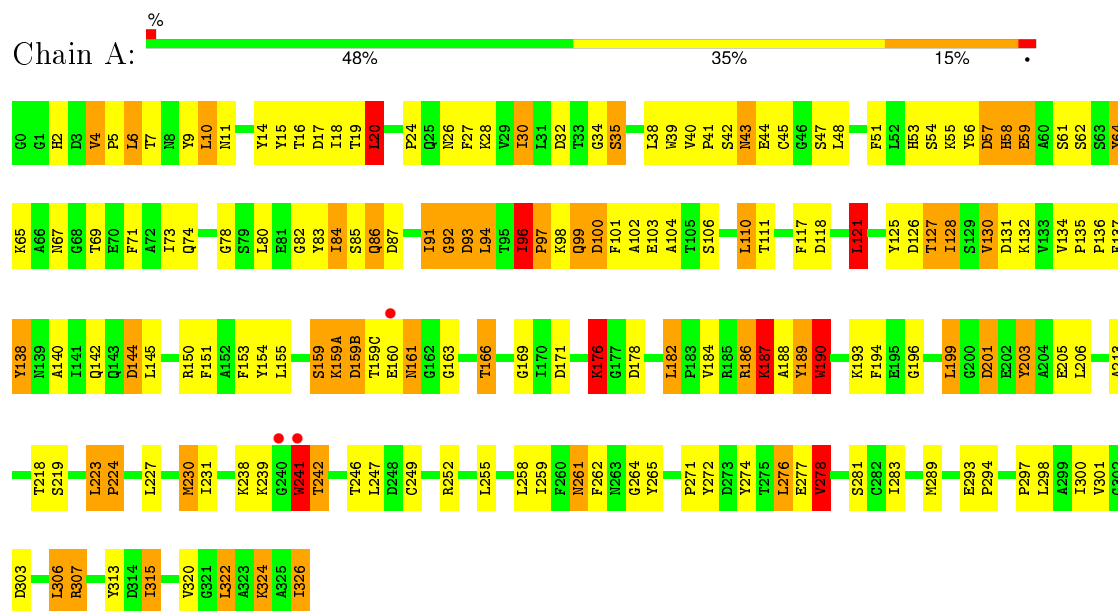
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	164	Total 164	O 164	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: SACCHAROPEPSIN



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	86.82Å 86.82Å 110.24Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	15.40 – 2.80 15.41 – 2.80	Depositor EDS
% Data completeness (in resolution range)	100.0 (15.40-2.80) 99.5 (15.41-2.80)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	7.48 (at 2.81Å)	Xtriage
Refinement program	RESTRAIN	Depositor
R, R_{free}	0.190 , 0.270 0.173 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	49.6	Xtriage
Anisotropy	0.344	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.24 , 70.1	EDS
Estimated twinning fraction	0.027 for -h,-k,l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 12107 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	2812	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.62% 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: 2Y4, BMA, NAG, 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	0.78	0/2592	1.89	86/3526 (2.4%)

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	6	0
2	A	4	0
All	All	10	0

There are no bond length outliers.

All (86) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	93	ASP	CB-CG-OD1	-11.98	107.51	118.30
1	A	130	VAL	CA-CB-CG1	9.93	125.80	110.90
1	A	189	TYR	CA-CB-CG	9.78	131.98	113.40
1	A	150	ARG	NE-CZ-NH2	-9.68	115.46	120.30
1	A	278	VAL	CA-CB-CG1	9.57	125.26	110.90
1	A	150	ARG	NE-CZ-NH1	8.95	124.78	120.30
1	A	186	ARG	CG-CD-NE	8.92	130.54	111.80
1	A	324	LYS	CA-CB-CG	8.85	132.87	113.40
1	A	239	LYS	CD-CE-NZ	8.54	131.35	111.70
1	A	289	MET	CG-SD-CE	8.50	113.80	100.20
1	A	230	MET	CG-SD-CE	-8.48	86.63	100.20
1	A	203	TYR	CB-CG-CD1	8.14	125.89	121.00
1	A	178	ASP	CB-CG-OD1	7.91	125.42	118.30
1	A	186	ARG	CA-CB-CG	7.90	130.79	113.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	4	VAL	CA-CB-CG1	7.82	122.63	110.90
1	A	301	VAL	CA-CB-CG1	7.76	122.54	110.90
1	A	64	TYR	CB-CG-CD1	-7.75	116.35	121.00
1	A	153	PHE	CB-CG-CD1	-7.63	115.46	120.80
1	A	96	ILE	C-N-CD	7.60	144.37	128.40
1	A	132	LYS	CD-CE-NZ	7.50	128.94	111.70
1	A	34	GLY	C-N-CA	7.49	140.41	121.70
1	A	62	SER	C-N-CA	7.37	140.12	121.70
1	A	166	THR	CA-CB-CG2	7.35	122.68	112.40
1	A	153	PHE	CG-CD2-CE2	-7.33	112.74	120.80
1	A	298	LEU	CA-CB-CG	7.33	132.15	115.30
1	A	138	TYR	CB-CG-CD1	7.32	125.39	121.00
1	A	201	ASP	CB-CG-OD1	7.26	124.84	118.30
1	A	131	ASP	CB-CG-OD1	7.26	124.83	118.30
1	A	100	ASP	CB-CG-OD1	7.18	124.76	118.30
1	A	92	GLY	C-N-CA	7.16	139.59	121.70
1	A	171	ASP	CB-CG-OD1	7.14	124.72	118.30
1	A	176	LYS	CD-CE-NZ	7.12	128.07	111.70
1	A	94	LEU	CA-CB-CG	7.04	131.50	115.30
1	A	57	ASP	CB-CG-OD1	-6.91	112.08	118.30
1	A	320	VAL	CG1-CB-CG2	6.88	121.91	110.90
1	A	154	TYR	CB-CG-CD2	-6.87	116.88	121.00
1	A	94	LEU	CB-CG-CD2	-6.80	99.44	111.00
1	A	322	LEU	CB-CG-CD2	6.78	122.52	111.00
1	A	307	ARG	CG-CD-NE	-6.75	97.62	111.80
1	A	150	ARG	CD-NE-CZ	6.70	132.98	123.60
1	A	138	TYR	CA-CB-CG	6.69	126.11	113.40
1	A	186	ARG	NE-CZ-NH1	-6.59	117.00	120.30
1	A	6	LEU	CB-CG-CD1	6.57	122.17	111.00
1	A	189	TYR	CA-C-N	6.48	131.46	117.20
1	A	252	ARG	CG-CD-NE	6.46	125.36	111.80
1	A	126	ASP	C-N-CA	6.41	137.72	121.70
1	A	20	LEU	C-N-CA	-6.25	109.17	122.30
1	A	96	ILE	CA-CB-CG2	6.25	123.40	110.90
1	A	199	LEU	CA-CB-CG	6.13	129.41	115.30
1	A	94	LEU	CB-CG-CD1	-6.13	100.58	111.00
1	A	187	LYS	CD-CE-NZ	-6.10	97.67	111.70
1	A	186	ARG	NE-CZ-NH2	6.03	123.32	120.30
1	A	306	LEU	CA-CB-CG	6.01	129.13	115.30
1	A	201	ASP	CB-CG-OD2	-5.99	112.91	118.30
1	A	258	LEU	CA-CB-CG	5.96	129.01	115.30
1	A	194	PHE	CB-CG-CD1	5.94	124.96	120.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	184	VAL	CA-CB-CG2	-5.93	102.00	110.90
1	A	315	ILE	CA-CB-CG1	5.81	122.04	111.00
1	A	276	LEU	CA-CB-CG	5.80	128.64	115.30
1	A	97	PRO	N-CD-CG	-5.77	94.55	103.20
1	A	272	TYR	C-N-CA	5.71	135.98	121.70
1	A	93	ASP	CB-CG-OD2	5.69	123.42	118.30
1	A	86	GLN	N-CA-CB	5.67	120.82	110.60
1	A	86	GLN	CB-CA-C	5.65	121.70	110.40
1	A	144	ASP	C-N-CA	5.63	135.78	121.70
1	A	252	ARG	NE-CZ-NH2	5.62	123.11	120.30
1	A	150	ARG	CB-CG-CD	-5.56	97.14	111.60
1	A	182	LEU	CB-CG-CD1	5.56	120.45	111.00
1	A	190	TRP	CA-C-O	5.52	131.68	120.10
1	A	47	SER	O-C-N	5.50	131.50	122.70
1	A	20	LEU	O-C-N	-5.47	113.91	123.20
1	A	306	LEU	N-CA-CB	5.44	121.29	110.40
1	A	189	TYR	CB-CG-CD1	5.42	124.25	121.00
1	A	93	ASP	CA-CB-CG	-5.38	101.57	113.40
1	A	78	GLY	C-N-CA	5.32	135.01	121.70
1	A	97	PRO	CA-N-CD	-5.25	104.14	111.50
1	A	203	TYR	CB-CG-CD2	-5.22	117.87	121.00
1	A	99	GLN	C-N-CA	-5.22	108.66	121.70
1	A	134	VAL	CA-CB-CG1	5.17	118.65	110.90
1	A	189	TYR	O-C-N	-5.16	114.44	122.70
1	A	84	ILE	N-CA-CB	5.12	122.58	110.80
1	A	227	LEU	CB-CG-CD2	-5.10	102.33	111.00
1	A	121	LEU	CA-CB-CG	5.10	127.03	115.30
1	A	155	LEU	CB-CG-CD1	-5.10	102.33	111.00
1	A	223	LEU	CA-CB-CG	5.10	127.03	115.30
1	A	255	LEU	CB-CG-CD2	-5.01	102.49	111.00

All (10) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	16	THR	CB
1	A	63	SER	CA
1	A	96	ILE	CB
1	A	166	THR	CB
1	A	241	TRP	CA
1	A	315	ILE	CB
2	A	333	NAG	C1
2	A	334	MAN	C3,C1

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Mol	Chain	Res	Type	Atom
2	A	335	MAN	C1

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	2528	0	2401	146	5
2	A	61	0	50	9	1
3	A	45	0	49	3	0
4	A	14	0	13	0	0
5	A	164	0	0	26	6
All	All	2812	0	2513	155	8

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 30.

All (155) 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:69:THR:HB	1:A:84:ILE:HD13	1.21	1.10
1:A:242:THR:CG2	1:A:242:THR:O	2.02	1.07
1:A:242:THR:HG22	1:A:242:THR:O	1.37	1.06
1:A:189:TYR:O	1:A:190:TRP:HB2	1.59	1.01
1:A:261:ASN:HD22	1:A:261:ASN:C	1.66	0.98
1:A:97:PRO:HD3	5:A:515:HOH:O	1.65	0.97
1:A:261:ASN:HD21	1:A:264:GLY:H	1.19	0.90
1:A:97:PRO:O	1:A:98:LYS:HB2	1.69	0.90
1:A:69:THR:HB	1:A:84:ILE:CD1	2.02	0.89
1:A:159(B):ASP:HB3	5:A:635:HOH:O	1.73	0.88
1:A:128:ILE:O	1:A:128:ILE:HG12	1.73	0.87
1:A:58:HIS:HB2	1:A:64:TYR:CD2	2.10	0.87
2:A:335:MAN:C2	2:A:336:BMA:C1	2.61	0.78
1:A:161:ASN:HA	5:A:661:HOH:O	1.84	0.77
1:A:121:LEU:C	1:A:121:LEU:HD12	2.05	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:176:LYS:HG2	1:A:326:ILE:HD12	1.68	0.75
1:A:201:ASP:HB2	5:A:638:HOH:O	1.86	0.75
1:A:84:ILE:HG22	1:A:85:SER:N	1.99	0.75
1:A:159:SER:HA	5:A:654:HOH:O	1.87	0.75
1:A:32:ASP:OD2	1:A:35:SER:HB2	1.87	0.74
1:A:84:ILE:CG2	1:A:85:SER:N	2.51	0.74
1:A:159(C):THR:OG1	1:A:160:GLU:N	2.20	0.73
1:A:53:HIS:HB3	1:A:118:ASP:OD1	1.88	0.72
1:A:84:ILE:CG2	1:A:85:SER:H	2.01	0.72
1:A:38:LEU:HD13	1:A:121:LEU:HB2	1.73	0.71
1:A:182:LEU:CD1	1:A:322:LEU:HD12	2.20	0.71
2:A:335:MAN:O2	2:A:335:MAN:C2	2.38	0.70
1:A:84:ILE:H	1:A:84:ILE:HD12	1.55	0.70
1:A:57:ASP:OD1	1:A:59:GLU:HG3	1.91	0.70
1:A:159(C):THR:HA	5:A:640:HOH:O	1.93	0.69
1:A:86:GLN:O	1:A:87:ASP:HB2	1.91	0.69
1:A:4:VAL:HG23	1:A:16:THR:HG21	1.73	0.69
2:A:334:MAN:C1	2:A:334:MAN:O2	2.40	0.69
1:A:16:THR:HG22	1:A:17:ASP:H	1.57	0.69
1:A:137:PHE:HE2	1:A:315:ILE:CG2	2.04	0.69
1:A:65:LYS:HE3	5:A:649:HOH:O	1.92	0.68
1:A:58:HIS:HB2	1:A:64:TYR:CG	2.28	0.68
1:A:64:TYR:CD1	1:A:65:LYS:N	2.62	0.68
2:A:334:MAN:H3	2:A:334:MAN:O2	1.93	0.68
1:A:48:LEU:O	1:A:48:LEU:HD12	1.92	0.68
1:A:313:TYR:N	1:A:313:TYR:CD1	2.62	0.67
2:A:334:MAN:C2	2:A:334:MAN:O2	2.43	0.67
1:A:261:ASN:HD21	1:A:264:GLY:N	1.92	0.66
2:A:335:MAN:O2	2:A:335:MAN:C1	2.44	0.66
1:A:261:ASN:HD22	1:A:262:PHE:N	1.93	0.65
1:A:303:ASP:O	1:A:307:ARG:HG3	1.96	0.65
1:A:96:ILE:HA	5:A:515:HOH:O	1.96	0.65
1:A:189:TYR:O	1:A:190:TRP:CB	2.31	0.65
1:A:182:LEU:HD11	1:A:322:LEU:HD12	1.79	0.64
1:A:261:ASN:ND2	1:A:261:ASN:C	2.42	0.64
1:A:161:ASN:C	1:A:161:ASN:HD22	2.00	0.63
1:A:51:PHE:HB3	5:A:555:HOH:O	1.99	0.62
1:A:16:THR:HG22	1:A:17:ASP:N	2.15	0.62
1:A:261:ASN:ND2	1:A:264:GLY:H	1.95	0.61
1:A:86:GLN:NE2	1:A:98:LYS:HA	2.16	0.61
1:A:30:ILE:HD12	1:A:117:PHE:CD2	2.36	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:84:ILE:N	1:A:84:ILE:HD12	2.16	0.60
2:A:334:MAN:C3	2:A:334:MAN:O2	2.49	0.60
1:A:86:GLN:HE21	1:A:98:LYS:HA	1.66	0.60
1:A:196:GLY:HA3	1:A:203:TYR:OH	2.03	0.59
2:A:335:MAN:O2	2:A:335:MAN:C3	2.50	0.59
1:A:6:LEU:HB2	1:A:163:GLY:CA	2.33	0.58
1:A:277:GLU:HG2	5:A:557:HOH:O	2.03	0.58
1:A:10:LEU:HD13	5:A:501:HOH:O	2.03	0.58
1:A:205:GLU:O	1:A:206:LEU:HD23	2.04	0.57
1:A:213:ALA:HB3	1:A:300:ILE:HG12	1.87	0.57
1:A:166:THR:HG22	1:A:169:GLY:O	2.05	0.56
1:A:84:ILE:CG2	1:A:100:ASP:HB3	2.36	0.56
1:A:110:LEU:HD12	5:A:642:HOH:O	2.04	0.56
1:A:145:LEU:HD22	5:A:524:HOH:O	2.04	0.56
1:A:65:LYS:HB3	1:A:86:GLN:HG2	1.87	0.56
1:A:271:PRO:HA	1:A:274:TYR:CE2	2.41	0.56
1:A:246:THR:HG22	1:A:283:ILE:HG22	1.89	0.56
1:A:9:TYR:CZ	1:A:10:LEU:HD22	2.42	0.55
1:A:97:PRO:O	1:A:98:LYS:CB	2.45	0.55
1:A:4:VAL:CG2	1:A:16:THR:HG21	2.37	0.55
1:A:151:PHE:HA	1:A:166:THR:O	2.08	0.54
1:A:84:ILE:HG21	1:A:100:ASP:HB3	1.88	0.54
1:A:140:ALA:O	1:A:145:LEU:HB2	2.08	0.54
1:A:67:ASN:C	1:A:67:ASN:OD1	2.46	0.54
1:A:48:LEU:HA	5:A:555:HOH:O	2.08	0.54
1:A:219:SER:HB2	3:A:400:2Y4:H3	1.89	0.54
1:A:6:LEU:HB2	1:A:163:GLY:HA3	1.90	0.54
1:A:188:ALA:HB3	5:A:566:HOH:O	2.07	0.53
1:A:276:LEU:HD23	1:A:283:ILE:HD11	1.91	0.53
1:A:223:LEU:HB3	1:A:224:PRO:HD2	1.91	0.53
1:A:15:TYR:CE2	1:A:28:LYS:HD3	2.44	0.53
1:A:315:ILE:HD12	5:A:532:HOH:O	2.08	0.53
1:A:249:CYS:HB3	5:A:550:HOH:O	2.08	0.53
1:A:86:GLN:O	1:A:87:ASP:CB	2.53	0.52
1:A:69:THR:CB	1:A:84:ILE:HD13	2.15	0.52
1:A:42:SER:O	1:A:55:LYS:HG2	2.10	0.52
1:A:38:LEU:HD12	1:A:39:TRP:H	1.76	0.50
1:A:6:LEU:HB2	1:A:163:GLY:C	2.31	0.50
1:A:84:ILE:HG23	1:A:85:SER:H	1.75	0.50
1:A:18:ILE:HG13	1:A:18:ILE:O	2.09	0.50
1:A:48:LEU:HD12	1:A:48:LEU:C	2.32	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:30:ILE:CD1	1:A:117:PHE:CD2	2.95	0.49
1:A:30:ILE:HD12	1:A:117:PHE:CG	2.47	0.49
1:A:218:THR:HG23	3:A:400:2Y4:H15	1.93	0.49
1:A:17:ASP:OD2	1:A:28:LYS:HE2	2.13	0.49
1:A:91:ILE:O	1:A:92:GLY:C	2.50	0.49
1:A:2:HIS:HD2	1:A:93:ASP:OD2	1.96	0.48
1:A:2:HIS:HE1	5:A:632:HOH:O	1.96	0.48
1:A:324:LYS:HD3	5:A:634:HOH:O	2.13	0.48
1:A:127:THR:O	1:A:127:THR:HG22	2.13	0.48
1:A:83:TYR:HD1	1:A:84:ILE:O	1.97	0.47
1:A:121:LEU:C	1:A:121:LEU:CD1	2.79	0.47
1:A:11:ASN:HB3	1:A:307:ARG:NH1	2.29	0.47
1:A:121:LEU:HD12	1:A:121:LEU:O	2.15	0.47
1:A:187:LYS:HZ3	1:A:187:LYS:HG3	1.27	0.46
1:A:278:VAL:O	1:A:278:VAL:HG12	2.16	0.46
1:A:41:PRO:HA	1:A:104:ALA:HB3	1.98	0.46
1:A:39:TRP:HE3	1:A:40:VAL:N	2.14	0.46
1:A:58:HIS:HB2	1:A:64:TYR:CE2	2.50	0.45
1:A:137:PHE:HE2	1:A:315:ILE:HG21	1.80	0.45
1:A:9:TYR:CE1	1:A:10:LEU:HD22	2.51	0.45
1:A:41:PRO:HG2	1:A:54:SER:O	2.17	0.45
1:A:51:PHE:CB	5:A:555:HOH:O	2.61	0.45
1:A:86:GLN:HG3	1:A:87:ASP:N	2.33	0.44
1:A:99:GLN:OE1	1:A:136:PRO:HA	2.17	0.44
1:A:176:LYS:HB2	1:A:176:LYS:HE3	1.59	0.44
1:A:84:ILE:HG23	1:A:101:PHE:O	2.17	0.44
1:A:159(A):LYS:HB2	1:A:159(A):LYS:HE2	1.82	0.44
1:A:28:LYS:HD2	5:A:546:HOH:O	2.17	0.44
1:A:326:ILE:OXT	1:A:326:ILE:HG22	2.17	0.43
1:A:5:PRO:O	1:A:16:THR:HG23	2.18	0.43
1:A:94:LEU:HD23	5:A:524:HOH:O	2.18	0.43
1:A:219:SER:CB	3:A:400:2Y4:H3	2.48	0.43
1:A:56:TYR:HD2	1:A:103:GLU:OE1	2.02	0.43
1:A:71:PHE:CE1	1:A:130:VAL:CG1	3.02	0.43
1:A:135:PRO:HG2	1:A:138:TYR:CD1	2.53	0.43
1:A:84:ILE:HG22	1:A:85:SER:H	1.68	0.43
1:A:128:ILE:HD12	1:A:128:ILE:HG21	1.64	0.43
1:A:7:THR:O	1:A:14:TYR:HA	2.19	0.43
2:A:335:MAN:C3	2:A:336:BMA:C1	2.96	0.42
1:A:294:PRO:HG2	5:A:664:HOH:O	2.18	0.42
1:A:145:LEU:HD11	5:A:515:HOH:O	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:137:PHE:HE2	1:A:315:ILE:HG23	1.80	0.42
1:A:261:ASN:HA	1:A:265:TYR:O	2.19	0.42
1:A:73:ILE:CG2	1:A:80:LEU:HD21	2.49	0.42
1:A:19:THR:HB	1:A:24:PRO:HB3	2.00	0.42
1:A:259:ILE:HG21	1:A:259:ILE:HD13	1.79	0.42
1:A:297:PRO:HD2	5:A:628:HOH:O	2.19	0.42
1:A:186:ARG:HD2	1:A:186:ARG:HH11	1.58	0.42
1:A:64:TYR:CD1	1:A:64:TYR:C	2.89	0.41
1:A:125:TYR:CD2	1:A:187:LYS:HB3	2.55	0.41
1:A:61:SER:HB3	1:A:64:TYR:HB2	2.03	0.41
1:A:28:LYS:HE3	5:A:546:HOH:O	2.19	0.41
1:A:82:GLY:HA3	1:A:103:GLU:O	2.20	0.41
1:A:73:ILE:CG2	1:A:80:LEU:CD2	2.98	0.41
1:A:101:PHE:HD1	1:A:102:ALA:N	2.19	0.40
1:A:20:LEU:HA	1:A:20:LEU:HD12	1.36	0.40
1:A:223:LEU:O	1:A:224:PRO:C	2.59	0.40
1:A:127:THR:O	1:A:127:THR:CG2	2.70	0.40

All (8) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:230:MET:CB	5:A:633:HOH:O[5_675]	1.17	1.03
1:A:230:MET:CG	5:A:633:HOH:O[5_675]	1.19	1.01
5:A:622:HOH:O	5:A:623:HOH:O[3_565]	1.89	0.31
2:A:336:BMA:O3	5:A:654:HOH:O[3_565]	1.93	0.27
1:A:230:MET:SD	5:A:633:HOH:O[5_675]	1.95	0.25
5:A:654:HOH:O	5:A:655:HOH:O[2_664]	1.99	0.21
1:A:43:ASN:O	1:A:241:TRP:CZ2[6_555]	1.99	0.21
1:A:45:CYS:O	1:A:241:TRP:CZ3[6_555]	2.07	0.13

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	327/329 (99%)	293 (90%)	31 (10%)	3 (1%)	21	55

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	241	TRP
1	A	127	THR
1	A	190	TRP

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	268/268 (100%)	228 (85%)	40 (15%)	4	11

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	LEU
1	A	20	LEU
1	A	26	ASN
1	A	27	PHE
1	A	30	ILE
1	A	35	SER
1	A	43	ASN
1	A	44	GLU
1	A	58	HIS
1	A	59	GLU
1	A	74	GLN
1	A	91	ILE
1	A	96	ILE
1	A	106	SER
1	A	110	LEU
1	A	111	THR
1	A	121	LEU

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Mol	Chain	Res	Type
1	A	128	ILE
1	A	142	GLN
1	A	144	ASP
1	A	159	SER
1	A	159(A)	LYS
1	A	159(B)	ASP
1	A	161	ASN
1	A	176	LYS
1	A	187	LYS
1	A	193	LYS
1	A	199	LEU
1	A	224	PRO
1	A	231	ILE
1	A	238	LYS
1	A	241	TRP
1	A	242	THR
1	A	247	LEU
1	A	261	ASN
1	A	278	VAL
1	A	281	SER
1	A	293	GLU
1	A	306	LEU
1	A	326	ILE

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	2	HIS
1	A	43	ASN
1	A	86	GLN
1	A	161	ASN
1	A	232	ASN
1	A	261	ASN
1	A	317	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 ⓘ

5 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	332	1,2	14,14,15	1.09	1 (7%)	15,19,21	2.25	3 (20%)
2	NAG	A	333	2	14,14,15	1.21	2 (14%)	15,19,21	2.62	5 (33%)
2	MAN	A	334	2	11,11,12	13.48	2 (18%)	14,15,17	7.22	5 (35%)
2	MAN	A	335	2	11,11,12	12.92	3 (27%)	14,15,17	7.35	9 (64%)
2	BMA	A	336	2	11,11,12	1.30	3 (27%)	14,15,17	2.69	4 (28%)

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	332	1,2	-	0/6/23/26	0/1/1/1
2	NAG	A	333	2	1/1/5/7	0/6/23/26	0/1/1/1
2	MAN	A	334	2	2/2/4/5	0/2/19/22	0/1/1/1
2	MAN	A	335	2	1/1/4/5	0/2/19/22	0/1/1/1
2	BMA	A	336	2	-	0/2/19/22	0/1/1/1

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	335	MAN	C2-C3	-4.03	1.47	1.52
2	A	335	MAN	C4-C3	-3.02	1.44	1.52
2	A	336	BMA	C2-C3	-2.42	1.49	1.52
2	A	336	BMA	O4-C4	2.17	1.48	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	334	MAN	O4-C4	2.30	1.48	1.43
2	A	333	NAG	C2-N2	2.34	1.50	1.46
2	A	336	BMA	O5-C1	2.47	1.47	1.43
2	A	333	NAG	O4-C4	2.48	1.48	1.43
2	A	332	NAG	C2-N2	2.53	1.50	1.46
2	A	335	MAN	O2-C2	42.48	2.38	1.43
2	A	334	MAN	O2-C2	44.57	2.43	1.43

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	334	MAN	O2-C2-C1	-19.37	70.37	109.21
2	A	334	MAN	O2-C2-C3	-17.92	74.07	110.12
2	A	335	MAN	O2-C2-C1	-17.80	73.52	109.21
2	A	335	MAN	O2-C2-C3	-16.58	76.77	110.12
2	A	332	NAG	C2-N2-C7	-5.49	115.99	123.04
2	A	335	MAN	O6-C6-C5	-4.45	96.62	111.33
2	A	333	NAG	C2-N2-C7	-2.75	119.51	123.04
2	A	333	NAG	C3-C4-C5	-2.70	105.49	110.20
2	A	332	NAG	O3-C3-C4	-2.51	104.68	110.34
2	A	334	MAN	O3-C3-C4	-2.42	104.89	110.34
2	A	333	NAG	O4-C4-C5	-2.31	103.11	109.24
2	A	334	MAN	O5-C5-C6	-2.24	102.49	107.35
2	A	336	BMA	C2-C3-C4	2.21	114.79	111.04
2	A	336	BMA	C6-C5-C4	2.39	118.92	113.02
2	A	335	MAN	C3-C4-C5	2.48	114.52	110.20
2	A	333	NAG	O7-C7-C8	2.84	127.28	122.06
2	A	335	MAN	C1-O5-C5	2.99	116.04	112.25
2	A	336	BMA	C1-O5-C5	3.64	116.87	112.25
2	A	334	MAN	C1-O5-C5	3.67	116.91	112.25
2	A	335	MAN	O5-C1-C2	3.92	117.21	110.86
2	A	335	MAN	C1-C2-C3	4.14	114.44	109.54
2	A	335	MAN	O3-C3-C2	5.16	119.32	110.00
2	A	332	NAG	C1-O5-C5	5.22	118.87	112.25
2	A	335	MAN	C2-C3-C4	7.48	123.75	111.04
2	A	336	BMA	C1-C2-C3	7.80	118.77	109.54
2	A	333	NAG	C1-O5-C5	7.91	122.28	112.25

All (4) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	333	NAG	C1

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Mol	Chain	Res	Type	Atom
2	A	334	MAN	C3
2	A	334	MAN	C1
2	A	335	MAN	C1

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	334	MAN	4	0
2	A	335	MAN	5	0
2	A	336	BMA	2	1

5.6 Ligand geometry [i](#)

2 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	NAG	A	337	1	14,14,15	1.39	3 (21%)	15,19,21	2.91	7 (46%)
3	2Y4	A	400	-	45,47,47	0.74	1 (2%)	55,63,63	1.67	12 (21%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	A	337	1	-	0/6/23/26	0/1/1/1
3	2Y4	A	400	-	-	0/51/70/70	0/3/3/3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	400	2Y4	O6-C31	2.23	1.27	1.22
4	A	337	NAG	O5-C5	2.53	1.49	1.43
4	A	337	NAG	O5-C1	2.53	1.48	1.43
4	A	337	NAG	O4-C4	2.54	1.49	1.43

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	337	NAG	C2-N2-C7	-6.46	114.75	123.04
3	A	400	2Y4	O6-C31-N5	-4.53	116.74	122.86
4	A	337	NAG	C1-O5-C5	-4.01	107.16	112.25
4	A	337	NAG	O5-C5-C6	-3.92	98.85	107.35
4	A	337	NAG	C8-C7-N2	-3.64	109.14	116.11
4	A	337	NAG	O6-C6-C5	-3.63	99.34	111.33
3	A	400	2Y4	F2-C30-C31	-3.35	105.20	109.78
3	A	400	2Y4	F2-C30-F1	-2.73	98.51	105.45
3	A	400	2Y4	F1-C30-C31	-2.71	106.08	109.78
3	A	400	2Y4	C26-C24-C23	-2.21	108.65	112.22
3	A	400	2Y4	C6-C7-N3	-2.20	111.71	116.78
3	A	400	2Y4	C18-C17-C15	-2.19	107.18	113.91
3	A	400	2Y4	C25-C23-C24	2.13	114.62	109.26
4	A	337	NAG	C4-C3-C2	2.89	115.73	111.23
3	A	400	2Y4	C21-N4-C16	3.01	128.72	123.13
4	A	337	NAG	O7-C7-C8	3.01	127.58	122.06
3	A	400	2Y4	C29-C21-N4	3.26	116.22	110.11
3	A	400	2Y4	C3-N1-C2	3.54	119.12	112.56
3	A	400	2Y4	C4-C3-N1	4.85	120.30	109.90

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	400	2Y4	3	0

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	329/329 (100%)	-0.87	3 (0%) 85 79	14, 26, 54, 116	4 (1%)

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	241	TRP	3.7
1	A	240	GLY	3.0
1	A	160	GLU	2.0

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

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

6.3 Carbohydrates ⓘ

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	332	14/15	0.94	0.14	0.02	34,49,72,72	0
2	BMA	A	336	11/12	0.84	0.23	-	53,67,124,149	0
2	NAG	A	333	14/15	0.96	0.12	-	18,33,47,52	0
2	MAN	A	334	11/12	0.89	0.21	-	25,42,54,56	0
2	MAN	A	335	11/12	0.94	0.24	-	28,40,44,55	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	2Y4	A	400	45/45	0.91	0.15	0.75	1,24,36,44	0
4	NAG	A	337	14/15	0.89	0.14	-	42,53,92,120	0

6.5 Other polymers [i](#)

There are no such residues in this entry.