



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 15, 2017 – 01:12 am GMT

PDB ID : 4BH1
Title : H5 (tyTy) Influenza Virus Haemagglutinin in Complex with Avian Receptor Analogue 3'-SLN
Authors : Xiong, X.; Coombs, P.J.; Martin, S.R.; Liu, J.; Xiao, H.; McCauley, J.W.; Locher, K.; Walker, P.A.; Collins, P.J.; Kawaoka, Y.; Skehel, J.J.; Gamblin, S.J.
Deposited on : 2013-03-29
Resolution : 2.15 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	:	1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix)	:	1.9-1692
EDS	:	trunk28620
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
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)	:	recalc28949

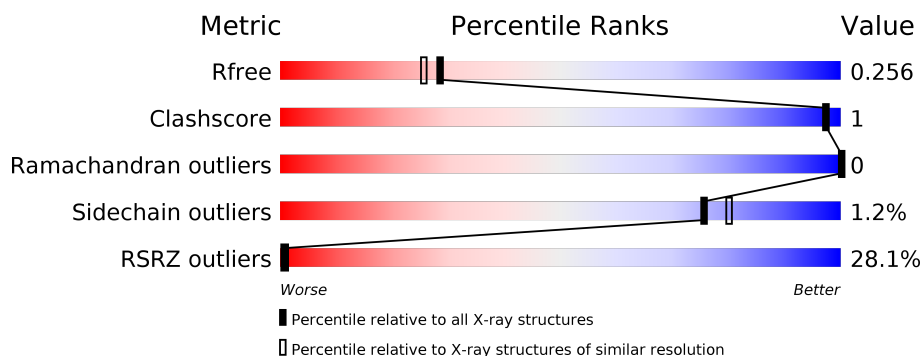
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.15 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}	100719	1170 (2.16-2.16)
Clashscore	112137	1278 (2.16-2.16)
Ramachandran outliers	110173	1256 (2.16-2.16)
Sidechain outliers	110143	1255 (2.16-2.16)
RSRZ outliers	101464	1175 (2.16-2.16)

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	326	<div> <div>9%</div> <div>94%</div> <div>• •</div> </div>
1	C	326	<div> <div>16%</div> <div>94%</div> <div>• •</div> </div>
1	E	326	<div> <div>10%</div> <div>93%</div> <div>• •</div> </div>
2	B	166	<div> <div>55%</div> <div>86%</div> <div>• 13%</div> </div>
2	D	166	<div> <div>51%</div> <div>84%</div> <div>• 12%</div> </div>
2	F	166	<div> <div>60%</div> <div>86%</div> <div>• 13%</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
4	GAL	A	1322	-	-	-	X
5	PO4	A	1325	-	-	-	X
5	PO4	C	1324	-	-	-	X
5	PO4	E	1324	-	-	-	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 11980 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 HEMAGGLUTININ.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	319	Total	C	N	O	S	0	0	0
			2526	1592	439	481	14			
1	C	319	Total	C	N	O	S	0	0	0
			2519	1587	439	479	14			
1	E	319	Total	C	N	O	S	0	0	0
			2519	1587	439	479	14			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	323	ARG	-	EXPRESSION TAG	UNP Q207Z6
A	324	GLU	-	EXPRESSION TAG	UNP Q207Z6
A	325	THR	-	EXPRESSION TAG	UNP Q207Z6
A	326	ARG	-	EXPRESSION TAG	UNP Q207Z6
C	323	ARG	-	EXPRESSION TAG	UNP Q207Z6
C	324	GLU	-	EXPRESSION TAG	UNP Q207Z6
C	325	THR	-	EXPRESSION TAG	UNP Q207Z6
C	326	ARG	-	EXPRESSION TAG	UNP Q207Z6
E	323	ARG	-	EXPRESSION TAG	UNP Q207Z6
E	324	GLU	-	EXPRESSION TAG	UNP Q207Z6
E	325	THR	-	EXPRESSION TAG	UNP Q207Z6
E	326	ARG	-	EXPRESSION TAG	UNP Q207Z6

- Molecule 2 is a protein called HEMAGGLUTININ.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	145	Total	C	N	O	S	0	0	0
			1129	697	200	224	8			
2	D	146	Total	C	N	O	S	0	0	0
			1130	698	201	223	8			
2	F	145	Total	C	N	O	S	0	0	0
			1129	697	200	224	8			

- 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	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	C	1	Total	C	N	O	0	0
			14	8	1	5		
3	E	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 4 is a polymer of unknown type called SUGAR (3-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	3	Total	C	N	O	0	0
			46	25	2	19		
4	C	3	Total	C	N	O	0	0
			46	25	2	19		
4	E	3	Total	C	N	O	0	0
			46	25	2	19		

- Molecule 5 is PHOSPHATE ION (three-letter code: PO4) (formula: O_4P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	O	P	0	0
			5	4	1		
5	A	1	Total	O	P	0	0
			5	4	1		
5	C	1	Total	O	P	0	0
			5	4	1		
5	E	1	Total	O	P	0	0
			5	4	1		

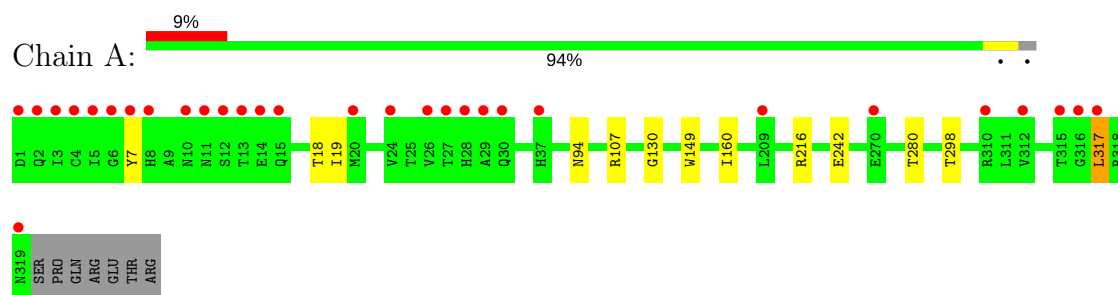
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	266	Total	O	0	0
			266	266		
6	B	36	Total	O	0	0
			36	36		
6	C	243	Total	O	0	0
			243	243		
6	D	35	Total	O	0	0
			35	35		
6	E	218	Total	O	0	0
			218	218		
6	F	30	Total	O	0	0
			30	30		

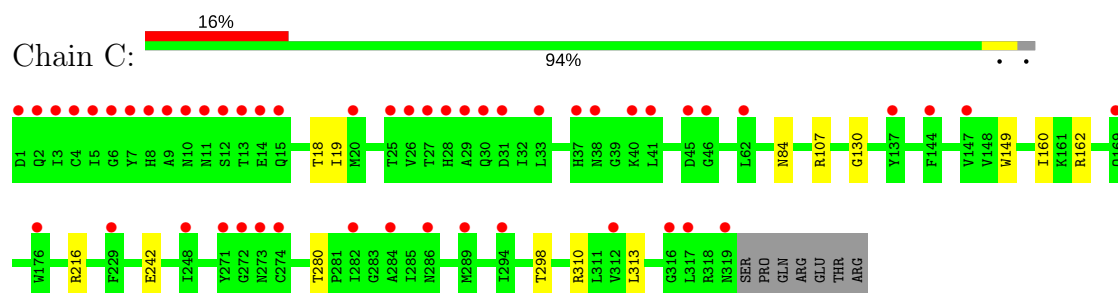
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

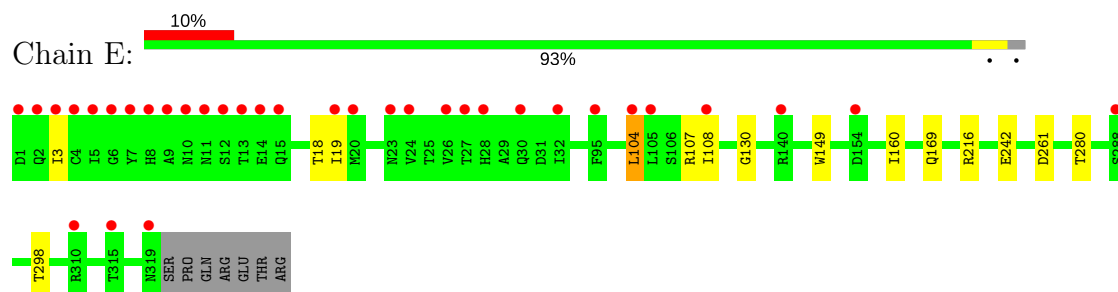
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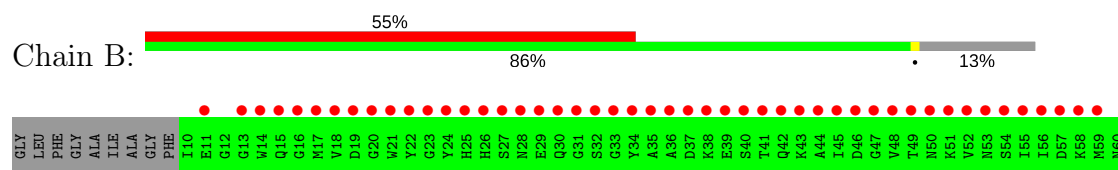
• Molecule 1: HEMAGGLUTININ

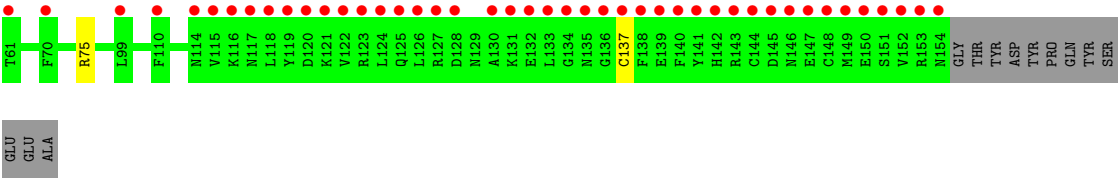


• Molecule 1: HEMAGGLUTININ

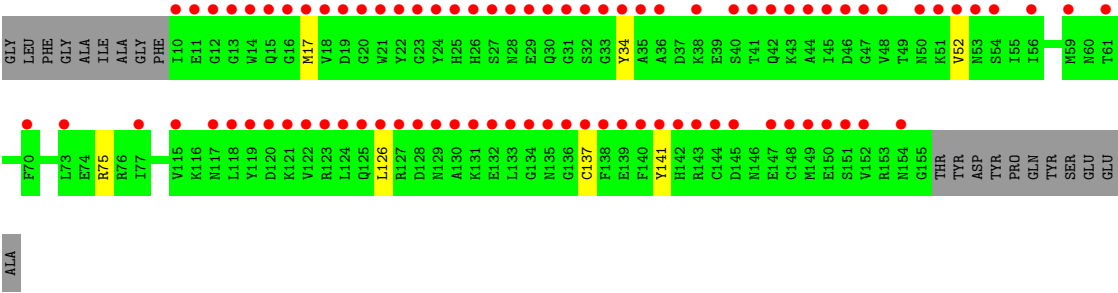
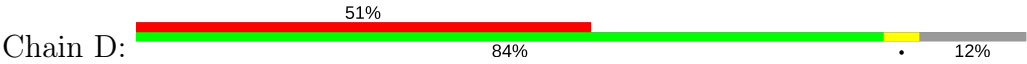


• Molecule 2: HEMAGGLUTININ

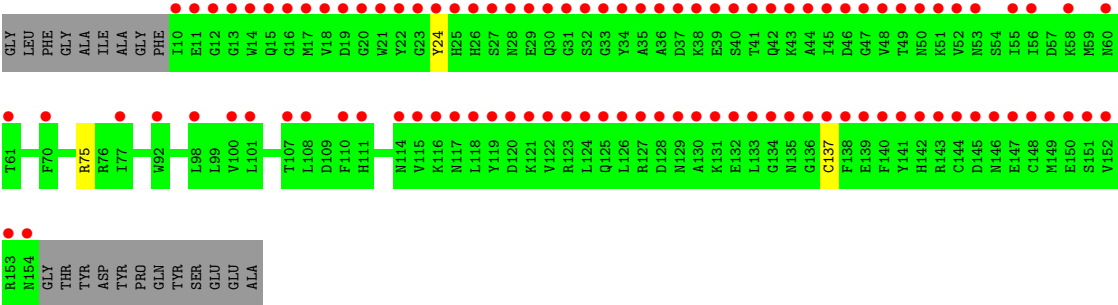
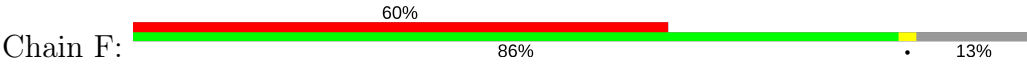




• Molecule 2: HEMAGGLUTININ



• Molecule 2: HEMAGGLUTININ



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	70.86Å 228.29Å 71.92Å 90.00° 113.71° 90.00°	Depositor
Resolution (Å)	62.48 – 2.15 43.13 – 2.16	Depositor EDS
% Data completeness (in resolution range)	98.0 (62.48-2.15) 97.8 (43.13-2.16)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.02 (at 2.16Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.205 , 0.224 0.261 , 0.256	Depositor DCC
R_{free} test set	5488 reflections (5.01%)	DCC
Wilson B-factor (Å ²)	35.4	Xtriage
Anisotropy	0.700	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 54.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.035 for l,-k,h	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	11980	wwPDB-VP
Average B, all atoms (Å ²)	71.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.31% 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, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: PO4, GAL, NAG, SIA

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	0.32	0/2586	0.55	2/3515 (0.1%)
1	C	0.32	0/2579	0.55	2/3506 (0.1%)
1	E	0.31	0/2579	0.55	1/3506 (0.0%)
2	B	0.29	0/1149	0.44	0/1551
2	D	0.30	0/1150	0.44	0/1552
2	F	0.29	0/1149	0.43	0/1551
All	All	0.31	0/11192	0.52	5/15181 (0.0%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	216	ARG	NE-CZ-NH1	5.59	123.10	120.30
1	C	216	ARG	NE-CZ-NH2	-5.57	117.51	120.30
1	C	216	ARG	NE-CZ-NH1	5.46	123.03	120.30
1	E	216	ARG	NE-CZ-NH1	5.42	123.01	120.30
1	A	216	ARG	NE-CZ-NH2	-5.18	117.71	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2526	0	2460	7	0
1	C	2519	0	2447	7	0
1	E	2519	0	2447	6	1
2	B	1129	0	1007	1	0
2	D	1130	0	1008	2	0
2	F	1129	0	1007	1	0
3	A	14	0	13	0	0
3	C	14	0	13	0	0
3	E	14	0	13	0	0
4	A	46	0	40	0	0
4	C	46	0	40	0	0
4	E	46	0	40	0	0
5	A	10	0	0	0	0
5	C	5	0	0	0	0
5	E	5	0	0	0	0
6	A	266	0	0	2	1
6	B	36	0	0	0	0
6	C	243	0	0	3	0
6	D	35	0	0	0	0
6	E	218	0	0	0	0
6	F	30	0	0	0	0
All	All	11980	0	10535	22	1

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 1.

All (22) 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:C:162:ARG:HD2	6:C:2149:HOH:O	2.03	0.58
1:C:84:ASN:ND2	6:C:2069:HOH:O	2.36	0.57
1:A:7:TYR:HB2	1:A:317:LEU:CD2	2.43	0.48
1:A:130:GLY:HA3	1:A:149:TRP:HB3	1.96	0.47
1:E:130:GLY:HA3	1:E:149:TRP:HB3	1.95	0.47
1:C:130:GLY:HA3	1:C:149:TRP:HB3	1.97	0.47
1:E:280:THR:HG22	1:E:298:THR:HG22	1.98	0.46
1:A:280:THR:HG22	1:A:298:THR:HG22	1.98	0.45
1:E:160:ILE:O	1:E:242:GLU:HA	2.17	0.45
1:A:94:ASN:ND2	6:A:2077:HOH:O	2.45	0.45
1:C:313:LEU:HD13	2:D:52:VAL:HG12	1.98	0.45
1:A:160:ILE:O	1:A:242:GLU:HA	2.17	0.44
1:C:160:ILE:O	1:C:242:GLU:HA	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:17:MET:HG3	2:D:34:TYR:HB3	1.99	0.44
1:A:18:THR:HG22	1:A:19:ILE:N	2.33	0.43
1:C:280:THR:HG22	1:C:298:THR:HG22	1.98	0.43
1:E:18:THR:HG22	1:E:19:ILE:N	2.34	0.43
1:E:3:ILE:HD11	2:F:24:TYR:HB3	2.01	0.42
1:A:94:ASN:HB2	6:A:2078:HOH:O	2.20	0.41
1:C:18:THR:HG22	1:C:19:ILE:N	2.34	0.41
2:B:75:ARG:HG3	6:C:2090:HOH:O	2.20	0.41
1:E:104:LEU:HD13	1:E:108:ILE:HD12	2.02	0.41

All (1) 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:E:169:GLN:NE2	6:A:2135:HOH:O[1_655]	2.09	0.11

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	317/326 (97%)	309 (98%)	8 (2%)	0	100	100
1	C	317/326 (97%)	309 (98%)	8 (2%)	0	100	100
1	E	317/326 (97%)	309 (98%)	8 (2%)	0	100	100
2	B	143/166 (86%)	137 (96%)	6 (4%)	0	100	100
2	D	144/166 (87%)	137 (95%)	7 (5%)	0	100	100
2	F	143/166 (86%)	137 (96%)	6 (4%)	0	100	100
All	All	1381/1476 (94%)	1338 (97%)	43 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	284/292 (97%)	282 (99%)	2 (1%)	87	91
1	C	282/292 (97%)	280 (99%)	2 (1%)	87	91
1	E	282/292 (97%)	279 (99%)	3 (1%)	78	82
2	B	113/141 (80%)	112 (99%)	1 (1%)	82	87
2	D	112/141 (79%)	108 (96%)	4 (4%)	40	38
2	F	113/141 (80%)	111 (98%)	2 (2%)	64	68
All	All	1186/1299 (91%)	1172 (99%)	14 (1%)	75	80

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	107	ARG
1	A	317	LEU
2	B	137	CYS
1	C	107	ARG
1	C	310	ARG
2	D	75	ARG
2	D	126	LEU
2	D	137	CYS
2	D	141	TYR
1	E	104	LEU
1	E	107	ARG
1	E	261	ASP
2	F	75	ARG
2	F	137	CYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	2	GLN
1	C	2	GLN
2	D	15	GLN

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Mol	Chain	Res	Type
1	E	2	GLN
1	E	87	ASN
1	E	110	HIS

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 ⓘ

9 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	$\# Z > 2$
4	SIA	A	1321	4	17,20,21	0.52	0	19,28,31	1.00	1 (5%)
4	GAL	A	1322	4	11,11,12	0.43	0	13,15,17	1.00	0
4	NAG	A	1323	4	15,15,15	0.36	0	21,21,21	1.36	4 (19%)
4	SIA	C	1321	4	17,20,21	0.55	0	19,28,31	0.96	1 (5%)
4	GAL	C	1322	4	11,11,12	0.34	0	13,15,17	1.22	2 (15%)
4	NAG	C	1323	4	15,15,15	0.43	0	21,21,21	1.23	2 (9%)
4	SIA	E	1321	4	17,20,21	0.35	0	19,28,31	1.00	2 (10%)
4	GAL	E	1322	4	11,11,12	0.26	0	13,15,17	0.87	1 (7%)
4	NAG	E	1323	4	15,15,15	0.38	0	21,21,21	1.26	2 (9%)

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	SIA	A	1321	4	-	0/14/34/38	0/1/1/1
4	GAL	A	1322	4	-	0/2/19/22	0/1/1/1
4	NAG	A	1323	4	-	0/6/26/26	0/1/1/1
4	SIA	C	1321	4	-	0/14/34/38	0/1/1/1
4	GAL	C	1322	4	-	0/2/19/22	0/1/1/1
4	NAG	C	1323	4	-	0/6/26/26	0/1/1/1
4	SIA	E	1321	4	-	0/14/34/38	0/1/1/1
4	GAL	E	1322	4	-	0/2/19/22	0/1/1/1
4	NAG	E	1323	4	-	0/6/26/26	0/1/1/1

There are no bond length outliers.

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	1323	NAG	C1-C2-N2	-4.25	105.80	110.73
4	A	1323	NAG	C1-C2-N2	-3.61	106.55	110.73
4	E	1323	NAG	C1-C2-N2	-3.52	106.65	110.73
4	A	1321	SIA	C6-C5-N5	-2.57	106.48	111.00
4	A	1323	NAG	C3-C2-N2	-2.54	105.74	110.61
4	E	1323	NAG	C3-C2-N2	-2.49	105.83	110.61
4	C	1321	SIA	C6-C5-N5	-2.42	106.75	111.00
4	E	1321	SIA	C4-C5-N5	-2.35	105.56	110.40
4	E	1321	SIA	C6-C5-N5	-2.18	107.16	111.00
4	E	1322	GAL	O3-C3-C2	-2.10	106.20	110.02
4	C	1323	NAG	C3-C2-N2	-2.09	106.60	110.61
4	C	1322	GAL	C1-O5-C5	2.05	114.99	112.17
4	A	1323	NAG	C1-C2-C3	2.25	113.61	110.54
4	A	1323	NAG	O5-C1-C2	2.52	112.05	109.52
4	C	1322	GAL	C1-C2-C3	2.62	112.97	109.65

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.6 Ligand geometry ⓘ

7 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	$\# Z > 2$
3	NAG	A	1320	1	14,14,15	0.46	0	15,19,21	1.38	2 (13%)
5	PO4	A	1324	-	4,4,4	0.77	0	6,6,6	0.48	0
5	PO4	A	1325	-	4,4,4	0.65	0	6,6,6	0.51	0
3	NAG	C	1320	1	14,14,15	0.53	0	15,19,21	1.37	3 (20%)
5	PO4	C	1324	-	4,4,4	0.88	0	6,6,6	0.45	0
3	NAG	E	1320	1	14,14,15	0.37	0	15,19,21	1.11	1 (6%)
5	PO4	E	1324	-	4,4,4	0.81	0	6,6,6	0.48	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	A	1320	1	-	0/6/23/26	0/1/1/1
5	PO4	A	1324	-	-	0/0/0/0	0/0/0/0
5	PO4	A	1325	-	-	0/0/0/0	0/0/0/0
3	NAG	C	1320	1	-	0/6/23/26	0/1/1/1
5	PO4	C	1324	-	-	0/0/0/0	0/0/0/0
3	NAG	E	1320	1	-	0/6/23/26	0/1/1/1
5	PO4	E	1324	-	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1320	NAG	C3-C4-C5	-2.14	106.44	110.22
3	C	1320	NAG	C3-C4-C5	-2.08	106.55	110.22
3	C	1320	NAG	C1-C2-N2	2.04	113.98	110.49
3	E	1320	NAG	C1-O5-C5	2.55	115.68	112.17
3	A	1320	NAG	C1-O5-C5	3.46	116.93	112.17
3	C	1320	NAG	C1-O5-C5	3.61	117.14	112.17

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

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	319/326 (97%)	1.18	30 (9%) 9 13	23, 38, 124, 285	0
1	C	319/326 (97%)	1.20	51 (15%) 2 3	22, 40, 101, 147	0
1	E	319/326 (97%)	1.04	34 (10%) 7 9	24, 41, 129, 233	0
2	B	145/166 (87%)	5.81	92 (63%) 0 0	25, 172, 220, 245	0
2	D	146/166 (87%)	3.57	85 (58%) 0 0	29, 126, 188, 208	0
2	F	145/166 (87%)	5.54	100 (68%) 0 0	24, 166, 204, 226	0
All	All	1393/1476 (94%)	2.34	392 (28%) 1 1	22, 47, 193, 285	0

All (392) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	23	GLY	33.3
2	D	138	PHE	32.9
2	B	138	PHE	31.9
2	F	138	PHE	24.0
1	A	3	ILE	22.9
2	D	140	PHE	21.4
2	F	122	VAL	21.2
1	A	4	CYS	21.0
2	B	136	GLY	18.1
2	F	22	TYR	17.8
1	E	3	ILE	17.8
2	B	33	GLY	17.7
2	B	32	SER	17.7
2	B	36	ALA	17.0
2	F	140	PHE	16.9
2	B	141	TYR	16.7
2	B	21	TRP	16.6
1	A	5	ILE	16.2
2	B	35	ALA	16.1

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Mol	Chain	Res	Type	RSRZ
2	F	27	SER	15.9
2	B	26	HIS	15.7
2	F	118	LEU	15.0
2	F	30	GLN	14.9
1	A	2	GLN	14.8
2	F	148	CYS	14.6
2	D	125	GLN	14.5
2	B	22	TYR	14.4
2	F	144	CYS	14.2
2	B	122	VAL	13.5
2	B	20	GLY	13.3
2	F	31	GLY	13.2
1	E	4	CYS	13.2
2	D	127	ARG	13.2
2	B	140	PHE	13.2
2	B	152	VAL	13.2
2	B	34	TYR	13.1
2	B	137	CYS	13.0
1	A	6	GLY	13.0
2	F	152	VAL	12.8
2	F	33	GLY	12.7
2	F	26	HIS	12.6
2	F	124	LEU	12.4
2	B	125	GLN	12.3
2	B	146	ASN	12.2
2	F	21	TRP	12.1
2	B	118	LEU	11.9
2	B	50	ASN	11.9
2	F	133	LEU	11.7
2	B	124	LEU	11.7
2	B	17	MET	11.2
2	B	133	LEU	11.1
2	F	24	TYR	11.0
2	B	154	ASN	10.9
2	B	19	ASP	10.8
2	B	126	LEU	10.8
2	B	24	TYR	10.5
2	B	25	HIS	10.3
2	B	148	CYS	10.3
2	F	132	GLU	10.3
2	F	32	SER	10.3
2	D	30	GLN	10.2

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Mol	Chain	Res	Type	RSRZ
2	F	150	GLU	10.1
2	B	30	GLN	10.1
2	B	16	GLY	9.9
2	F	123	ARG	9.9
2	F	126	LEU	9.9
2	F	146	ASN	9.8
2	B	14	TRP	9.8
2	F	35	ALA	9.8
2	D	131	LYS	9.8
2	D	126	LEU	9.8
2	B	28	ASN	9.7
2	F	125	GLN	9.6
2	F	11	GLU	9.2
2	D	33	GLY	9.2
2	B	15	GLN	9.1
2	B	147	GLU	9.1
2	B	27	SER	9.1
2	F	40	SER	9.1
2	F	47	GLY	9.1
1	E	2	GLN	9.0
2	B	128	ASP	8.9
2	B	153	ARG	8.9
1	A	1	ASP	8.9
1	C	6	GLY	8.8
2	F	147	GLU	8.8
2	F	128	ASP	8.8
2	F	25	HIS	8.7
2	F	154	ASN	8.7
2	F	149	MET	8.7
2	F	137	CYS	8.7
2	F	44	ALA	8.7
2	F	20	GLY	8.6
2	F	29	GLU	8.6
2	D	19	ASP	8.6
2	B	145	ASP	8.5
2	F	16	GLY	8.5
2	F	130	ALA	8.5
1	E	7	TYR	8.5
2	F	141	TYR	8.4
2	D	128	ASP	8.4
2	B	131	LYS	8.4
1	A	13	THR	8.3

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Mol	Chain	Res	Type	RSRZ
2	F	36	ALA	8.3
2	B	29	GLU	8.2
2	F	120	ASP	8.1
2	F	136	GLY	8.0
2	F	45	ILE	8.0
2	B	144	CYS	8.0
2	F	50	ASN	8.0
2	B	47	GLY	8.0
2	D	141	TYR	7.8
2	D	136	GLY	7.8
2	D	27	SER	7.8
2	F	134	GLY	7.7
2	B	132	GLU	7.7
2	D	26	HIS	7.7
2	F	142	HIS	7.7
2	F	145	ASP	7.6
2	B	51	LYS	7.6
2	F	18	VAL	7.5
2	F	42	GLN	7.4
2	F	14	TRP	7.4
2	F	28	ASN	7.3
2	B	44	ALA	7.3
2	B	48	VAL	7.2
1	C	11	ASN	7.2
2	B	151	SER	7.1
2	F	23	GLY	7.1
1	E	12	SER	7.1
2	B	121	LYS	7.1
2	D	21	TRP	7.0
2	D	119	TYR	7.0
2	B	149	MET	7.0
2	F	127	ARG	7.0
2	D	129	ASN	6.9
2	F	151	SER	6.9
1	E	1	ASP	6.9
2	F	46	ASP	6.9
2	D	130	ALA	6.9
2	B	54	SER	6.9
2	B	31	GLY	6.8
2	B	143	ARG	6.7
1	E	319	ASN	6.7
2	B	139	GLU	6.7

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Mol	Chain	Res	Type	RSRZ
2	F	34	TYR	6.6
2	D	47	GLY	6.6
2	B	119	TYR	6.6
2	F	139	GLU	6.6
2	B	142	HIS	6.5
2	B	40	SER	6.5
2	D	143	ARG	6.5
1	A	8	HIS	6.3
2	F	43	LYS	6.3
1	A	11	ASN	6.3
2	D	16	GLY	6.3
1	E	27	THR	6.2
2	B	150	GLU	6.2
2	D	134	GLY	6.2
2	D	139	GLU	6.1
1	C	316	GLY	6.0
1	E	11	ASN	6.0
2	F	17	MET	6.0
2	D	52	VAL	6.0
1	E	13	THR	6.0
1	C	3	ILE	5.9
2	B	18	VAL	5.9
2	F	56	ILE	5.9
2	D	154	ASN	5.9
2	B	120	ASP	5.9
2	F	51	LYS	5.9
1	A	30	GLN	5.8
2	D	31	GLY	5.8
2	F	115	VAL	5.8
2	B	117	ASN	5.7
2	F	38	LYS	5.7
1	C	5	ILE	5.7
2	B	38	LYS	5.6
1	C	13	THR	5.6
2	B	56	ILE	5.6
2	D	142	HIS	5.5
2	F	10	ILE	5.5
2	D	42	GLN	5.5
2	B	42	GLN	5.5
2	F	19	ASP	5.5
2	D	28	ASN	5.4
2	D	46	ASP	5.4

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Mol	Chain	Res	Type	RSRZ
2	F	153	ARG	5.4
2	D	122	VAL	5.4
2	B	45	ILE	5.4
1	E	6	GLY	5.4
2	D	118	LEU	5.4
2	F	116	LYS	5.4
2	B	49	THR	5.4
2	F	131	LYS	5.3
2	B	43	LYS	5.3
2	F	117	ASN	5.3
2	D	152	VAL	5.2
2	B	135	ASN	5.2
1	E	5	ILE	5.2
2	D	18	VAL	5.2
2	D	24	TYR	5.2
1	A	317	LEU	5.2
1	E	10	ASN	5.1
2	B	46	ASP	5.1
2	F	110	PHE	5.1
1	E	9	ALA	5.0
2	D	36	ALA	5.0
1	E	26	VAL	5.0
1	C	12	SER	5.0
2	D	132	GLU	5.0
2	D	14	TRP	4.9
2	F	48	VAL	4.9
2	D	38	LYS	4.9
1	A	12	SER	4.9
2	B	55	ILE	4.9
1	A	29	ALA	4.9
2	D	56	ILE	4.8
1	C	4	CYS	4.8
2	B	127	ARG	4.8
1	A	10	ASN	4.7
1	C	28	HIS	4.7
2	F	119	TYR	4.7
2	D	151	SER	4.7
2	D	121	LYS	4.7
2	D	137	CYS	4.7
1	A	316	GLY	4.6
2	F	121	LYS	4.6
1	A	7	TYR	4.6

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Mol	Chain	Res	Type	RSRZ
1	C	7	TYR	4.5
1	E	8	HIS	4.5
1	C	9	ALA	4.5
1	E	28	HIS	4.4
2	D	23	GLY	4.4
1	E	19	ILE	4.3
2	D	15	GLN	4.3
2	F	135	ASN	4.3
1	E	20	MET	4.3
2	D	150	GLU	4.3
2	F	41	THR	4.2
1	C	294	ILE	4.2
2	D	50	ASN	4.2
2	D	20	GLY	4.1
2	B	11	GLU	4.1
1	A	27	THR	4.1
2	B	123	ARG	4.1
2	B	52	VAL	4.1
2	F	61	THR	4.1
2	F	39	GLU	4.0
2	F	52	VAL	4.0
1	E	15	GLN	4.0
1	C	15	GLN	4.0
2	F	15	GLN	3.9
1	C	33	LEU	3.9
1	C	2	GLN	3.9
2	D	51	LYS	3.9
2	F	55	ILE	3.9
2	F	114	ASN	3.8
2	B	134	GLY	3.8
2	D	17	MET	3.8
1	C	8	HIS	3.7
2	D	29	GLU	3.7
2	D	145	ASP	3.7
1	C	10	ASN	3.7
2	F	129	ASN	3.6
2	F	37	ASP	3.6
2	D	123	ARG	3.6
2	B	37	ASP	3.6
2	D	117	ASN	3.5
2	D	35	ALA	3.5
1	E	315	THR	3.5

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Mol	Chain	Res	Type	RSRZ
1	C	312	VAL	3.5
2	F	53	ASN	3.4
2	D	32	SER	3.4
1	C	27	THR	3.4
1	E	140	ARG	3.3
2	D	61	THR	3.3
2	D	43	LYS	3.3
2	D	48	VAL	3.3
1	C	229	PHE	3.3
2	D	133	LEU	3.3
2	D	115	VAL	3.3
2	F	12	GLY	3.2
2	F	107	THR	3.2
2	D	135	ASN	3.1
2	D	124	LEU	3.1
2	F	13	GLY	3.1
1	C	29	ALA	3.1
2	B	59	MET	3.1
1	A	15	GLN	3.1
2	D	59	MET	3.1
1	C	37	HIS	3.0
1	A	319	ASN	3.0
1	C	14	GLU	3.0
2	D	73	LEU	3.0
2	F	143	ARG	3.0
1	C	40	LYS	3.0
2	D	41	THR	3.0
1	C	38	ASN	3.0
2	D	54	SER	3.0
2	D	25	HIS	2.9
1	A	312	VAL	2.9
1	C	20	MET	2.9
2	B	53	ASN	2.9
2	D	148	CYS	2.9
1	C	30	GLN	2.8
2	D	13	GLY	2.8
2	D	77	ILE	2.8
2	D	149	MET	2.8
2	F	111	HIS	2.8
1	E	104	LEU	2.7
2	B	114	ASN	2.7
2	B	39	GLU	2.7

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Mol	Chain	Res	Type	RSRZ
2	D	11	GLU	2.7
1	C	46	GLY	2.7
1	C	282	ILE	2.7
1	A	14	GLU	2.7
1	E	30	GLN	2.7
1	C	144	PHE	2.6
1	C	1	ASP	2.6
1	C	284	ALA	2.6
1	A	310	ARG	2.6
2	F	60	ASN	2.6
2	B	61	THR	2.6
2	D	147	GLU	2.6
1	A	315	THR	2.6
1	E	310	ARG	2.5
2	D	12	GLY	2.5
1	E	14	GLU	2.5
1	A	28	HIS	2.5
1	C	319	ASN	2.5
2	D	70	PHE	2.5
2	F	101	LEU	2.5
1	C	169	GLN	2.4
1	C	147	VAL	2.4
1	C	273	ASN	2.4
1	C	272	GLY	2.4
1	E	108	ILE	2.4
1	A	209	LEU	2.4
1	E	154	ASP	2.4
1	A	26	VAL	2.4
2	B	13	GLY	2.4
2	F	100	VAL	2.4
1	C	271	TYR	2.4
1	C	274	CYS	2.4
2	D	144	CYS	2.4
2	D	44	ALA	2.4
2	D	53	ASN	2.4
1	E	24	VAL	2.3
2	D	22	TYR	2.3
2	D	34	TYR	2.3
2	F	98	LEU	2.3
2	D	45	ILE	2.3
2	D	40	SER	2.3
1	A	24	VAL	2.3

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Mol	Chain	Res	Type	RSRZ
1	C	286	ASN	2.3
1	C	62	LEU	2.3
1	C	25	THR	2.3
1	E	23	ASN	2.3
1	C	45	ASP	2.2
2	F	58	LYS	2.2
2	B	41	THR	2.2
2	B	99	LEU	2.2
2	F	49	THR	2.2
2	B	70	PHE	2.2
1	C	289	MET	2.2
1	C	31	ASP	2.2
1	E	32	ILE	2.2
2	F	108	LEU	2.2
2	B	110	PHE	2.1
1	E	288	SER	2.1
2	F	77	ILE	2.1
1	C	317	LEU	2.1
1	C	176	TRP	2.1
2	F	92	TRP	2.1
1	C	137	TYR	2.1
2	D	10	ILE	2.1
1	A	20	MET	2.1
2	F	70	PHE	2.1
1	A	37	HIS	2.1
1	C	26	VAL	2.1
2	B	115	VAL	2.1
2	B	130	ALA	2.1
1	A	270	GLU	2.1
2	B	116	LYS	2.1
1	E	95	PHE	2.1
2	D	120	ASP	2.0
1	C	41	LEU	2.0
1	C	248	ILE	2.0
2	B	58	LYS	2.0
1	E	105	LEU	2.0
2	B	57	ASP	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 [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
4	GAL	A	1322	11/12	0.83	0.16	2.71	30,36,37,40	0
4	SIA	E	1321	20/21	0.91	0.13	-0.26	34,40,46,46	0
4	SIA	C	1321	20/21	0.90	0.13	-1.60	23,25,29,30	0
4	SIA	A	1321	20/21	0.91	0.12	-1.74	26,31,33,34	0
4	GAL	E	1322	11/12	0.90	0.11	-2.16	47,50,54,56	0
4	NAG	A	1323	15/15	0.87	0.20	-	46,56,64,66	0
4	NAG	C	1323	15/15	0.84	0.26	-	43,56,66,68	0
4	GAL	C	1322	11/12	0.88	0.13	-	28,32,34,36	0
4	NAG	E	1323	15/15	0.88	0.26	-	62,73,76,78	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(Å ²)	Q<0.9
5	PO4	E	1324	5/5	0.85	0.38	10.54	64,71,74,75	0
5	PO4	A	1325	5/5	0.77	0.39	9.88	66,71,78,80	0
5	PO4	C	1324	5/5	0.83	0.34	6.55	53,65,67,67	0
5	PO4	A	1324	5/5	0.93	0.22	1.02	63,63,67,68	0
3	NAG	C	1320	14/15	0.83	0.16	0.12	37,40,47,49	0
3	NAG	E	1320	14/15	0.88	0.15	-0.11	35,37,46,50	0
3	NAG	A	1320	14/15	0.89	0.15	-	36,39,47,49	0

6.5 Other polymers [i](#)

There are no such residues in this entry.