



wwPDB X-ray Structure Validation Summary Report ⓘ

Nov 6, 2023 – 07:17 PM EST

PDB ID : 4LAJ
Title : Crystal structure of HIV-1 YU2 envelope gp120 glycoprotein in complex with CD4-mimetic miniprotein, M48U1, and llama single-domain, broadly neutralizing, co-receptor binding site antibody, JM4
Authors : Acharya, P.; Luongo, T.S.; Kwong, P.D.
Deposited on : 2013-06-20
Resolution : 2.14 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.36
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36

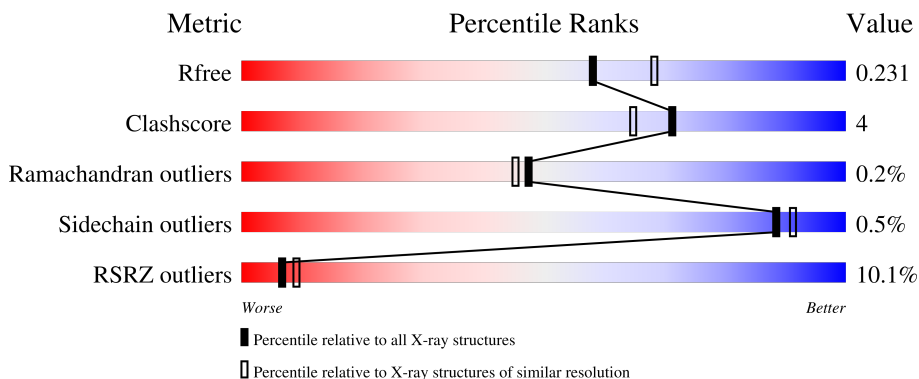
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.14 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	2523 (2.16-2.12)
Clashscore	141614	2653 (2.16-2.12)
Ramachandran outliers	138981	2618 (2.16-2.12)
Sidechain outliers	138945	2617 (2.16-2.12)
RSRZ outliers	127900	2485 (2.16-2.12)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. 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	376	 10% 86% 5% 9%
1	B	376	 8% 83% 7% 9%
1	F	376	 8% 85% 11% 9%
1	J	376	 9% 85% 5% 10%
2	C	28	 4% 82% 14% 9%

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Mol	Chain	Length	Quality of chain
2	D	28	
2	G	28	
2	K	28	
3	H	129	
3	I	129	
3	L	129	
3	M	129	

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	NAG	F	503	-	-	-	X
5	EDO	J	510	-	-	-	X

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 16384 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 HIV-1 YU2 gp120 envelope glycoprotein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	J	337	2660	1669	465	506	20	0	6	0
1	F	334	2633	1654	461	498	20	0	5	0
1	A	343	2702	1693	474	515	20	0	4	0
1	B	341	2680	1681	470	509	20	0	3	0

- Molecule 2 is a protein called CD4-mimetic miniprotein M48U1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	K	28	214	136	39	33	6	0	1	1
2	D	28	209	133	38	32	6	0	0	1
2	G	28	209	133	38	32	6	0	0	1
2	C	28	209	133	38	32	6	0	0	1

- Molecule 3 is a protein called Llama single domain antibody, JM4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	H	124	944	589	156	193	6	0	0	0
3	M	123	939	586	155	192	6	0	0	0
3	L	123	939	586	155	192	6	0	0	0
3	I	125	949	592	157	194	6	0	0	0

- Molecule 4 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



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

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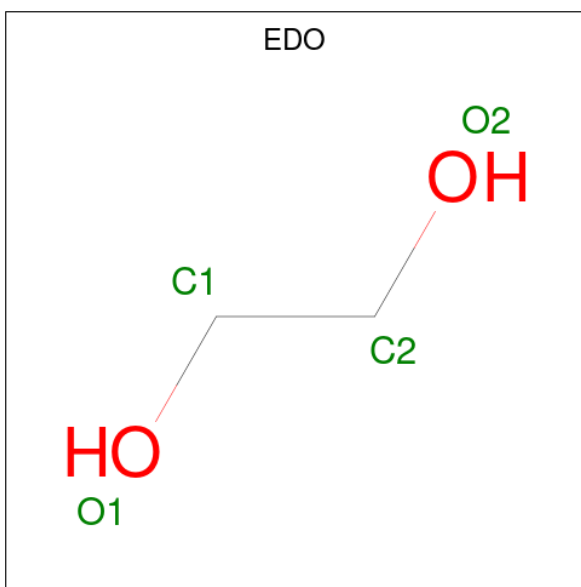
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	F	1	Total	C	N	O	0	0
			14	8	1	5		
4	F	1	Total	C	N	O	0	0
			14	8	1	5		
4	F	1	Total	C	N	O	0	0
			14	8	1	5		
4	F	1	Total	C	N	O	0	0
			14	8	1	5		
4	F	1	Total	C	N	O	0	0
			14	8	1	5		
4	A	1	Total	C	N	O	0	0
			14	8	1	5		
4	A	1	Total	C	N	O	0	0
			14	8	1	5		
4	A	1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

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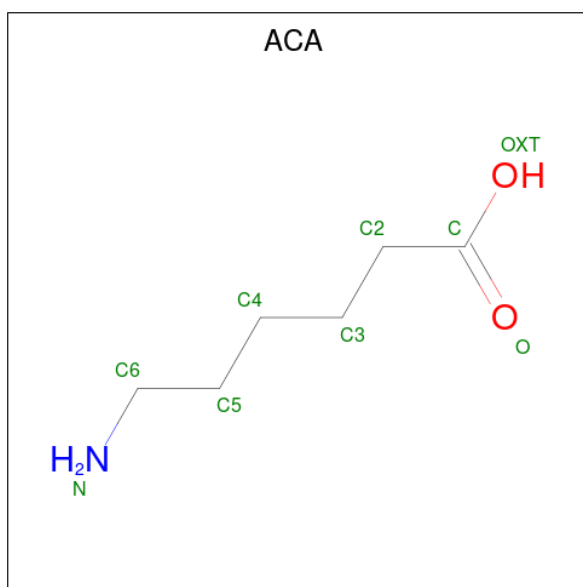
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	B	1	Total	C	N	O	0	0
			14	8	1	5		
4	B	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 5 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	J	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	B	1	Total	C	O	0	0
			4	2	2		
5	B	1	Total	C	O	0	0
			4	2	2		
5	B	1	Total	C	O	0	0
			4	2	2		

- Molecule 6 is 6-AMINOHEXANOIC ACID (three-letter code: ACA) (formula: C₆H₁₃NO₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	B	1	Total	C	N	O	0	0
			9	6	1	2		

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	J	73	Total	O	0	0
			73	73		
7	F	111	Total	O	0	0
			111	111		
7	A	114	Total	O	0	0
			114	114		
7	B	139	Total	O	0	0
			139	139		
7	K	4	Total	O	0	0
			4	4		
7	D	10	Total	O	0	0
			10	10		
7	G	3	Total	O	0	0
			3	3		
7	C	10	Total	O	0	0
			10	10		
7	H	37	Total	O	0	0
			37	37		
7	M	9	Total	O	0	0
			9	9		
7	L	25	Total	O	0	0
			25	25		

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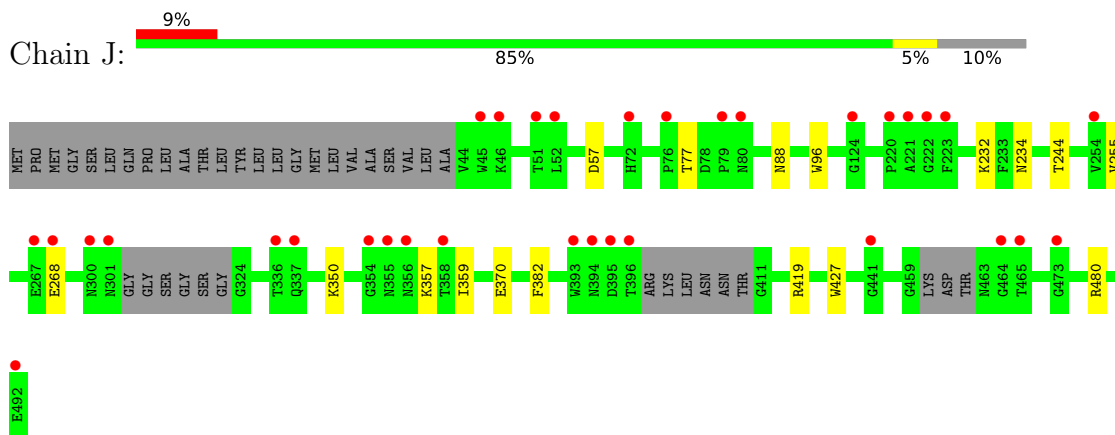
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	I	35	Total	O	0	0
			35	35		

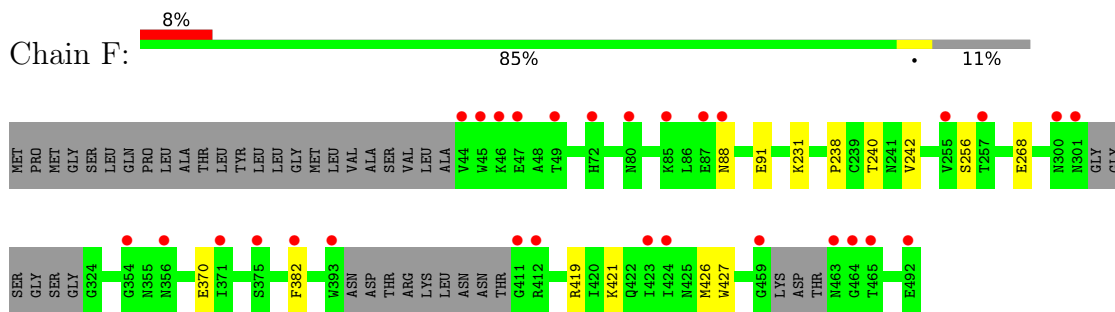
3 Residue-property plots [i](#)

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

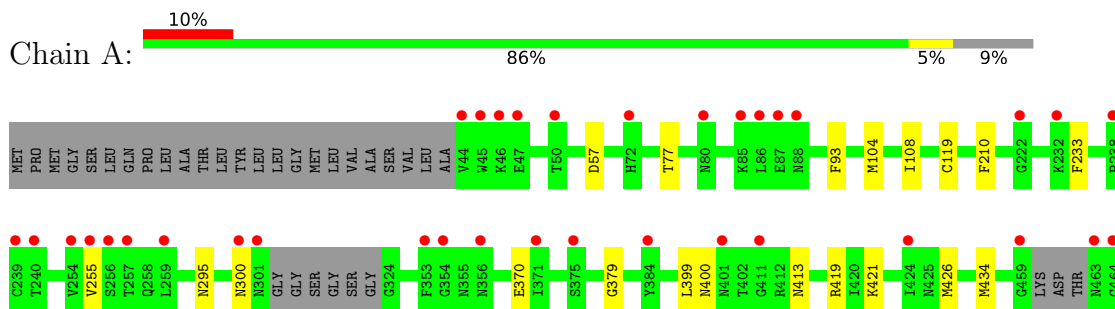
- Molecule 1: HIV-1 YU2 gp120 envelope glycoprotein



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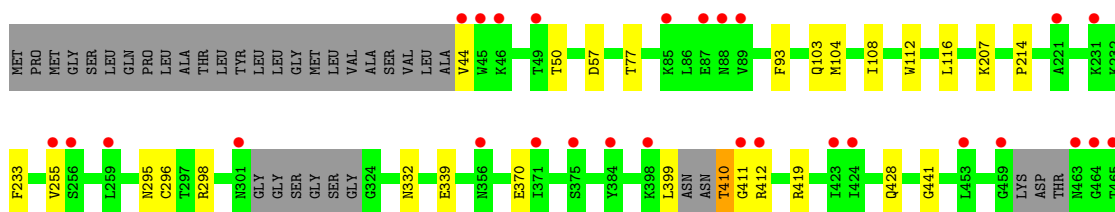
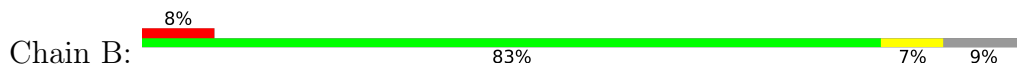


- Molecule 1: HIV-1 YU2 gp120 envelope glycoprotein

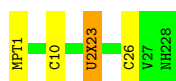
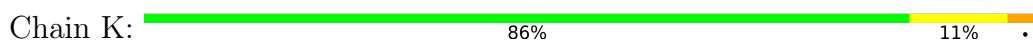




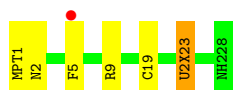
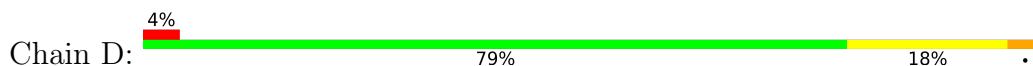
- Molecule 1: HIV-1 YU2 gp120 envelope glycoprotein



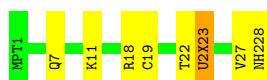
- Molecule 2: CD4-mimetic miniprotein M48U1



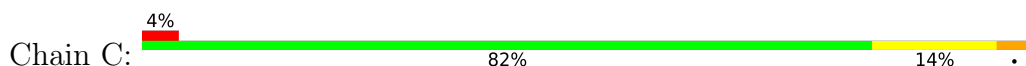
- Molecule 2: CD4-mimetic miniprotein M48U1



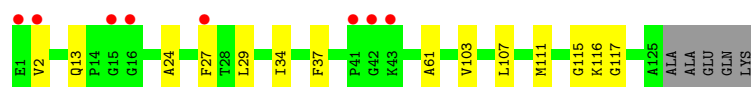
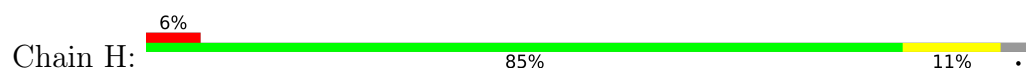
- Molecule 2: CD4-mimetic miniprotein M48U1



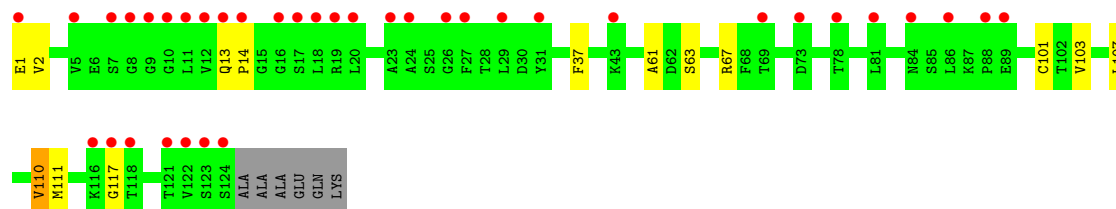
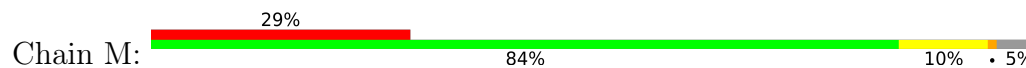
- Molecule 2: CD4-mimetic miniprotein M48U1



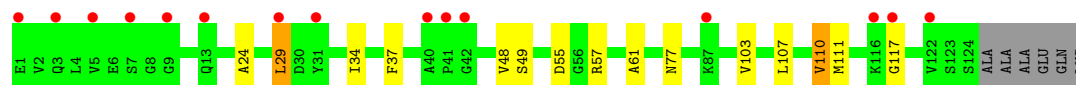
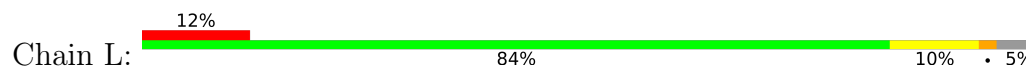
- Molecule 3: Llama single domain antibody, JM4



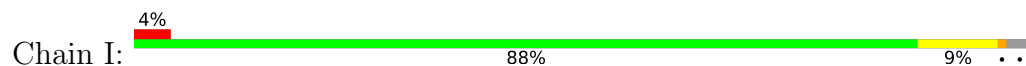
- Molecule 3: Llama single domain antibody, JM4



- Molecule 3: Llama single domain antibody, JM4



- Molecule 3: Llama single domain antibody, JM4



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	96.77Å 86.22Å 144.85Å 90.00° 102.88° 90.00°	Depositor
Resolution (Å)	38.80 – 2.14 38.80 – 2.13	Depositor EDS
% Data completeness (in resolution range)	76.7 (38.80-2.14) 71.7 (38.80-2.13)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.84 (at 2.14Å)	Xtrriage
Refinement program	PHENIX (phenix.refine: 1.8.2_1309)	Depositor
R, R_{free}	0.185 , 0.227 0.192 , 0.231	Depositor DCC
R_{free} test set	4979 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	29.5	Xtrriage
Anisotropy	0.350	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 43.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	16384	wwPDB-VP
Average B, all atoms (Å ²)	60.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.68% 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: DPR, U2X, MPT, NAG, NH2, EDO, ACA

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	0/2783	0.50	0/3775
1	B	0.28	0/2754	0.49	0/3734
1	F	0.27	0/2706	0.47	0/3670
1	J	0.28	0/2746	0.49	0/3724
2	C	0.53	1/176 (0.6%)	0.71	1/231 (0.4%)
2	D	0.55	1/176 (0.6%)	0.59	0/231
2	G	0.52	1/176 (0.6%)	0.46	0/231
2	K	0.35	0/184	0.58	0/242
3	H	0.30	0/962	0.56	0/1304
3	I	0.28	0/967	0.53	0/1311
3	L	0.30	0/957	0.56	0/1297
3	M	0.27	0/957	0.51	0/1297
All	All	0.29	3/15544 (0.0%)	0.51	1/21047 (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
2	D	0	1
2	K	0	2
3	H	0	1
3	I	0	1
3	L	0	1
3	M	0	1
All	All	0	7

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	19	CYS	C-N	6.18	1.48	1.34
2	G	19	CYS	C-N	5.92	1.47	1.34
2	C	19	CYS	C-N	5.88	1.47	1.34

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	19	CYS	C-N-CA	5.82	136.24	121.70

There are no chirality outliers.

5 of 7 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	D	1	MPT	Mainchain
3	H	117	GLY	Peptide
2	K	1	MPT	Peptide
3	L	117	GLY	Peptide
3	M	117	GLY	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	2702	0	2627	17	0
1	B	2680	0	2605	22	0
1	F	2633	0	2557	14	0
1	J	2660	0	2584	13	0
2	C	209	0	212	5	0
2	D	209	0	212	3	0
2	G	209	0	212	10	0
2	K	214	0	220	7	0
3	H	944	0	901	8	0
3	I	949	0	906	10	0
3	L	939	0	896	9	0
3	M	939	0	896	8	0
4	A	126	0	117	2	0
4	B	112	0	104	2	0
4	F	126	0	117	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	J	126	0	117	4	0
5	A	12	0	18	4	0
5	B	12	0	18	3	0
5	J	4	0	6	1	0
6	B	9	0	6	1	0
7	A	114	0	0	0	0
7	B	139	0	0	2	0
7	C	10	0	0	1	0
7	D	10	0	0	0	0
7	F	111	0	0	1	0
7	G	3	0	0	0	0
7	H	37	0	0	0	0
7	I	35	0	0	3	0
7	J	73	0	0	0	0
7	K	4	0	0	0	0
7	L	25	0	0	0	0
7	M	9	0	0	0	0
All	All	16384	0	15331	109	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 4.

The worst 5 of 109 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)
2:K:10:CYS:CB	2:K:26:CYS:SG	2.24	1.25
2:K:10:CYS:SG	2:K:26:CYS:SG	1.05	0.96
2:K:10:CYS:SG	2:K:26:CYS:CB	2.63	0.86
1:B:296:CYS:SG	7:B:717:HOH:O	2.33	0.85
1:J:370:GLU:HG2	2:K:23:U2X:H71	1.59	0.85

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	342/376 (91%)	326 (95%)	15 (4%)	1 (0%)	41	36
1	B	336/376 (89%)	323 (96%)	13 (4%)	0	100	100
1	F	330/376 (88%)	317 (96%)	12 (4%)	1 (0%)	41	36
1	J	336/376 (89%)	322 (96%)	14 (4%)	0	100	100
2	C	24/28 (86%)	24 (100%)	0	0	100	100
2	D	24/28 (86%)	24 (100%)	0	0	100	100
2	G	24/28 (86%)	24 (100%)	0	0	100	100
2	K	25/28 (89%)	23 (92%)	2 (8%)	0	100	100
3	H	122/129 (95%)	118 (97%)	4 (3%)	0	100	100
3	I	123/129 (95%)	118 (96%)	4 (3%)	1 (1%)	19	11
3	L	121/129 (94%)	116 (96%)	4 (3%)	1 (1%)	19	11
3	M	121/129 (94%)	117 (97%)	4 (3%)	0	100	100
All	All	1928/2132 (90%)	1852 (96%)	72 (4%)	4 (0%)	47	45

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	F	88	ASN
1	A	413	ASN
3	I	24	ALA
3	L	29	LEU

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	307/328 (94%)	307 (100%)	0	100	100
1	B	303/328 (92%)	302 (100%)	1 (0%)	92	94
1	F	297/328 (90%)	297 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	J	302/328 (92%)	301 (100%)	1 (0%)	92	94
2	C	20/20 (100%)	20 (100%)	0	100	100
2	D	20/20 (100%)	19 (95%)	1 (5%)	24	20
2	G	20/20 (100%)	20 (100%)	0	100	100
2	K	21/20 (105%)	21 (100%)	0	100	100
3	H	103/106 (97%)	102 (99%)	1 (1%)	76	79
3	I	103/106 (97%)	101 (98%)	2 (2%)	57	59
3	L	103/106 (97%)	102 (99%)	1 (1%)	76	79
3	M	103/106 (97%)	102 (99%)	1 (1%)	76	79
All	All	1702/1816 (94%)	1694 (100%)	8 (0%)	88	91

5 of 8 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	I	124	SER
3	I	48	VAL
3	M	110	VAL
3	H	13	GLN
3	L	110	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
2	C	2	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

8 non-standard protein/DNA/RNA residues 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	U2X	C	23	2	19,20,21	2.35	5 (26%)	22,25,27	1.63	4 (18%)
2	U2X	G	23	2	19,20,21	2.38	4 (21%)	22,25,27	1.64	5 (22%)
2	U2X	K	23	2	19,20,21	2.41	4 (21%)	22,25,27	1.52	5 (22%)
2	U2X	D	23	2	19,20,21	2.32	4 (21%)	22,25,27	1.59	5 (22%)

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	U2X	C	23	2	-	3/10/19/21	0/2/2/2
2	U2X	G	23	2	-	3/10/19/21	0/2/2/2
2	U2X	K	23	2	-	5/10/19/21	0/2/2/2
2	U2X	D	23	2	-	3/10/19/21	0/2/2/2

The worst 5 of 17 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	K	23	U2X	CE2-CD2	7.62	1.52	1.38
2	G	23	U2X	CE2-CD2	7.48	1.52	1.38
2	C	23	U2X	CE2-CD2	7.38	1.52	1.38
2	D	23	U2X	CE2-CD2	7.26	1.52	1.38
2	K	23	U2X	CE1-CD1	4.60	1.47	1.38

The worst 5 of 19 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	23	U2X	OH-C7-C3	3.50	117.35	108.21
2	K	23	U2X	OH-C7-C3	3.40	117.09	108.21
2	C	23	U2X	OH-C7-C3	3.38	117.04	108.21
2	D	23	U2X	C2-C3-C7	3.23	119.39	111.28
2	G	23	U2X	C2-C3-C7	3.22	119.37	111.28

There are no chirality outliers.

5 of 14 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	K	23	U2X	C2-C3-C7-OH
2	D	23	U2X	C2-C3-C7-OH
2	G	23	U2X	C2-C3-C7-OH
2	G	23	U2X	C4-C3-C7-OH
2	C	23	U2X	C2-C3-C7-OH

There are no ring outliers.

4 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	23	U2X	3	0
2	G	23	U2X	4	0
2	K	23	U2X	3	0
2	D	23	U2X	2	0

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

43 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	505	1	14,14,15	0.19	0	17,19,21	0.46	0
4	NAG	B	506	1	14,14,15	0.55	0	17,19,21	0.39	0
4	NAG	B	502	-	14,14,15	0.34	0	17,19,21	0.62	0
4	NAG	B	505	1	14,14,15	0.30	0	17,19,21	0.43	0
4	NAG	J	507	1	14,14,15	0.30	0	17,19,21	0.40	0
4	NAG	A	506	1	14,14,15	0.39	0	17,19,21	0.47	0
4	NAG	F	502	1	14,14,15	0.29	0	17,19,21	0.60	0
4	NAG	J	508	1	14,14,15	0.30	0	17,19,21	0.39	0
4	NAG	B	507	1	14,14,15	0.48	0	17,19,21	0.41	0
5	EDO	A	510	-	3,3,3	0.42	0	2,2,2	0.28	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	NAG	F	504	1	14,14,15	0.28	0	17,19,21	0.65	0
4	NAG	F	501	1	14,14,15	0.25	0	17,19,21	0.54	0
4	NAG	B	508	1	14,14,15	0.41	0	17,19,21	0.71	0
4	NAG	J	501	1	14,14,15	0.38	0	17,19,21	0.45	0
5	EDO	J	510	-	3,3,3	0.46	0	2,2,2	0.33	0
4	NAG	F	509	-	14,14,15	0.95	1 (7%)	17,19,21	0.65	0
4	NAG	A	508	1	14,14,15	0.48	0	17,19,21	0.93	1 (5%)
5	EDO	A	512	-	3,3,3	0.39	0	2,2,2	0.21	0
6	ACA	B	512	-	8,8,8	0.69	0	8,8,8	0.95	0
5	EDO	A	511	-	3,3,3	0.32	0	2,2,2	0.53	0
4	NAG	J	504	1	14,14,15	0.19	0	17,19,21	0.53	0
4	NAG	F	506	1	14,14,15	0.48	0	17,19,21	0.74	1 (5%)
4	NAG	A	503	1	14,14,15	0.22	0	17,19,21	0.52	0
4	NAG	F	507	1	14,14,15	0.19	0	17,19,21	0.50	0
4	NAG	J	509	1	14,14,15	0.24	0	17,19,21	0.46	0
4	NAG	J	502	1	14,14,15	0.13	0	17,19,21	0.61	0
4	NAG	J	503	1	14,14,15	0.24	0	17,19,21	1.02	1 (5%)
5	EDO	B	511	-	3,3,3	0.46	0	2,2,2	0.33	0
4	NAG	A	504	1	14,14,15	0.26	0	17,19,21	0.63	0
4	NAG	F	508	1	14,14,15	0.16	0	17,19,21	0.44	0
4	NAG	F	505	1	14,14,15	0.34	0	17,19,21	0.45	0
4	NAG	B	504	1	14,14,15	0.20	0	17,19,21	0.58	0
4	NAG	J	505	-	14,14,15	0.25	0	17,19,21	0.34	0
4	NAG	B	503	1	14,14,15	0.36	0	17,19,21	0.52	0
4	NAG	A	509	1	14,14,15	0.26	0	17,19,21	0.36	0
4	NAG	F	503	1	14,14,15	0.33	0	17,19,21	0.50	0
5	EDO	B	510	-	3,3,3	0.46	0	2,2,2	0.33	0
5	EDO	B	509	-	3,3,3	0.46	0	2,2,2	0.33	0
4	NAG	A	501	1	14,14,15	0.14	0	17,19,21	0.56	0
4	NAG	J	506	1	14,14,15	0.24	0	17,19,21	0.41	0
4	NAG	B	501	1	14,14,15	0.18	0	17,19,21	0.65	0
4	NAG	A	502	-	14,14,15	0.30	0	17,19,21	0.55	0
4	NAG	A	507	1	14,14,15	0.29	0	17,19,21	0.41	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
4	NAG	A	505	1	-	0/6/23/26	0/1/1/1
4	NAG	B	506	1	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	B	502	-	-	2/6/23/26	0/1/1/1
4	NAG	B	505	1	-	1/6/23/26	0/1/1/1
4	NAG	J	507	1	-	1/6/23/26	0/1/1/1
4	NAG	A	506	1	-	1/6/23/26	0/1/1/1
4	NAG	F	502	1	-	2/6/23/26	0/1/1/1
4	NAG	J	508	1	-	2/6/23/26	0/1/1/1
4	NAG	B	507	1	-	0/6/23/26	0/1/1/1
5	EDO	A	510	-	-	1/1/1/1	-
4	NAG	F	504	1	-	0/6/23/26	0/1/1/1
4	NAG	F	501	1	-	2/6/23/26	0/1/1/1
4	NAG	B	508	1	-	2/6/23/26	0/1/1/1
4	NAG	J	501	1	-	2/6/23/26	0/1/1/1
5	EDO	J	510	-	-	0/1/1/1	-
4	NAG	F	509	-	-	0/6/23/26	0/1/1/1
4	NAG	A	508	1	-	4/6/23/26	0/1/1/1
5	EDO	A	512	-	-	1/1/1/1	-
6	ACA	B	512	-	-	0/6/6/6	-
5	EDO	A	511	-	-	1/1/1/1	-
4	NAG	J	504	1	-	0/6/23/26	0/1/1/1
4	NAG	F	506	1	-	0/6/23/26	0/1/1/1
4	NAG	A	503	1	-	3/6/23/26	0/1/1/1
4	NAG	F	507	1	-	0/6/23/26	0/1/1/1
4	NAG	J	509	1	-	2/6/23/26	0/1/1/1
4	NAG	J	502	1	-	2/6/23/26	0/1/1/1
4	NAG	J	503	1	-	5/6/23/26	0/1/1/1
5	EDO	B	511	-	-	0/1/1/1	-
4	NAG	A	504	1	-	0/6/23/26	0/1/1/1
4	NAG	F	508	1	-	2/6/23/26	0/1/1/1
4	NAG	F	505	1	-	0/6/23/26	0/1/1/1
4	NAG	B	504	1	-	0/6/23/26	0/1/1/1
4	NAG	J	505	-	-	2/6/23/26	0/1/1/1
4	NAG	B	503	1	-	2/6/23/26	0/1/1/1
4	NAG	A	509	1	-	2/6/23/26	0/1/1/1
4	NAG	F	503	1	-	1/6/23/26	0/1/1/1
5	EDO	B	510	-	-	0/1/1/1	-
5	EDO	B	509	-	-	0/1/1/1	-
4	NAG	A	501	1	-	2/6/23/26	0/1/1/1
4	NAG	J	506	1	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	B	501	1	-	2/6/23/26	0/1/1/1
4	NAG	A	502	-	-	3/6/23/26	0/1/1/1
4	NAG	A	507	1	-	0/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	F	509	NAG	O5-C1	-3.03	1.38	1.43

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	508	NAG	C1-O5-C5	2.87	116.08	112.19
4	J	503	NAG	C2-N2-C7	2.83	126.94	122.90
4	F	506	NAG	C1-O5-C5	2.78	115.96	112.19

There are no chirality outliers.

5 of 52 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	508	NAG	C4-C5-C6-O6
4	A	508	NAG	O5-C5-C6-O6
4	J	501	NAG	C4-C5-C6-O6
4	A	509	NAG	O5-C5-C6-O6
4	B	508	NAG	O5-C5-C6-O6

There are no ring outliers.

12 monomers are involved in 20 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	502	NAG	2	0
5	A	510	EDO	1	0
5	J	510	EDO	1	0
4	F	509	NAG	3	0
5	A	512	EDO	3	0
6	B	512	ACA	1	0
4	J	502	NAG	1	0
4	J	503	NAG	2	0
4	J	505	NAG	1	0
5	B	510	EDO	2	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	509	EDO	1	0
4	A	502	NAG	2	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	343/376 (91%)	0.70	38 (11%) 5 6	33, 50, 99, 160	0
1	B	341/376 (90%)	0.54	29 (8%) 10 14	29, 45, 88, 140	0
1	F	334/376 (88%)	0.56	29 (8%) 10 13	36, 53, 97, 135	0
1	J	337/376 (89%)	0.68	33 (9%) 7 10	36, 60, 102, 133	0
2	C	24/28 (85%)	0.40	1 (4%) 36 44	21, 49, 70, 72	0
2	D	24/28 (85%)	0.27	1 (4%) 36 44	21, 51, 70, 76	0
2	G	24/28 (85%)	0.43	0 100 100	21, 66, 81, 88	0
2	K	24/28 (85%)	0.32	0 100 100	21, 60, 79, 86	0
3	H	124/129 (96%)	0.39	8 (6%) 18 23	35, 53, 83, 131	0
3	I	125/129 (96%)	0.51	5 (4%) 38 46	34, 54, 83, 117	0
3	L	123/129 (95%)	0.66	15 (12%) 4 5	38, 63, 87, 113	0
3	M	123/129 (95%)	1.54	37 (30%) 0 0	47, 92, 134, 156	0
All	All	1946/2132 (91%)	0.65	196 (10%) 7 9	21, 55, 103, 160	0

The worst 5 of 196 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	44	VAL	15.4
3	M	123	SER	8.3
1	B	45	TRP	7.8
3	M	124	SER	7.2
1	A	44	VAL	7.2

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

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,

median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	DPR	D	21	7/8	0.92	0.11	19,19,24,25	0
2	DPR	C	21	7/8	0.92	0.14	19,19,24,25	0
2	DPR	G	21	7/8	0.93	0.14	19,19,24,25	0
2	DPR	K	21	7/8	0.93	0.14	19,19,24,25	0
2	U2X	K	23	19/20	0.94	0.22	35,46,57,59	0
2	U2X	C	23	19/20	0.95	0.20	25,38,46,50	0
2	U2X	G	23	19/20	0.96	0.18	35,41,50,58	0
2	U2X	D	23	19/20	0.96	0.23	32,39,47,47	0

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

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. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	NAG	F	503	14/15	0.61	0.44	115,130,142,143	0
4	NAG	B	502	14/15	0.61	0.32	47,79,94,94	0
5	EDO	J	510	4/4	0.68	0.43	78,79,79,83	0
5	EDO	B	510	4/4	0.68	0.36	80,81,81,84	0
4	NAG	J	505	14/15	0.71	0.29	98,109,122,128	0
4	NAG	A	508	14/15	0.73	0.23	72,81,90,92	0
5	EDO	B	509	4/4	0.78	0.29	58,59,61,61	0
4	NAG	A	502	14/15	0.79	0.29	57,75,97,100	0
4	NAG	F	506	14/15	0.79	0.32	84,89,93,94	0
4	NAG	J	509	14/15	0.80	0.33	84,93,103,104	0
4	NAG	J	503	14/15	0.81	0.38	97,104,107,107	0
4	NAG	F	508	14/15	0.81	0.20	57,74,84,87	0
4	NAG	J	501	14/15	0.82	0.27	90,98,103,104	0
4	NAG	J	502	14/15	0.82	0.19	91,95,98,103	0
6	ACA	B	512	9/9	0.84	0.31	52,59,69,70	0
5	EDO	A	510	4/4	0.85	0.36	62,63,64,66	0
4	NAG	B	503	14/15	0.86	0.16	73,84,89,91	0
4	NAG	A	501	14/15	0.86	0.27	66,75,80,86	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	NAG	F	509	14/15	0.86	0.20	56,67,90,95	0
4	NAG	A	509	14/15	0.88	0.22	50,62,82,84	0
4	NAG	B	501	14/15	0.88	0.20	61,69,76,80	0
5	EDO	A	512	4/4	0.88	0.21	55,55,57,59	0
4	NAG	A	506	14/15	0.88	0.22	52,62,72,76	0
4	NAG	A	503	14/15	0.88	0.21	82,85,94,94	0
4	NAG	B	508	14/15	0.88	0.16	50,72,78,83	0
4	NAG	F	505	14/15	0.89	0.24	67,71,80,83	0
4	NAG	J	507	14/15	0.90	0.29	81,85,96,98	0
4	NAG	B	506	14/15	0.90	0.16	49,60,68,69	0
5	EDO	B	511	4/4	0.90	0.16	53,59,63,70	0
4	NAG	F	507	14/15	0.90	0.14	52,58,65,68	0
4	NAG	A	507	14/15	0.91	0.14	36,45,52,52	0
4	NAG	A	505	14/15	0.91	0.18	62,67,76,79	0
4	NAG	B	505	14/15	0.91	0.12	47,52,70,70	0
5	EDO	A	511	4/4	0.92	0.16	49,50,50,52	0
4	NAG	J	506	14/15	0.93	0.26	90,94,104,106	0
4	NAG	J	508	14/15	0.94	0.08	45,50,59,61	0
4	NAG	B	504	14/15	0.95	0.13	29,40,44,45	0
4	NAG	F	502	14/15	0.95	0.12	52,62,71,74	0
4	NAG	F	501	14/15	0.95	0.18	62,70,79,87	0
4	NAG	B	507	14/15	0.96	0.11	34,44,48,50	0
4	NAG	J	504	14/15	0.96	0.13	52,55,64,64	0
4	NAG	A	504	14/15	0.97	0.12	30,40,46,47	0
4	NAG	F	504	14/15	0.97	0.15	34,43,55,58	0

6.5 Other polymers [\(i\)](#)

There are no such residues in this entry.