



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 28, 2020 – 04:29 PM EDT

PDB ID : 6UTB
Title : CRYSTAL STRUCTURE OF UNLIGANDED HIV-1 LM/HT CLADE A/E
CRF01 GP120 CORE
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Deposited on : 2019-10-29
Resolution : 2.50 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

<https://www.wwpdb.org/validation/2017/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	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.14.6
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)	:	2.14.6

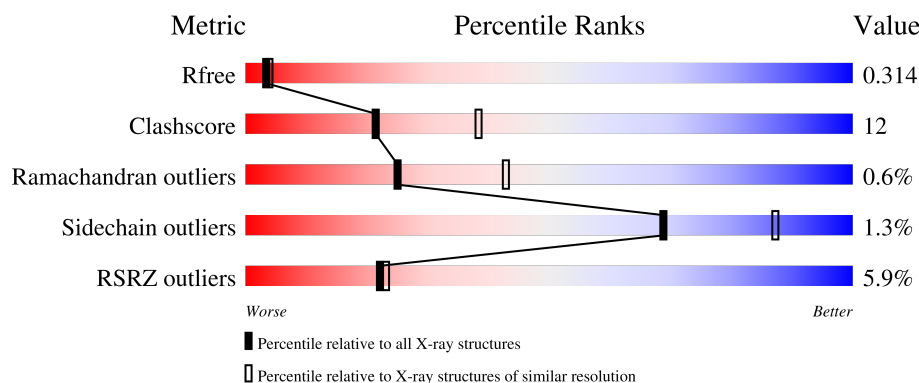
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION


The reported resolution of this entry is 2.50 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	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

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 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	355	

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	504	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	NAG	A	510	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 2875 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 LM/HT Clade A/E CRF01 gp120 core.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	341	Total	C	N	O	S	0	0	0
			2674	1680	461	510	23			

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	42	VAL	-	expression tag	UNP A0A0M3KKW9
A	43	PRO	-	expression tag	UNP A0A0M3KKW9
A	61	TYR	HIS	engineered mutation	UNP A0A0M3KKW9
A	105	HIS	GLN	engineered mutation	UNP A0A0M3KKW9
A	108	ILE	VAL	engineered mutation	UNP A0A0M3KKW9
A	375	THR	HIS	engineered mutation	UNP A0A0M3KKW9
A	474	ASP	ASN	engineered mutation	UNP A0A0M3KKW9
A	475	MET	ILE	engineered mutation	UNP A0A0M3KKW9
A	476	ARG	LYS	engineered mutation	UNP A0A0M3KKW9

- Molecule 2 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: C₈H₁₅NO₆).



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

- Molecule 3 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (three-letter code: EPE) (formula: $C_8H_{18}N_2O_4S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			15	8	2	4	1		

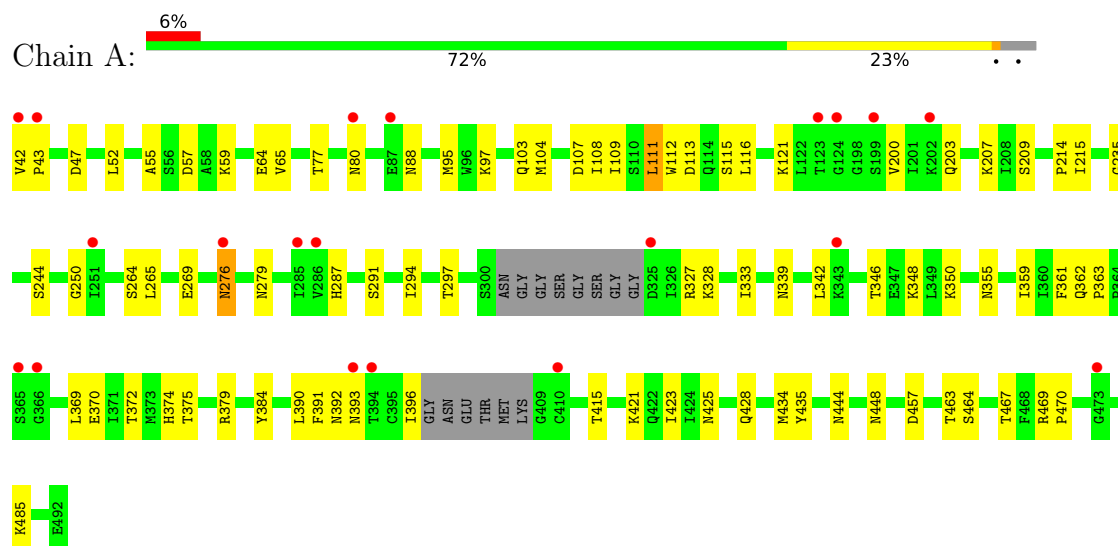
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	46	Total	O	0	0
			46	46		

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 LM/HT Clade A/E CRF01 gp120 core



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	66.90Å 67.69Å 86.47Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	28.51 – 2.50 28.51 – 2.50	Depositor EDS
% Data completeness (in resolution range)	95.0 (28.51-2.50) 96.5 (28.51-2.50)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.03 (at 2.51Å)	Xtriage
Refinement program	PHENIX (1.15.2_3472: ???)	Depositor
R, R_{free}	0.263 , 0.305 0.265 , 0.314	Depositor DCC
R_{free} test set	653 reflections (4.78%)	wwPDB-VP
Wilson B-factor (Å ²)	44.4	Xtriage
Anisotropy	0.684	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 32.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.018 for k,h,-l	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	2875	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.07% 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: EPE, NAG

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.35	0/2731	0.59	1/3708 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	111	LEU	CA-CB-CG	5.34	127.59	115.30

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	2674	0	2603	66	0
2	A	140	0	130	1	0
3	A	15	0	17	2	0
4	A	46	0	0	18	0
All	All	2875	0	2750	67	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 12.

All (67) 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:342:LEU:HD12	1:A:396:ILE:HD11	1.17	1.14
1:A:250:GLY:O	4:A:601:HOH:O	1.88	0.92
1:A:207:LYS:O	4:A:602:HOH:O	1.88	0.90
1:A:52:LEU:H	1:A:103:GLN:HE22	1.28	0.81
1:A:370:GLU:OE2	1:A:384:TYR:OH	2.00	0.78
1:A:203:GLN:O	4:A:603:HOH:O	2.05	0.75
1:A:485:LYS:NZ	4:A:614:HOH:O	2.23	0.71
1:A:379:ARG:NH2	1:A:444:ASN:H	1.90	0.70
1:A:276:ASN:OD1	1:A:279:ASN:N	2.26	0.68
1:A:297:THR:OG1	1:A:444:ASN:ND2	2.27	0.68
1:A:116:LEU:O	4:A:608:HOH:O	2.12	0.67
1:A:339:ASN:ND2	4:A:615:HOH:O	2.26	0.67
1:A:423:ILE:O	4:A:607:HOH:O	2.12	0.67
1:A:327:ARG:HD2	4:A:627:HOH:O	1.95	0.67
1:A:421:LYS:O	4:A:606:HOH:O	2.12	0.66
1:A:342:LEU:CD1	1:A:396:ILE:HD11	2.11	0.65
1:A:264:SER:O	1:A:287:HIS:NE2	2.25	0.65
1:A:435:TYR:OH	4:A:605:HOH:O	2.11	0.64
1:A:65:VAL:HB	1:A:115:SER:HB3	1.81	0.62
1:A:291:SER:HB2	1:A:448:ASN:HB3	1.82	0.62
1:A:350:LYS:NZ	1:A:355:ASN:O	2.33	0.61
1:A:59:LYS:NZ	4:A:616:HOH:O	2.30	0.60
1:A:370:GLU:OE1	1:A:425:ASN:HB2	2.02	0.58
1:A:370:GLU:HA	1:A:375:THR:HG21	1.87	0.57
1:A:113:ASP:OD1	4:A:609:HOH:O	2.18	0.56
1:A:463:THR:HG22	1:A:464:SER:H	1.70	0.56
1:A:269:GLU:OE2	1:A:348:LYS:NZ	2.39	0.56
1:A:57:ASP:O	1:A:59:LYS:HE2	2.06	0.55
1:A:342:LEU:HD11	1:A:361:PHE:CE2	2.41	0.55
1:A:333:ILE:HD12	1:A:390:LEU:HD21	1.88	0.55
1:A:457:ASP:OD2	1:A:467:THR:OG1	2.25	0.55
1:A:42:VAL:HB	1:A:43:PRO:HD3	1.89	0.54
1:A:423:ILE:HG12	1:A:434:MET:HG3	1.88	0.54
1:A:103:GLN:NE2	3:A:511:EPE:O8	2.28	0.54
1:A:265:LEU:HD21	1:A:291:SER:HB3	1.90	0.53
1:A:64:GLU:HA	1:A:209:SER:HB3	1.91	0.51
1:A:342:LEU:O	1:A:346:THR:HG23	2.11	0.50
1:A:342:LEU:HD13	1:A:342:LEU:O	2.13	0.49
1:A:363:PRO:O	1:A:469:ARG:NH1	2.36	0.48
1:A:350:LYS:HG2	1:A:355:ASN:HA	1.96	0.47
1:A:428:GLN:HG2	4:A:609:HOH:O	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:415:THR:OG1	4:A:610:HOH:O	2.20	0.47
1:A:346:THR:HG22	1:A:361:PHE:CE2	2.50	0.47
1:A:369:LEU:HA	1:A:372:THR:HG22	1.98	0.46
1:A:350:LYS:HA	1:A:359:ILE:HD11	1.97	0.46
1:A:374:HIS:NE2	4:A:604:HOH:O	2.08	0.46
1:A:97:LYS:HD2	1:A:97:LYS:HA	1.66	0.45
1:A:107:ASP:OD1	3:A:511:EPE:H52	2.16	0.45
2:A:510:NAG:O3	4:A:611:HOH:O	2.21	0.45
1:A:391:PHE:CE2	1:A:470:PRO:HG3	2.52	0.45
1:A:88:ASN:ND2	4:A:622:HOH:O	2.50	0.44
1:A:109:ILE:HG23	4:A:609:HOH:O	2.17	0.44
1:A:392:ASN:OD1	1:A:393:ASN:N	2.51	0.44
1:A:121:LYS:HE3	1:A:121:LYS:HB2	1.77	0.43
1:A:95:MET:SD	1:A:235:GLY:HA3	2.58	0.43
1:A:346:THR:CG2	1:A:361:PHE:HE2	2.31	0.43
1:A:350:LYS:CA	1:A:359:ILE:HD11	2.48	0.43
1:A:111:LEU:O	1:A:115:SER:OG	2.22	0.42
1:A:421:LYS:HE2	1:A:421:LYS:HB3	1.81	0.42
1:A:214:PRO:O	1:A:215:ILE:HD13	2.19	0.42
1:A:104:MET:O	1:A:108:ILE:HG12	2.20	0.42
1:A:294:ILE:HD12	1:A:333:ILE:HD11	2.02	0.41
1:A:207:LYS:HD2	1:A:207:LYS:HA	1.68	0.41
1:A:112:TRP:CE3	1:A:116:LEU:CD1	3.04	0.41
1:A:55:ALA:HB1	1:A:77:THR:HA	2.03	0.40
1:A:361:PHE:O	1:A:362:GLN:NE2	2.53	0.40
1:A:362:GLN:HB3	1:A:469:ARG:NH1	2.37	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	335/355 (94%)	311 (93%)	22 (7%)	2 (1%)	25	43

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	276	ASN
1	A	200	VAL

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	305/313 (97%)	301 (99%)	4 (1%)	69	87

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	47	ASP
1	A	80	ASN
1	A	244	SER
1	A	328	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	103	GLN
1	A	444	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

11 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$
2	NAG	A	506	1	14,14,15	0.46	0	17,19,21	0.49	0
2	NAG	A	501	1	14,14,15	0.27	0	17,19,21	0.62	0
3	EPE	A	511	-	15,15,15	0.67	0	18,20,20	2.17	7 (38%)
2	NAG	A	503	1	14,14,15	0.26	0	17,19,21	0.61	1 (5%)
2	NAG	A	509	1	14,14,15	0.87	1 (7%)	17,19,21	0.66	0
2	NAG	A	507	1	14,14,15	1.11	2 (14%)	17,19,21	0.97	1 (5%)
2	NAG	A	510	1	14,14,15	0.30	0	17,19,21	1.00	1 (5%)
2	NAG	A	504	1	14,14,15	0.68	0	17,19,21	0.64	0
2	NAG	A	505	1	14,14,15	0.80	1 (7%)	17,19,21	0.92	1 (5%)
2	NAG	A	508	1	14,14,15	0.46	0	17,19,21	0.64	1 (5%)
2	NAG	A	502	1	14,14,15	0.82	1 (7%)	17,19,21	0.53	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	A	506	1	-	2/6/23/26	0/1/1/1
2	NAG	A	501	1	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	EPE	A	511	-	-	6/9/19/19	0/1/1/1
2	NAG	A	503	1	-	2/6/23/26	0/1/1/1
2	NAG	A	509	1	-	2/6/23/26	0/1/1/1
2	NAG	A	507	1	-	0/6/23/26	0/1/1/1
2	NAG	A	510	1	-	0/6/23/26	0/1/1/1
2	NAG	A	504	1	-	0/6/23/26	0/1/1/1
2	NAG	A	505	1	-	1/6/23/26	0/1/1/1
2	NAG	A	508	1	-	2/6/23/26	0/1/1/1
2	NAG	A	502	1	-	0/6/23/26	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	507	NAG	O5-C1	-3.05	1.38	1.43
2	A	505	NAG	O5-C1	-2.66	1.39	1.43
2	A	509	NAG	O5-C1	-2.65	1.39	1.43
2	A	507	NAG	C1-C2	2.36	1.55	1.52
2	A	502	NAG	O5-C1	-2.25	1.40	1.43

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	511	EPE	C5-N4-C3	3.88	117.56	108.83
3	A	511	EPE	O1S-S-C10	3.69	111.36	106.92
2	A	510	NAG	C1-O5-C5	3.69	117.19	112.19
3	A	511	EPE	C7-N4-C5	3.53	120.27	111.23
3	A	511	EPE	C7-N4-C3	3.12	119.22	111.23
2	A	505	NAG	C1-O5-C5	-3.11	107.97	112.19
2	A	507	NAG	C4-C3-C2	2.80	115.12	111.02
3	A	511	EPE	O2S-S-C10	2.77	110.25	106.92
3	A	511	EPE	C9-N1-C6	-2.60	104.58	111.23
3	A	511	EPE	C9-N1-C2	-2.23	105.53	111.23
2	A	508	NAG	C1-O5-C5	2.20	115.18	112.19
2	A	503	NAG	C1-O5-C5	2.12	115.06	112.19

There are no chirality outliers.

All (15) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	511	EPE	C8-C7-N4-C5

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Mol	Chain	Res	Type	Atoms
3	A	511	EPE	S-C10-C9-N1
3	A	511	EPE	C9-C10-S-O1S
2	A	506	NAG	O5-C5-C6-O6
2	A	509	NAG	O5-C5-C6-O6
2	A	509	NAG	C4-C5-C6-O6
2	A	506	NAG	C4-C5-C6-O6
3	A	511	EPE	C9-C10-S-O3S
2	A	508	NAG	O5-C5-C6-O6
2	A	505	NAG	O5-C5-C6-O6
3	A	511	EPE	C9-C10-S-O2S
3	A	511	EPE	C8-C7-N4-C3
2	A	503	NAG	C4-C5-C6-O6
2	A	503	NAG	O5-C5-C6-O6
2	A	508	NAG	C4-C5-C6-O6

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	511	EPE	2	0
2	A	510	NAG	1	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	341/355 (96%)	0.61	20 (5%) 22 23	30, 43, 66, 78	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	42	VAL	6.1
1	A	394	THR	5.0
1	A	473	GLY	3.6
1	A	123	THR	3.6
1	A	285	ILE	3.1
1	A	366	GLY	3.1
1	A	124	GLY	2.8
1	A	410	CYS	2.7
1	A	276	ASN	2.5
1	A	325	ASP	2.5
1	A	365	SER	2.4
1	A	251	ILE	2.3
1	A	43	PRO	2.3
1	A	87	GLU	2.2
1	A	80	ASN	2.2
1	A	393	ASN	2.1
1	A	199	SER	2.1
1	A	286	VAL	2.1
1	A	202	LYS	2.0
1	A	343	LYS	2.0

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

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

6.3 Carbohydrates [i](#)

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
2	NAG	A	509	14/15	0.45	0.40	62,71,81,81	0
2	NAG	A	504	14/15	0.57	0.43	63,71,76,81	0
2	NAG	A	507	14/15	0.66	0.27	58,68,83,88	0
2	NAG	A	502	14/15	0.72	0.32	59,69,79,81	0
2	NAG	A	510	14/15	0.76	0.48	52,72,77,77	0
2	NAG	A	505	14/15	0.81	0.22	41,51,57,66	0
2	NAG	A	506	14/15	0.82	0.28	54,57,58,58	0
2	NAG	A	501	14/15	0.83	0.18	35,48,58,58	0
2	NAG	A	508	14/15	0.86	0.27	43,53,58,66	0
2	NAG	A	503	14/15	0.86	0.17	28,35,47,51	0
3	EPE	A	511	15/15	0.93	0.20	34,38,41,43	0

6.5 Other polymers [i](#)

There are no such residues in this entry.