



Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 04:32 AM GMT

PDB ID : 1VPS
Title : POLYOMAVIRUS VP1 PENTAMER COMPLEXED WITH A DISIALY-
LATED HEXASACCHARIDE
Authors : Stehle, T.; Harrison, S.C.
Deposited on : 1997-03-07
Resolution : 1.90 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

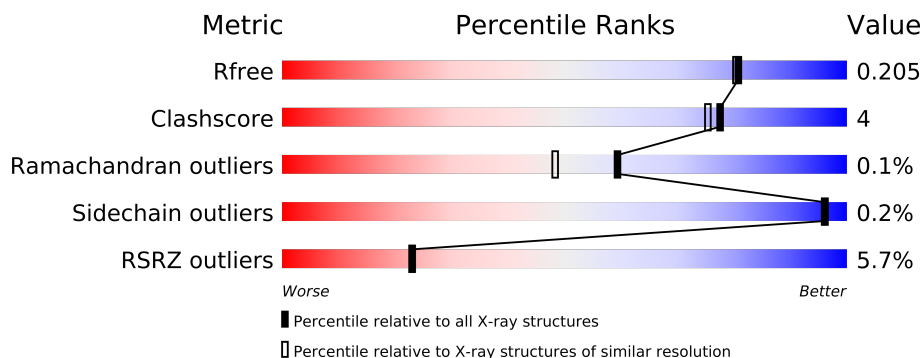
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 1.90 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}	66092	3684 (1.90-1.90)
Clashscore	79885	4465 (1.90-1.90)
Ramachandran outliers	78287	4413 (1.90-1.90)
Sidechain outliers	78261	4414 (1.90-1.90)
RSRZ outliers	66119	3686 (1.90-1.90)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	289	
1	B	289	
1	C	289	
1	D	289	
1	E	289	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 13591 atoms, of which 0 are hydrogen and 0 are deuterium.

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 POLYOMAVIRUS VP1 PENTAMER.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	285	Total	C	N	O	S	0	0	0
			2220	1404	373	432	11			
1	B	289	Total	C	N	O	S	0	0	0
			2254	1427	378	438	11			
1	C	285	Total	C	N	O	S	0	0	0
			2220	1404	373	432	11			
1	D	285	Total	C	N	O	S	0	0	0
			2220	1404	373	432	11			
1	E	285	Total	C	N	O	S	0	0	0
			2220	1404	373	432	11			

- Molecule 2 is a polymer of unknown type called SUGAR (4-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	4	Total	C	N	O	0	0
			66	36	3	27		
2	B	4	Total	C	N	O	0	0
			66	36	3	27		
2	C	4	Total	C	N	O	0	0
			66	36	3	27		
2	D	4	Total	C	N	O	0	0
			66	36	3	27		
2	E	4	Total	C	N	O	0	0
			66	36	3	27		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	391	Total	O	0	0
			391	391		
3	B	444	Total	O	0	0
			444	444		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	C	452	Total 452	O 452	0	0
3	D	436	Total 436	O 436	0	0
3	E	404	Total 404	O 404	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	221.50Å 221.50Å 99.80Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 1.90 24.74 – 1.92	Depositor EDS
% Data completeness (in resolution range)	85.6 (20.00-1.90) 88.7 (24.74-1.92)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.79 (at 1.92Å)	Xtriage
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.177 , 0.205 0.178 , 0.205	Depositor DCC
R_{free} test set	7533 reflections (4.16%)	DCC
Wilson B-factor (Å ²)	20.5	Xtriage
Anisotropy	0.100	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 59.4	EDS
Estimated twinning fraction	0.019 for -h,-k,l	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 188609 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	13591	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.21% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: SIA, GAL, 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	0.38	0/2273	0.63	0/3098
1	B	0.39	0/2310	0.65	0/3151
1	C	0.40	0/2273	0.65	0/3098
1	D	0.42	1/2273 (0.0%)	0.64	0/3098
1	E	0.38	0/2273	0.63	0/3098
All	All	0.39	1/11402 (0.0%)	0.64	0/15543

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	293	ASN	C-N	-5.84	1.20	1.34

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2220	0	2187	15	0
1	B	2254	0	2216	20	0
1	C	2220	0	2187	20	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	2220	0	2187	18	0
1	E	2220	0	2187	16	0
2	A	66	0	56	0	0
2	B	66	0	56	1	0
2	C	66	0	56	0	0
2	D	66	0	56	1	0
2	E	66	0	56	0	0
3	A	391	0	0	3	0
3	B	444	0	0	6	0
3	C	452	0	0	3	0
3	D	436	0	0	2	0
3	E	404	0	0	6	0
All	All	13591	0	11244	83	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 4.

All (83) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:273:CYS:SG	3:A:651:HOH:O	2.50	0.69
1:C:35:GLU:HB3	1:C:316:LYS:HB2	1.75	0.68
1:D:35:GLU:HB3	1:D:316:LYS:HB2	1.75	0.66
1:E:273:CYS:SG	3:E:705:HOH:O	2.50	0.66
1:B:35:GLU:HB3	1:B:316:LYS:HB2	1.85	0.57
1:A:33:GLY:O	1:A:316:LYS:HB3	2.04	0.57
1:E:35:GLU:HB3	1:E:316:LYS:HB2	1.87	0.57
1:C:106:PRO:HG2	1:C:120:TRP:CZ2	2.41	0.56
1:D:102:LYS:NZ	1:D:277:GLY:HA3	2.20	0.56
1:E:262:VAL:HG13	3:E:692:HOH:O	2.06	0.56
1:B:153:ILE:HA	1:C:81:LEU:HD21	1.87	0.56
1:C:305:TYR:HE1	3:C:502:HOH:O	1.88	0.56
1:D:39:LEU:HD23	1:D:313:ARG:HD2	1.89	0.55
1:C:305:TYR:HD2	3:C:717:HOH:O	1.90	0.55
1:B:224:VAL:HG11	1:B:283:VAL:HG21	1.90	0.54
3:B:626:HOH:O	1:C:294:TYR:HA	2.07	0.54
1:D:102:LYS:HZ1	1:D:277:GLY:HA3	1.73	0.54
1:A:81:LEU:HD21	1:E:153:ILE:HA	1.90	0.53
1:D:121:GLU:OE1	1:D:313:ARG:HD3	2.09	0.53
1:B:115:ASP:HB2	3:B:651:HOH:O	2.08	0.52
1:C:153:ILE:HA	1:D:81:LEU:HD21	1.91	0.52
1:B:106:PRO:HG2	1:B:120:TRP:CZ2	2.44	0.52
1:A:121:GLU:OE1	1:A:313:ARG:HD3	2.10	0.51

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
3:B:749:HOH:O	1:C:242:ASN:HB2	2.09	0.51
1:C:121:GLU:OE1	1:C:313:ARG:HD3	2.10	0.51
1:C:236:ASN:ND2	1:C:272:LEU:O	2.44	0.51
1:B:102:LYS:NZ	1:B:277:GLY:HA3	2.27	0.49
1:E:208:LEU:HD13	3:E:694:HOH:O	2.12	0.49
1:E:121:GLU:OE1	1:E:313:ARG:HD3	2.13	0.49
1:E:40:VAL:HB	1:E:312:LYS:HB2	1.94	0.49
1:E:238:ARG:HD3	3:E:637:HOH:O	2.13	0.48
1:C:53:LEU:HG	3:C:759:HOH:O	2.13	0.48
1:D:293:ASN:HB3	3:D:733:HOH:O	2.13	0.48
1:B:236:ASN:ND2	1:B:272:LEU:O	2.46	0.48
1:A:65:SER:HB3	1:A:68:GLU:HB2	1.95	0.48
1:D:106:PRO:HG2	1:D:120:TRP:CZ2	2.49	0.48
1:A:39:LEU:HD23	1:A:313:ARG:HD2	1.96	0.47
1:C:236:ASN:ND2	1:C:274:LYS:HE3	2.29	0.47
1:A:106:PRO:HG2	1:A:120:TRP:CZ2	2.49	0.47
1:C:33:GLY:O	1:C:316:LYS:HB3	2.14	0.47
1:C:236:ASN:HD22	1:C:274:LYS:HE3	1.79	0.47
1:E:236:ASN:ND2	1:E:272:LEU:O	2.48	0.47
1:E:106:PRO:HG2	1:E:120:TRP:CZ2	2.50	0.47
1:E:224:VAL:HG11	1:E:283:VAL:HG21	1.97	0.46
1:E:97:THR:OG1	1:E:220:GLY:HA2	2.15	0.46
1:E:33:GLY:O	1:E:316:LYS:HB3	2.16	0.46
1:D:40:VAL:HB	1:D:312:LYS:HB2	1.97	0.46
1:B:121:GLU:OE1	1:B:313:ARG:HD3	2.16	0.45
1:B:44:ASP:O	1:B:312:LYS:HE3	2.17	0.45
3:A:679:HOH:O	1:B:208:LEU:HD13	2.17	0.45
1:A:40:VAL:HB	1:A:312:LYS:HB2	1.99	0.45
3:B:751:HOH:O	1:C:208:LEU:HD13	2.17	0.44
1:B:319:TYR:HA	1:B:320:PRO:HD3	1.80	0.44
1:E:38:ASP:HB2	3:E:573:HOH:O	2.17	0.44
1:B:83:THR:HG23	2:B:4:SIA:H91	2.00	0.44
1:D:224:VAL:HG11	1:D:283:VAL:HG21	2.00	0.44
1:D:97:THR:OG1	1:D:220:GLY:HA2	2.18	0.43
1:B:68:GLU:HG2	3:B:752:HOH:O	2.18	0.43
1:A:153:ILE:HA	1:B:81:LEU:HD21	2.00	0.43
1:D:52:PHE:HB2	3:E:694:HOH:O	2.18	0.43
1:B:39:LEU:HD23	1:B:313:ARG:HD2	2.01	0.43
1:B:40:VAL:HB	1:B:312:LYS:HB2	2.01	0.43
1:C:40:VAL:HB	1:C:312:LYS:HB2	1.99	0.42
1:B:251:PRO:HG3	3:B:727:HOH:O	2.19	0.42
1:D:236:ASN:ND2	1:D:272:LEU:O	2.50	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:250:PRO:HA	1:A:251:PRO:HD3	1.95	0.42
1:A:122:ALA:HB3	1:A:271:PRO:HD2	2.02	0.42
1:D:83:THR:HG23	2:D:4:SIA:H91	2.02	0.42
1:E:106:PRO:HG2	1:E:120:TRP:HZ2	1.83	0.42
1:D:276:GLU:HB3	3:D:722:HOH:O	2.20	0.41
1:C:274:LYS:HE2	1:C:274:LYS:HB3	1.92	0.41
1:A:251:PRO:HD2	1:B:245:GLY:HA3	2.02	0.41
1:A:97:THR:OG1	1:A:220:GLY:HA2	2.19	0.41
1:C:142:ASN:HB2	1:C:154:SER:OG	2.20	0.41
1:D:106:PRO:HG2	1:D:120:TRP:HZ2	1.84	0.41
1:E:274:LYS:HE2	1:E:274:LYS:HB3	1.89	0.41
1:C:251:PRO:HD2	1:D:245:GLY:HA3	2.03	0.40
1:D:250:PRO:HA	1:D:251:PRO:HD3	1.96	0.40
1:A:236:ASN:ND2	1:A:272:LEU:O	2.55	0.40
1:C:44:ASP:O	1:C:312:LYS:HE3	2.21	0.40
1:B:250:PRO:HA	1:B:251:PRO:HD3	1.89	0.40
1:B:97:THR:OG1	1:B:220:GLY:HA2	2.21	0.40
1:A:52:PHE:HB3	3:A:656:HOH:O	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	283/289 (98%)	273 (96%)	9 (3%)	1 (0%)	43	29
1	B	287/289 (99%)	273 (95%)	14 (5%)	0	100	100
1	C	283/289 (98%)	268 (95%)	15 (5%)	0	100	100
1	D	283/289 (98%)	272 (96%)	11 (4%)	0	100	100
1	E	283/289 (98%)	271 (96%)	12 (4%)	0	100	100
All	All	1419/1445 (98%)	1357 (96%)	61 (4%)	1 (0%)	59	48

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	34	MET

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	249/253 (98%)	247 (99%)	2 (1%)	89	89
1	B	253/253 (100%)	253 (100%)	0	100	100
1	C	249/253 (98%)	249 (100%)	0	100	100
1	D	249/253 (98%)	249 (100%)	0	100	100
1	E	249/253 (98%)	249 (100%)	0	100	100
All	All	1249/1265 (99%)	1247 (100%)	2 (0%)	96	96

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	188	GLU
1	A	236	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

20 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	SIA	A	1	2	20,20,21	3.96	2 (10%)	23,28,31	2.50	3 (13%)
2	GAL	A	2	2	10,11,12	0.53	0	11,15,17	0.54	0
2	NAG	A	3	2	15,15,15	0.84	1 (6%)	21,21,21	1.28	5 (23%)
2	SIA	A	4	2	20,20,21	4.25	5 (25%)	23,28,31	2.64	3 (13%)
2	SIA	B	1	2	20,20,21	3.90	2 (10%)	23,28,31	2.48	1 (4%)
2	GAL	B	2	2	10,11,12	0.45	0	11,15,17	0.53	0
2	NAG	B	3	2	15,15,15	0.47	0	21,21,21	0.52	0
2	SIA	B	4	2	20,20,21	4.05	2 (10%)	23,28,31	2.47	2 (8%)
2	SIA	C	1	2	20,20,21	3.87	2 (10%)	23,28,31	2.61	2 (8%)
2	GAL	C	2	2	10,11,12	0.64	0	11,15,17	0.54	0
2	NAG	C	3	2	15,15,15	0.90	1 (6%)	21,21,21	1.11	1 (4%)
2	SIA	C	4	2	20,20,21	4.14	3 (15%)	23,28,31	2.74	3 (13%)
2	SIA	D	1	2	20,20,21	3.96	2 (10%)	23,28,31	2.52	2 (8%)
2	GAL	D	2	2	10,11,12	0.47	0	11,15,17	0.51	0
2	NAG	D	3	2	15,15,15	0.51	0	21,21,21	0.80	0
2	SIA	D	4	2	20,20,21	4.38	2 (10%)	23,28,31	2.42	2 (8%)
2	SIA	E	1	2	20,20,21	4.00	2 (10%)	23,28,31	2.55	3 (13%)
2	GAL	E	2	2	10,11,12	0.52	0	11,15,17	0.60	0
2	NAG	E	3	2	15,15,15	0.73	0	21,21,21	0.91	1 (4%)
2	SIA	E	4	2	20,20,21	3.86	4 (20%)	23,28,31	2.72	2 (8%)

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	SIA	A	1	2	-	0/15/34/38	0/1/1/1
2	GAL	A	2	2	-	0/2/19/22	0/1/1/1
2	NAG	A	3	2	-	0/6/26/26	0/1/1/1
2	SIA	A	4	2	-	0/15/34/38	0/1/1/1
2	SIA	B	1	2	-	0/15/34/38	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GAL	B	2	2	-	0/2/19/22	0/1/1/1
2	NAG	B	3	2	-	0/6/26/26	0/1/1/1
2	SIA	B	4	2	-	0/15/34/38	0/1/1/1
2	SIA	C	1	2	-	0/15/34/38	0/1/1/1
2	GAL	C	2	2	-	0/2/19/22	0/1/1/1
2	NAG	C	3	2	-	0/6/26/26	0/1/1/1
2	SIA	C	4	2	-	0/15/34/38	0/1/1/1
2	SIA	D	1	2	-	0/15/34/38	0/1/1/1
2	GAL	D	2	2	-	0/2/19/22	0/1/1/1
2	NAG	D	3	2	-	0/6/26/26	0/1/1/1
2	SIA	D	4	2	-	0/15/34/38	0/1/1/1
2	SIA	E	1	2	-	0/15/34/38	0/1/1/1
2	GAL	E	2	2	-	0/2/19/22	0/1/1/1
2	NAG	E	3	2	-	0/6/26/26	0/1/1/1
2	SIA	E	4	2	-	0/15/34/38	0/1/1/1

All (28) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	4	SIA	O6-C2	19.00	1.45	1.28
2	C	4	SIA	O6-C2	17.78	1.44	1.28
2	B	4	SIA	O6-C2	17.66	1.44	1.28
2	E	1	SIA	O6-C2	17.53	1.44	1.28
2	A	4	SIA	O6-C2	17.45	1.44	1.28
2	D	1	SIA	O6-C2	17.38	1.44	1.28
2	A	1	SIA	O6-C2	17.31	1.44	1.28
2	B	1	SIA	O6-C2	17.07	1.43	1.28
2	C	1	SIA	O6-C2	16.96	1.43	1.28
2	E	4	SIA	O6-C2	16.55	1.43	1.28
2	A	4	SIA	C3-C4	5.22	1.58	1.52
2	A	4	SIA	C4-C5	3.31	1.56	1.53
2	C	4	SIA	C4-C5	3.03	1.55	1.53
2	A	4	SIA	C3-C2	2.58	1.55	1.46
2	C	4	SIA	C3-C2	2.54	1.55	1.46
2	E	4	SIA	C4-C5	2.50	1.55	1.53
2	D	4	SIA	C3-C2	2.48	1.55	1.46
2	E	1	SIA	C3-C2	2.33	1.54	1.46
2	C	1	SIA	C3-C2	2.26	1.54	1.46
2	B	4	SIA	C3-C2	2.26	1.54	1.46
2	E	4	SIA	C3-C4	2.25	1.55	1.52
2	E	4	SIA	C3-C2	2.18	1.54	1.46
2	A	3	NAG	C1-C2	2.17	1.55	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	1	SIA	C3-C2	2.11	1.53	1.46
2	A	1	SIA	C3-C2	2.04	1.53	1.46
2	C	3	NAG	C1-C2	2.04	1.55	1.53
2	A	4	SIA	C6-C5	2.03	1.56	1.53
2	B	1	SIA	C3-C2	2.02	1.53	1.46

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	4	SIA	O6-C2-C3	-12.46	108.63	124.91
2	E	4	SIA	O6-C2-C3	-12.23	108.93	124.91
2	A	4	SIA	O6-C2-C3	-11.69	109.64	124.91
2	C	1	SIA	O6-C2-C3	-11.51	109.88	124.91
2	E	1	SIA	O6-C2-C3	-11.35	110.09	124.91
2	D	1	SIA	O6-C2-C3	-11.30	110.14	124.91
2	B	1	SIA	O6-C2-C3	-11.22	110.25	124.91
2	A	1	SIA	O6-C2-C3	-11.22	110.25	124.91
2	B	4	SIA	O6-C2-C3	-11.06	110.45	124.91
2	D	4	SIA	O6-C2-C3	-10.69	110.94	124.91
2	C	1	SIA	C7-C6-C5	-3.03	109.77	114.24
2	E	4	SIA	C8-C7-C6	-2.92	107.33	112.99
2	C	3	NAG	C1-C2-N2	-2.86	107.53	110.85
2	D	4	SIA	C3-C4-C5	-2.80	106.57	110.41
2	A	4	SIA	C8-C7-C6	-2.78	107.59	112.99
2	A	3	NAG	O5-C1-C2	2.77	112.41	109.61
2	B	4	SIA	C8-C7-C6	-2.58	107.99	112.99
2	E	1	SIA	C7-C6-C5	-2.53	110.52	114.24
2	A	3	NAG	C3-C4-C5	-2.22	106.24	110.20
2	C	4	SIA	C8-C7-C6	-2.22	108.70	112.99
2	A	1	SIA	C3-C2-C1	-2.19	108.84	121.08
2	A	3	NAG	C6-C5-C4	2.19	118.28	113.00
2	A	3	NAG	C1-C2-C3	2.18	113.72	110.59
2	A	3	NAG	C1-C2-N2	-2.16	108.34	110.85
2	E	3	NAG	C6-C5-C4	2.13	118.14	113.00
2	D	1	SIA	C7-C6-C5	-2.11	111.13	114.24
2	A	4	SIA	C3-C2-C1	-2.09	109.38	121.08
2	A	1	SIA	C8-C7-C6	-2.08	108.96	112.99
2	C	4	SIA	O6-C6-C7	2.03	109.50	105.73
2	E	1	SIA	C3-C2-C1	-2.03	109.75	121.08

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

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	285/289 (98%)	0.07	18 (6%)	19 19	16, 25, 59, 95	0
1	B	289/289 (100%)	-0.15	10 (3%)	42 43	14, 22, 53, 82	0
1	C	285/289 (98%)	-0.06	16 (5%)	24 24	14, 20, 54, 92	0
1	D	285/289 (98%)	-0.03	17 (5%)	21 21	13, 20, 57, 92	0
1	E	285/289 (98%)	0.05	17 (5%)	21 21	17, 25, 59, 94	0
All	All	1429/1445 (98%)	-0.02	78 (5%)	23 24	13, 22, 58, 95	0

All (78) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	112	LEU	13.5
1	A	112	LEU	13.0
1	D	112	LEU	12.2
1	D	33	GLY	10.9
1	A	113	THR	10.9
1	E	113	THR	10.9
1	E	112	LEU	10.8
1	C	113	THR	9.7
1	C	33	GLY	9.5
1	B	32	GLY	9.3
1	D	113	THR	8.9
1	D	32	GLY	8.8
1	A	33	GLY	8.3
1	E	33	GLY	8.2
1	B	33	GLY	8.2
1	E	32	GLY	7.7
1	A	111	ASP	7.6
1	E	110	GLU	7.6
1	D	34	MET	7.3
1	C	111	ASP	7.3

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Mol	Chain	Res	Type	RSRZ
1	E	114	CYS	7.2
1	A	110	GLU	6.9
1	E	111	ASP	6.8
1	D	111	ASP	6.8
1	A	34	MET	6.8
1	A	114	CYS	6.6
1	D	114	CYS	6.2
1	D	110	GLU	6.2
1	B	320	PRO	6.0
1	A	32	GLY	5.9
1	C	114	CYS	5.8
1	E	34	MET	5.8
1	C	32	GLY	5.5
1	C	110	GLU	5.3
1	C	34	MET	4.8
1	E	316	LYS	4.6
1	E	109	ASN	4.5
1	E	115	ASP	4.5
1	A	109	ASN	4.4
1	C	107	MET	4.3
1	C	115	ASP	4.3
1	D	276	GLU	4.2
1	D	109	ASN	4.2
1	E	107	MET	4.2
1	A	115	ASP	4.1
1	C	276	GLU	4.0
1	B	110	GLU	3.9
1	A	107	MET	3.8
1	D	107	MET	3.8
1	C	316	LYS	3.6
1	E	276	GLU	3.5
1	D	316	LYS	3.4
1	B	107	MET	3.4
1	B	276	GLU	3.4
1	D	115	ASP	3.3
1	A	294	TYR	3.3
1	C	109	ASN	3.2
1	B	317	ASN	3.2
1	B	34	MET	3.2
1	A	276	GLU	3.1
1	B	319	TYR	3.0
1	A	316	LYS	3.0

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Mol	Chain	Res	Type	RSRZ
1	D	108	LEU	2.9
1	A	117	LEU	2.8
1	E	294	TYR	2.8
1	E	117	LEU	2.6
1	C	35	GLU	2.6
1	A	108	LEU	2.5
1	C	108	LEU	2.4
1	E	108	LEU	2.4
1	E	35	GLU	2.4
1	A	35	GLU	2.4
1	D	274	LYS	2.3
1	C	117	LEU	2.3
1	A	274	LYS	2.3
1	D	294	TYR	2.2
1	B	318	PRO	2.2
1	D	117	LEU	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	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	SIA	B	4	20/21	0.54	28.18	80,82,83,84	0
2	SIA	E	4	20/21	0.50	21.45	78,81,82,82	0
2	SIA	C	4	20/21	0.56	18.48	67,71,77,77	0
2	SIA	A	4	20/21	0.45	17.88	66,71,73,73	0
2	SIA	D	4	20/21	0.22	10.41	46,56,62,63	0
2	NAG	D	3	15/15	0.26	7.12	41,51,53,54	0
2	GAL	E	2	11/12	0.13	3.56	40,46,51,53	0
2	GAL	A	2	11/12	0.11	2.61	37,44,46,48	0
2	SIA	B	1	20/21	0.10	1.75	24,27,31,35	0
2	SIA	A	1	20/21	0.11	1.44	28,32,36,40	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	SIA	E	1	20/21	0.11	1.13	31,34,37,38	0
2	GAL	C	2	11/12	0.08	1.04	28,32,35,38	0
2	GAL	B	2	11/12	0.08	0.91	32,37,41,44	0
2	SIA	D	1	20/21	0.11	0.71	23,25,29,33	0
2	SIA	C	1	20/21	0.09	0.21	22,24,29,34	0
2	GAL	D	2	11/12	0.07	0.05	28,31,33,36	0
2	NAG	B	3	15/15	0.39	-	53,66,72,77	0
2	NAG	A	3	15/15	0.37	-	54,62,66,69	0
2	NAG	E	3	15/15	0.40	-	59,69,74,78	0
2	NAG	C	3	15/15	0.41	-	46,58,63,67	0

6.4 Ligands ⓘ

There are no ligands in this entry.

6.5 Other polymers ⓘ

There are no such residues in this entry.