



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:28 PM GMT

PDB ID : 4P26
Title : Structure of the P domain from a GI.7 Norovirus variant in complex with A-type 2 HBGA
Authors : Shanker, S.; Czako, R.; Sankaran, B.; Atmar, R.; Estes, M.; Prasad, B.V.V.
Deposited on : 2014-03-01
Resolution : 1.90 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

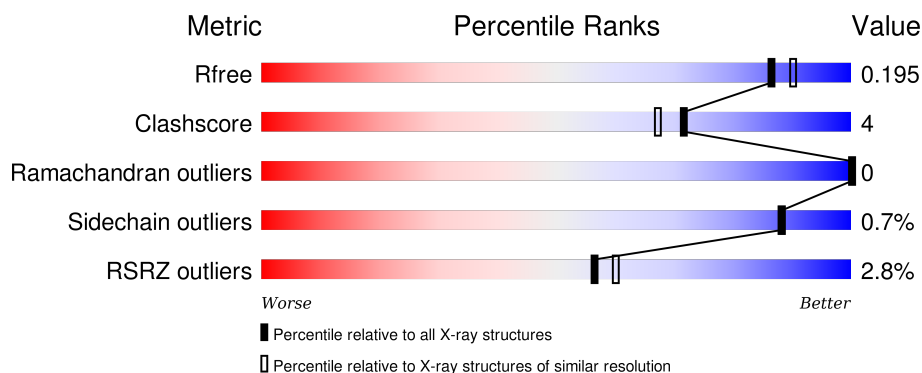
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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}	91344	4755 (1.90-1.90)
Clashscore	102246	5398 (1.90-1.90)
Ramachandran outliers	100387	5338 (1.90-1.90)
Sidechain outliers	100360	5339 (1.90-1.90)
RSRZ outliers	91569	4766 (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. 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	301	<div> <div>3%</div> <div>91%</div> <div>6%</div> <div>.</div> </div>
1	B	301	<div> <div>2%</div> <div>86%</div> <div>8%</div> <div>5%</div> </div>
1	C	301	<div> <div>3%</div> <div>88%</div> <div>8%</div> <div>.</div> </div>
1	D	301	<div> <div>3%</div> <div>85%</div> <div>9%</div> <div>6%</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
2	GAL	A	602	-	-	X	-
2	NAG	A	604	-	-	X	-
2	A2G	B	601	-	-	-	X

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9866 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 P domain of VP1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	292	Total	C	N	O	S	0	0	0
			2226	1417	378	423	8			
1	B	285	Total	C	N	O	S	0	0	0
			2163	1379	364	412	8			
1	C	289	Total	C	N	O	S	0	0	0
			2198	1401	370	419	8			
1	D	284	Total	C	N	O	S	0	0	0
			2156	1376	364	408	8			

- 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	1	0
			50	28	2	20		
2	B	4	Total	C	N	O	3	0
			50	28	2	20		
2	D	4	Total	C	N	O	1	0
			50	28	2	20		

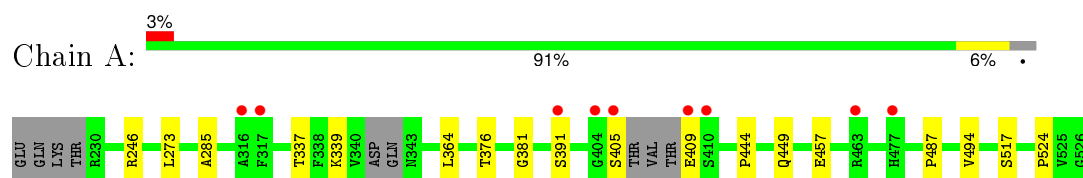
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	228	Total	O	0	0
			228	228		
3	B	258	Total	O	0	0
			258	258		
3	C	250	Total	O	0	0
			250	250		
3	D	237	Total	O	0	0
			237	237		

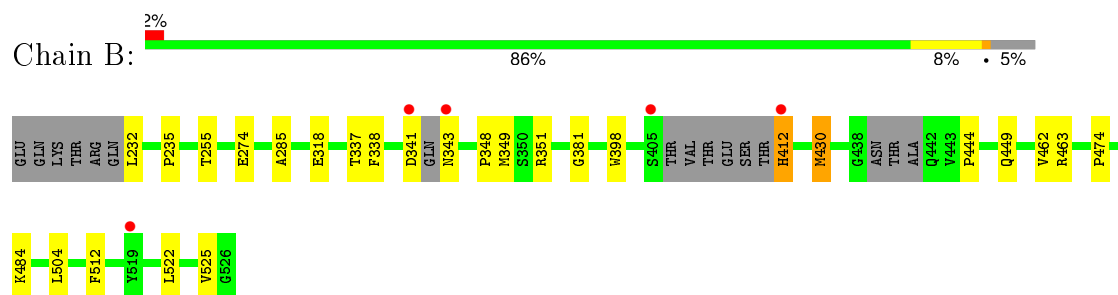
3 Residue-property plots [i](#)

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

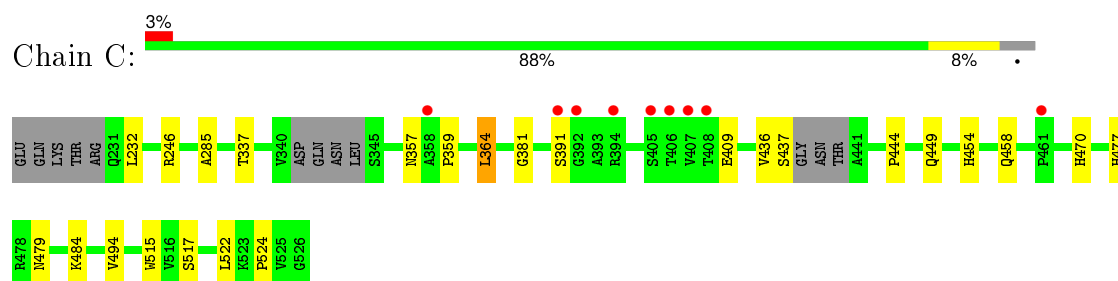
- Molecule 1: P domain of VP1



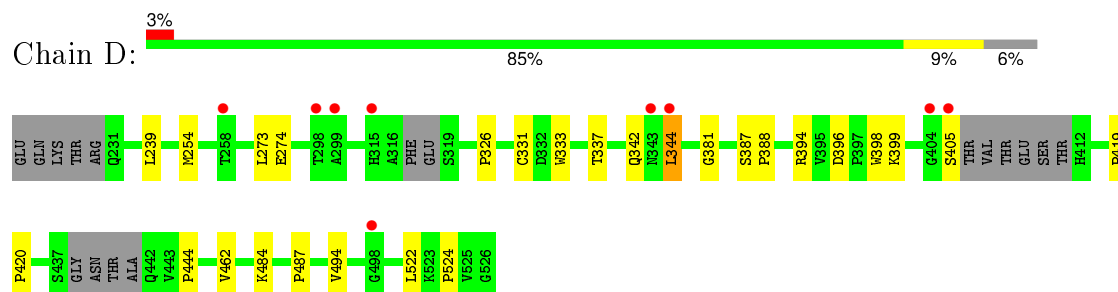
- Molecule 1: P domain of VP1



- Molecule 1: P domain of VP1



- Molecule 1: P domain of VP1



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	62.58 Å 63.26 Å 90.43 Å 99.25° 97.82° 119.62°	Depositor
Resolution (Å)	34.02 – 1.90 34.02 – 1.90	Depositor EDS
% Data completeness (in resolution range)	96.8 (34.02-1.90) 84.2 (34.02-1.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.95 (at 1.89 Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.1_1168)	Depositor
R, R_{free}	0.160 , 0.189 0.166 , 0.195	Depositor DCC
R_{free} test set	4403 reflections (5.47%)	DCC
Wilson B-factor (Å ²)	14.0	Xtriage
Anisotropy	0.209	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 54.1	EDS
Estimated twinning fraction	0.009 for -k,-h,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 87902 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	9866	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.39% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: FUC, GAL, NAG, A2G

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	0.34	0/2291	0.53	0/3132
1	B	0.32	0/2226	0.53	1/3044 (0.0%)
1	C	0.33	0/2263	0.55	0/3096
1	D	0.36	0/2219	0.53	0/3035
All	All	0.34	0/8999	0.53	1/12307 (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
1	B	0	1

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	B	430	MET	CG-SD-CE	5.45	108.92	100.20

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	351	ARG	Sidechain

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2226	0	2148	12	0
1	B	2163	0	2080	19	1
1	C	2198	0	2116	17	0
1	D	2156	0	2077	16	0
2	A	50	0	45	13	1
2	B	50	0	45	1	0
2	D	50	0	45	1	0
3	A	228	0	0	0	0
3	B	258	0	0	1	0
3	C	250	0	0	0	0
3	D	237	0	0	1	0
All	All	9866	0	8556	75	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 4.

All (75) 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:A:604:NAG:H61	2:A:604:NAG:O1	1.56	1.05
1:B:348:PRO:HG2	1:B:349:MET:HE3	1.48	0.93
1:D:396:ASP:HB3	1:D:399:LYS:HE2	1.66	0.77
1:C:470:HIS:CD2	1:C:479:ASN:HD21	2.04	0.74
1:A:405:SER:N	1:A:409:GLU:OE1	2.19	0.74
2:A:601:A2G:C1	2:A:602:GAL:O4	2.35	0.72
1:B:255:THR:HG22	1:B:430:MET:HE2	1.74	0.70
2:A:604:NAG:C6	2:A:604:NAG:O1	2.38	0.70
1:A:339:LYS:NZ	1:A:376:THR:OG1	2.25	0.70
1:B:318:GLU:O	1:B:412:HIS:N	2.28	0.67
1:C:454:HIS:NE2	1:C:458:GLN:NE2	2.43	0.66
1:A:246:ARG:HD3	1:A:444:PRO:O	1.96	0.65
1:A:273:LEU:HG	1:A:487:PRO:HA	1.79	0.65
2:A:602:GAL:O3	2:A:603:FUC:C1	2.45	0.64
2:A:602:GAL:H1	2:A:604:NAG:O3	1.97	0.64
1:A:337:THR:HG22	1:A:381:GLY:HA3	1.81	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:337:THR:HG22	1:B:381:GLY:HA3	1.81	0.61
1:D:337:THR:HG22	1:D:381:GLY:HA3	1.83	0.61
1:C:454:HIS:CD2	1:C:458:GLN:HE21	2.18	0.60
2:A:602:GAL:C3	2:A:603:FUC:O2	2.52	0.58
2:A:602:GAL:H3	2:A:603:FUC:O2	2.04	0.56
1:B:504:LEU:HD11	1:B:525:VAL:HG21	1.87	0.56
1:B:232:LEU:N	3:B:946:HOH:O	2.39	0.56
2:A:604:NAG:H83	2:A:604:NAG:O1	2.07	0.55
2:D:601:A2G:H8	3:D:812:HOH:O	2.07	0.54
1:C:470:HIS:CG	1:C:479:ASN:HD21	2.26	0.54
1:B:255:THR:CG2	1:B:430:MET:HE2	2.39	0.52
1:C:357:ASN:OD1	1:C:359:PRO:HD2	2.09	0.52
1:B:232:LEU:HD23	1:B:512:PHE:CD1	2.45	0.52
2:B:602:GAL:O2	2:B:604:NAG:H62	2.11	0.51
1:D:394:ARG:HH11	1:D:394:ARG:HG3	1.77	0.50
1:D:342:GLN:HB3	1:D:344:LEU:CD1	2.42	0.50
1:B:348:PRO:CG	1:B:349:MET:HE3	2.31	0.49
1:C:285:ALA:HA	1:C:449:GLN:HG2	1.94	0.49
1:C:494:VAL:O	1:C:524:PRO:HA	2.12	0.49
1:B:338:PHE:HB3	1:B:349:MET:CE	2.42	0.49
1:C:470:HIS:CG	1:C:479:ASN:ND2	2.81	0.48
2:A:602:GAL:O3	2:A:603:FUC:O2	2.30	0.48
1:B:338:PHE:HB3	1:B:349:MET:HE1	1.94	0.48
1:A:246:ARG:CD	1:A:444:PRO:O	2.62	0.48
1:D:484:LYS:HG3	1:D:522:LEU:HD21	1.94	0.48
2:A:603:FUC:H5	2:A:604:NAG:H62	1.95	0.48
1:A:457:GLU:OE1	1:B:235:PRO:HA	2.14	0.48
1:A:364:LEU:HG	1:A:409:GLU:OE2	2.15	0.47
1:D:494:VAL:O	1:D:524:PRO:HA	2.13	0.47
1:C:357:ASN:CG	1:C:359:PRO:HD2	2.35	0.47
1:B:474:PRO:HA	1:C:477:HIS:CE1	2.50	0.46
1:C:484:LYS:HG3	1:C:522:LEU:HD21	1.97	0.46
1:C:436:VAL:O	1:C:437:SER:HB3	2.16	0.46
1:D:273:LEU:HG	1:D:487:PRO:HA	1.97	0.45
1:D:331:CYS:HB2	1:D:387:SER:O	2.15	0.45
1:D:239:LEU:HD21	1:D:254:MET:HG3	1.98	0.45
2:A:604:NAG:H61	2:A:604:NAG:HO1	1.75	0.44
1:D:419:PRO:HA	1:D:420:PRO:HD3	1.81	0.44
1:D:398:TRP:CE2	1:D:444:PRO:HG3	2.53	0.44
1:C:337:THR:HG22	1:C:381:GLY:HA3	1.99	0.44
1:B:285:ALA:HA	1:B:449:GLN:HG2	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:601:A2G:O	2:A:602:GAL:H4	2.17	0.43
1:B:274:GLU:HG3	1:B:462:VAL:HG23	2.00	0.43
1:C:364:LEU:HD13	1:C:409:GLU:OE2	2.18	0.43
1:D:274:GLU:HG3	1:D:462:VAL:HG23	2.00	0.43
1:C:454:HIS:CD2	1:C:458:GLN:NE2	2.84	0.43
1:A:285:ALA:HA	1:A:449:GLN:HG2	2.01	0.43
1:D:326:PRO:HB2	1:D:333:TRP:CH2	2.53	0.43
1:C:246:ARG:HD3	1:C:444:PRO:O	2.19	0.43
1:A:494:VAL:O	1:A:524:PRO:HA	2.19	0.43
1:D:387:SER:HB2	1:D:388:PRO:CD	2.49	0.43
1:D:274:GLU:CG	1:D:462:VAL:HG23	2.49	0.43
1:A:391:SER:HA	2:A:601:A2G:O6	2.18	0.42
1:B:484:LYS:HG3	1:B:522:LEU:HD21	2.01	0.42
1:C:232:LEU:HD22	1:C:515:TRP:CH2	2.55	0.42
1:D:396:ASP:HB3	1:D:399:LYS:CE	2.45	0.41
1:B:398:TRP:CE2	1:B:444:PRO:HG3	2.55	0.41
1:B:341:ASP:O	1:B:343:ASN:N	2.53	0.41
1:A:285:ALA:HB3	1:B:285:ALA:HB3	2.02	0.40

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:B:463:ARG:NH2	2:A:604:NAG:O6[1_565]	1.92	0.28

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	286/301 (95%)	274 (96%)	12 (4%)	0	100	100
1	B	277/301 (92%)	266 (96%)	11 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	283/301 (94%)	276 (98%)	7 (2%)	0	100	100
1	D	276/301 (92%)	264 (96%)	12 (4%)	0	100	100
All	All	1122/1204 (93%)	1080 (96%)	42 (4%)	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	247/257 (96%)	246 (100%)	1 (0%)	93	94
1	B	240/257 (93%)	239 (100%)	1 (0%)	93	94
1	C	244/257 (95%)	241 (99%)	3 (1%)	78	76
1	D	239/257 (93%)	237 (99%)	2 (1%)	86	86
All	All	970/1028 (94%)	963 (99%)	7 (1%)	88	88

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	517	SER
1	B	412	HIS
1	C	364	LEU
1	C	391	SER
1	C	517	SER
1	D	344	LEU
1	D	405	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	363	HIS
1	A	521	GLN

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Mol	Chain	Res	Type
1	B	343	ASN
1	C	458	GLN
1	C	479	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

12 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	A2G	A	601	2	14,14,15	0.48	0	15,19,21	0.91	1 (6%)
2	GAL	A	602	2	12,12,12	0.48	0	17,17,17	0.77	1 (5%)
2	FUC	A	603	2	10,10,11	0.37	0	14,14,16	0.96	1 (7%)
2	NAG	A	604	2	13,14,15	0.70	0	14,19,21	1.19	2 (14%)
2	A2G	B	601	2	14,14,15	0.61	0	15,19,21	1.16	2 (13%)
2	GAL	B	602	2	12,12,12	0.48	0	17,17,17	0.78	1 (5%)
2	FUC	B	603	2	10,10,11	0.36	0	14,14,16	0.96	1 (7%)
2	NAG	B	604	2	13,14,15	0.70	0	14,19,21	1.18	2 (14%)
2	A2G	D	601	2	14,14,15	0.47	0	15,19,21	0.91	1 (6%)
2	GAL	D	602	2	12,12,12	0.48	0	17,17,17	0.78	1 (5%)
2	FUC	D	603	2	10,10,11	0.36	0	14,14,16	0.96	1 (7%)
2	NAG	D	604	2	13,14,15	0.71	0	14,19,21	1.18	2 (14%)

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	A2G	A	601	2	-	0/6/23/26	0/1/1/1
2	GAL	A	602	2	-	0/2/22/22	0/1/1/1
2	FUC	A	603	2	-	0/0/17/20	0/1/1/1
2	NAG	A	604	2	-	0/6/22/26	1/1/1/1
2	A2G	B	601	2	-	0/6/23/26	0/1/1/1
2	GAL	B	602	2	-	0/2/22/22	0/1/1/1
2	FUC	B	603	2	-	0/0/17/20	0/1/1/1
2	NAG	B	604	2	-	0/6/22/26	0/1/1/1
2	A2G	D	601	2	-	0/6/23/26	0/1/1/1
2	GAL	D	602	2	-	0/2/22/22	0/1/1/1
2	FUC	D	603	2	-	0/0/17/20	0/1/1/1
2	NAG	D	604	2	-	0/6/22/26	0/1/1/1

There are no bond length outliers.

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	601	A2G	C1-O-C5	-2.18	109.49	112.25
2	D	601	A2G	C1-O-C5	-2.17	109.49	112.25
2	B	604	NAG	C1-O5-C5	-2.17	109.48	113.47
2	B	601	A2G	C1-O-C5	-2.16	109.50	112.25
2	D	604	NAG	C1-O5-C5	-2.16	109.49	113.47
2	B	602	GAL	C1-O5-C5	-2.16	109.48	113.47
2	D	602	GAL	C1-O5-C5	-2.15	109.50	113.47
2	A	602	GAL	C1-O5-C5	-2.14	109.51	113.47
2	A	604	NAG	C1-O5-C5	-2.03	109.73	113.47
2	A	603	FUC	O5-C5-C6	2.03	109.48	106.13
2	B	603	FUC	O5-C5-C6	2.03	109.48	106.13
2	D	603	FUC	O5-C5-C6	2.05	109.52	106.13
2	B	604	NAG	O5-C5-C6	2.34	109.49	106.62
2	A	604	NAG	O5-C5-C6	2.35	109.49	106.62
2	D	604	NAG	O5-C5-C6	2.36	109.50	106.62
2	B	601	A2G	O4-C4-C5	2.73	116.48	109.24

There are no chirality outliers.

There are no torsion outliers.

All (1) ring outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	604	NAG	C1-C2-C3-C4-C5-O5

7 monomers are involved in 16 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	601	A2G	3	0
2	A	602	GAL	7	0
2	A	603	FUC	5	0
2	A	604	NAG	6	1
2	B	602	GAL	1	0
2	B	604	NAG	1	0
2	D	601	A2G	1	0

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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

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	292/301 (97%)	-0.13	9 (3%) 52 56	8, 16, 35, 51	0
1	B	285/301 (94%)	-0.23	5 (1%) 71 74	7, 15, 28, 49	0
1	C	289/301 (96%)	-0.20	9 (3%) 52 56	8, 15, 35, 50	0
1	D	284/301 (94%)	-0.07	9 (3%) 51 54	9, 16, 37, 57	0
All	All	1150/1204 (95%)	-0.16	32 (2%) 56 60	7, 15, 35, 57	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	407	VAL	6.8
1	D	405	SER	6.3
1	D	299	ALA	5.3
1	C	358	ALA	4.6
1	B	343	ASN	4.2
1	A	405	SER	4.1
1	D	343	ASN	3.5
1	A	409	GLU	3.4
1	B	341	ASP	3.4
1	D	298	THR	3.3
1	B	519	TYR	3.2
1	D	498	GLY	3.0
1	C	392	GLY	2.9
1	D	344	LEU	2.9
1	A	391	SER	2.8
1	D	404	GLY	2.7
1	C	394	ARG	2.5
1	A	410	SER	2.5
1	C	406	THR	2.5
1	B	412	HIS	2.5
1	A	317	PHE	2.5

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Mol	Chain	Res	Type	RSRZ
1	C	408	THR	2.5
1	A	477	HIS	2.3
1	C	461	PRO	2.3
1	C	391	SER	2.2
1	B	405	SER	2.2
1	A	316	ALA	2.2
1	D	258	THR	2.1
1	A	404	GLY	2.0
1	A	463	ARG	2.0
1	C	405	SER	2.0
1	D	315	HIS	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](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	A2G	B	601	14/15	0.84	0.17	5.06	13,17,18,19	0
2	A2G	A	601	14/15	0.81	0.20	1.65	18,23,29,32	0
2	A2G	D	601	14/15	0.84	0.16	1.39	15,18,20,21	0
2	NAG	A	604	14/15	0.38	0.58	-	66,70,76,77	8
2	GAL	B	602	12/12	0.81	0.21	-	22,31,38,40	0
2	NAG	D	604	14/15	0.77	0.28	-	46,49,54,54	9
2	GAL	D	602	12/12	0.83	0.19	-	26,38,44,46	0
2	FUC	B	603	10/11	0.90	0.11	-	21,21,22,22	0
2	NAG	B	604	14/15	0.69	0.27	-	36,39,48,48	7
2	FUC	D	603	10/11	0.86	0.15	-	29,30,31,32	0
2	GAL	A	602	12/12	0.75	0.34	-	41,56,59,64	0
2	FUC	A	603	10/11	0.65	0.27	-	59,60,61,62	0

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.