Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 10:11 PM GMT

PDB ID	:	4X06
Title	:	Crystal structure of P domain from norovirus strain Saga4 in complex with HBGA type B (triglycan)
Authors	:	Singh, B.K.; Hansman, G.S.
Deposited on	:	2014-11-20
Resolution	:	1.22 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

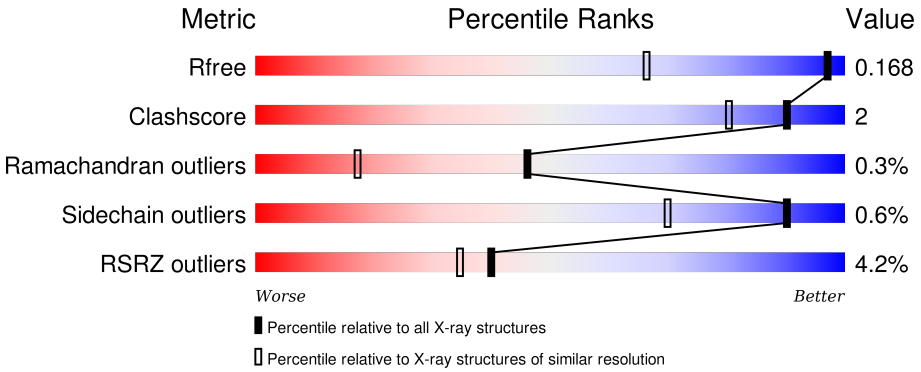
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)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

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.22 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	1197 (1.26-1.18)
Clashscore	102246	1295 (1.26-1.18)
Ramachandran outliers	100387	1239 (1.26-1.18)
Sidechain outliers	100360	1237 (1.26-1.18)
RSRZ outliers	91569	1201 (1.26-1.18)

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 <=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	309	4% 88% 10%
1	B	309	4% 89% 10%

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	FUC	A	601[A]	-	-	-	X
2	FUC	A	601[B]	-	-	-	X
2	FUC	B	601[A]	-	-	-	X
2	FUC	B	601[B]	-	-	-	X
3	ACT	A	604	-	-	-	X
3	ACT	B	604	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 10037 atoms, of which 4316 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 VP1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	309	Total	C	H	N	O	S	0	8	0
			4605	1547	2168	417	463	10			
1	B	309	Total	C	H	N	O	S	0	9	0
			4576	1547	2148	409	462	10			

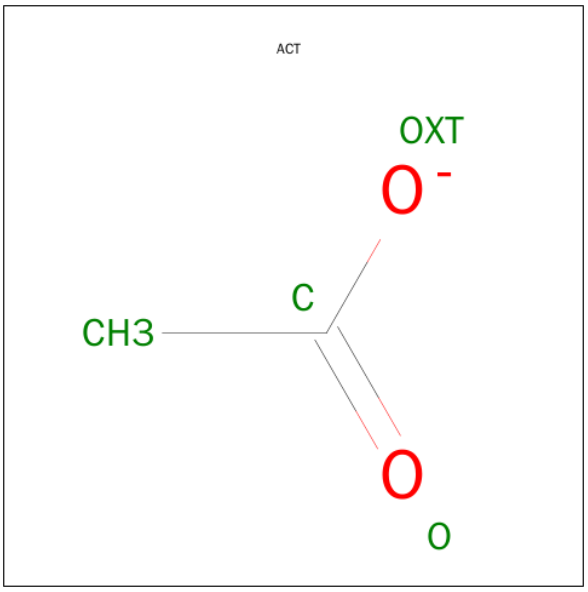
There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	222	THR	-	expression tag	UNP B5BTR7
A	223	GLY	-	expression tag	UNP B5BTR7
A	224	SER	-	expression tag	UNP B5BTR7
B	222	THR	-	expression tag	UNP B5BTR7
B	223	GLY	-	expression tag	UNP B5BTR7
B	224	SER	-	expression tag	UNP B5BTR7

- Molecule 2 is a polymer of unknown type called SUGAR (3-MER).

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	3	Total	C	O	0	3
			66	36	30		
2	B	3	Total	C	O	0	3
			66	36	30		

- Molecule 3 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).



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

- Molecule 4 is water.

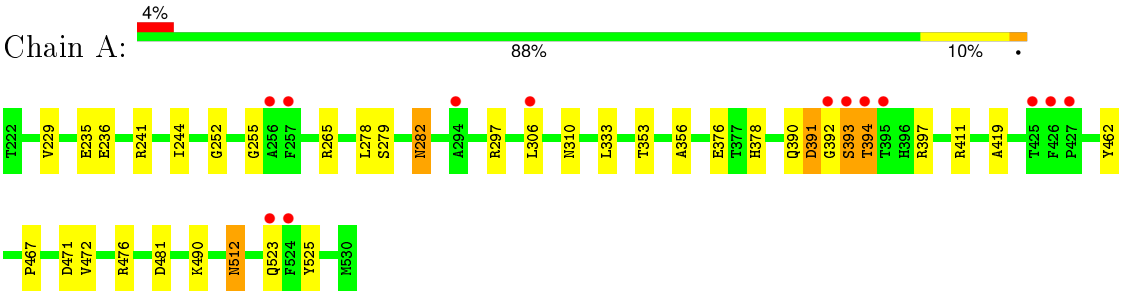
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	358	Total O 358 358	0	0
4	B	358	Total O 358 358	0	0

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.

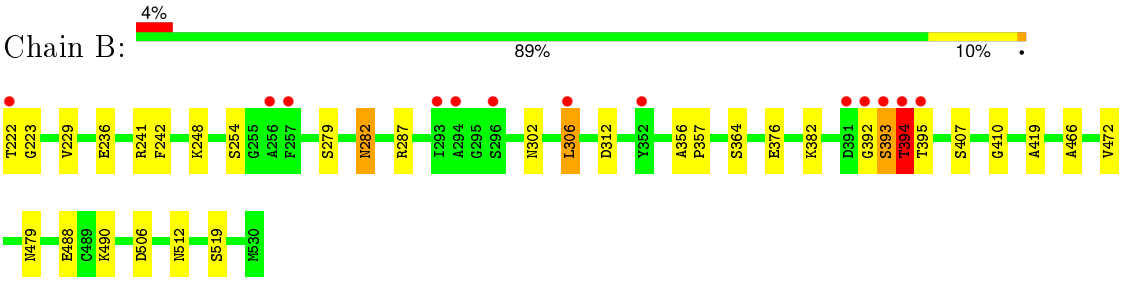
- Molecule 1: VP1

Chain A:



- Molecule 1: VP1

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	113.83Å 58.65Å 97.31Å 90.00° 107.34° 90.00°	Depositor
Resolution (Å)	32.32 – 1.22 48.10 – 1.22	Depositor EDS
% Data completeness (in resolution range)	96.7 (32.32-1.22) 96.7 (48.10-1.22)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.73 (at 1.22Å)	Xtriage
Refinement program	phenix	Depositor
R, R_{free}	0.134 , 0.161 0.149 , 0.168	Depositor DCC
R_{free} test set	8823 reflections (5.26%)	DCC
Wilson B-factor (Å^2)	13.0	Xtriage
Anisotropy	0.249	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.41 , 49.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Outliers	0 of 176472 reflections	Xtriage
F_o,F_c correlation	0.98	EDS
Total number of atoms	10037	wwPDB-VP
Average B, all atoms (Å^2)	17.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 76.14 % of the origin peak, indicating pseudo translational symmetry.The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 1.1028e-06.The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: GLA, FUC, ACT

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	1.41	12/2515 (0.5%)	1.09	19/3442 (0.6%)
1	B	1.42	16/2510 (0.6%)	0.95	7/3438 (0.2%)
All	All	1.41	28/5025 (0.6%)	1.02	26/6880 (0.4%)

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	A	0	4
1	B	0	3
All	All	0	7

All (28) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	236	GLU	CD-OE1	-6.62	1.18	1.25
1	B	302[A]	ASN	CA-C	6.25	1.69	1.52
1	B	302[B]	ASN	CA-C	6.25	1.69	1.52
1	B	519	SER	CB-OG	-6.12	1.34	1.42
1	B	254	SER	CB-OG	-6.10	1.34	1.42
1	B	488	GLU	CD-OE1	-6.09	1.19	1.25
1	B	229	VAL	C-O	-5.94	1.12	1.23
1	A	356	ALA	C-O	-5.86	1.12	1.23
1	B	242	PHE	C-O	-5.80	1.12	1.23
1	B	407[A]	SER	CA-C	5.73	1.67	1.52
1	B	407[B]	SER	CA-C	5.73	1.67	1.52
1	A	252	GLY	C-O	-5.57	1.14	1.23
1	A	236	GLU	CD-OE1	-5.52	1.19	1.25
1	A	279	SER	C-O	-5.51	1.12	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	419	ALA	C-O	-5.43	1.13	1.23
1	A	419	ALA	C-O	-5.41	1.13	1.23
1	A	376	GLU	CD-OE1	-5.39	1.19	1.25
1	A	229	VAL	C-O	-5.33	1.13	1.23
1	B	376	GLU	CD-OE1	-5.29	1.19	1.25
1	A	244	ILE	C-O	-5.26	1.13	1.23
1	B	466	ALA	C-O	-5.24	1.13	1.23
1	B	410	GLY	C-O	-5.18	1.15	1.23
1	B	512	ASN	C-O	-5.16	1.13	1.23
1	A	353	THR	C-O	-5.15	1.13	1.23
1	A	525	TYR	CE1-CZ	-5.08	1.31	1.38
1	A	467	PRO	C-O	-5.08	1.13	1.23
1	B	364	SER	CB-OG	-5.06	1.35	1.42
1	A	255	GLY	C-O	-5.02	1.15	1.23

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	397	ARG	NE-CZ-NH2	-12.95	113.82	120.30
1	A	397	ARG	NE-CZ-NH1	12.70	126.65	120.30
1	A	394[A]	THR	N-CA-C	9.09	135.54	111.00
1	A	394[B]	THR	N-CA-C	9.09	135.54	111.00
1	B	241	ARG	NE-CZ-NH2	-8.12	116.24	120.30
1	A	297	ARG	NE-CZ-NH2	-8.07	116.27	120.30
1	B	306	LEU	CB-CG-CD2	7.62	123.94	111.00
1	A	393[A]	SER	N-CA-C	6.46	128.46	111.00
1	A	393[B]	SER	N-CA-C	6.46	128.46	111.00
1	A	411	ARG	NE-CZ-NH2	-6.23	117.19	120.30
1	B	312	ASP	CB-CG-OD1	6.04	123.73	118.30
1	B	312	ASP	CB-CG-OD2	-5.98	112.92	118.30
1	A	265	ARG	NE-CZ-NH1	5.68	123.14	120.30
1	A	392[A]	GLY	N-CA-C	-5.61	99.07	113.10
1	A	392[B]	GLY	N-CA-C	-5.61	99.07	113.10
1	A	512[A]	ASN	CB-CA-C	-5.59	99.22	110.40
1	A	512[B]	ASN	CB-CA-C	-5.59	99.22	110.40
1	B	287	ARG	NE-CZ-NH2	-5.59	117.51	120.30
1	A	241	ARG	NE-CZ-NH2	-5.55	117.52	120.30
1	B	306	LEU	CB-CG-CD1	5.53	120.40	111.00
1	A	391[A]	ASP	CB-CG-OD1	5.43	123.18	118.30
1	A	391[B]	ASP	CB-CG-OD1	5.43	123.18	118.30
1	A	512[A]	ASN	CA-C-O	5.22	131.06	120.10
1	A	512[B]	ASN	CA-C-O	5.22	131.06	120.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	476	ARG	NE-CZ-NH2	-5.11	117.74	120.30
1	B	506	ASP	CB-CG-OD1	5.03	122.83	118.30

There are no chirality outliers.

All (7) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	282[B]	ASN	Mainchain
1	A	390	GLN	Peptide
1	A	393[A]	SER	Peptide
1	A	394[A]	THR	Mainchain
1	B	282[A]	ASN	Mainchain
1	B	282[B]	ASN	Mainchain
1	B	394[A]	THR	Mainchain

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	2437	2168	2327	10	1
1	B	2428	2148	2312	11	1
2	A	66	0	60	1	0
2	B	66	0	60	0	0
3	A	4	0	3	0	0
3	B	4	0	3	0	0
4	A	358	0	0	3	0
4	B	358	0	0	4	1
All	All	5721	4316	4765	22	2

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 2.

All (22) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:282[B]:ASN:OD1	4:A:1028:HOH:O	1.89	0.90

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:310:ASN:ND2	4:A:701:HOH:O	2.17	0.73
1:A:282[B]:ASN:HD21	1:A:306:LEU:HD13	1.56	0.71
2:A:603[B]:GLA:H2	4:A:713:HOH:O	1.91	0.69
1:B:223:GLY:HA2	4:B:771:HOH:O	2.01	0.60
1:B:248:LYS:NZ	4:B:1024:HOH:O	2.36	0.59
1:A:481:ASP:OD1	1:A:512[B]:ASN:ND2	2.37	0.56
1:A:282[B]:ASN:HD21	1:A:306:LEU:CD1	2.19	0.56
1:A:282[B]:ASN:ND2	1:A:306:LEU:HD13	2.23	0.53
1:A:278:LEU:HD11	1:A:462:TYR:CE2	2.46	0.51
1:A:471:ASP:OD2	1:A:523:GLN:OE1	2.30	0.49
1:B:279:SER:HB3	1:B:282[B]:ASN:HB2	1.93	0.49
1:B:356:ALA:HB3	1:B:357:PRO:HD3	1.95	0.48
1:B:282[B]:ASN:HD21	1:B:306:LEU:HD13	1.80	0.47
1:B:222:THR:N	4:B:702:HOH:O	2.50	0.44
1:B:382:LYS:HD2	4:B:1023:HOH:O	2.19	0.43
1:B:356:ALA:N	1:B:357:PRO:CD	2.82	0.42
1:A:235[B]:GLU:CD	1:A:235[B]:GLU:H	2.22	0.42
1:B:472:VAL:HG11	1:B:490:LYS:HG2	2.02	0.41
1:A:472:VAL:HG11	1:A:490:LYS:HG2	2.02	0.41
1:B:392[B]:GLY:O	1:B:393[B]:SER:C	2.59	0.41
1:B:393[B]:SER:O	1:B:394[B]:THR:CB	2.70	0.40

All (2) 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:A:378:HIS:HB3	1:B:479:ASN:HD21[4_955]	1.31	0.29
4:B:749:HOH:O	4:B:755:HOH:O[2_1056]	2.15	0.05

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	315/309 (102%)	309 (98%)	6 (2%)	0	100 100
1	B	316/309 (102%)	307 (97%)	5 (2%)	4 (1%)	15 1
All	All	631/618 (102%)	616 (98%)	11 (2%)	4 (1%)	46 6

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	393[A]	SER
1	B	393[B]	SER
1	B	394[A]	THR
1	B	394[B]	THR

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	270/268 (101%)	267 (99%)	3 (1%)	80 49
1	B	268/268 (100%)	267 (100%)	1 (0%)	93 79
All	All	538/536 (100%)	534 (99%)	4 (1%)	90 66

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

Mol	Chain	Res	Type
1	A	333	LEU
1	A	391[A]	ASP
1	A	391[B]	ASP
1	B	395	THR

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

5.3.3 RNA (i)

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)

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	FUC	A	601[A]	2	10,10,11	0.74	0	14,14,16	0.94	1 (7%)
2	FUC	A	601[B]	2	10,10,11	0.29	0	14,14,16	0.79	1 (7%)
2	GLA	A	602[A]	2	12,12,12	0.42	0	17,17,17	0.95	2 (11%)
2	GLA	A	602[B]	2	12,12,12	0.60	0	17,17,17	0.95	1 (5%)
2	GLA	A	603[A]	2	11,11,12	0.72	0	14,15,17	1.00	1 (7%)
2	GLA	A	603[B]	2	11,11,12	0.81	0	14,15,17	1.20	1 (7%)
2	FUC	B	601[A]	2	10,10,11	0.75	0	14,14,16	0.92	1 (7%)
2	FUC	B	601[B]	2	10,10,11	0.30	0	14,14,16	0.79	1 (7%)
2	GLA	B	602[A]	2	12,12,12	0.58	0	17,17,17	0.97	1 (5%)
2	GLA	B	602[B]	2	12,12,12	0.59	0	17,17,17	0.97	1 (5%)
2	GLA	B	603[A]	2	11,11,12	0.72	0	14,15,17	1.19	1 (7%)
2	GLA	B	603[B]	2	11,11,12	0.83	0	14,15,17	1.23	1 (7%)

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	FUC	A	601[A]	2	-	0/0/17/20	0/1/1/1
2	FUC	A	601[B]	2	-	0/0/17/20	0/1/1/1
2	GLA	A	602[A]	2	-	0/2/22/22	0/1/1/1
2	GLA	A	602[B]	2	-	0/2/22/22	0/1/1/1
2	GLA	A	603[A]	2	-	0/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GLA	A	603[B]	2	-	0/2/19/22	0/1/1/1
2	FUC	B	601[A]	2	-	0/0/17/20	0/1/1/1
2	FUC	B	601[B]	2	-	0/0/17/20	0/1/1/1
2	GLA	B	602[A]	2	-	0/2/22/22	0/1/1/1
2	GLA	B	602[B]	2	-	0/2/22/22	0/1/1/1
2	GLA	B	603[A]	2	-	0/2/19/22	0/1/1/1
2	GLA	B	603[B]	2	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	603[B]	GLA	C1-O5-C5	-3.83	107.39	112.25
2	A	603[B]	GLA	C1-O5-C5	-3.44	107.88	112.25
2	B	603[A]	GLA	C1-O5-C5	-3.35	108.00	112.25
2	A	602[B]	GLA	C1-O5-C5	-3.15	107.65	113.47
2	B	602[B]	GLA	C1-O5-C5	-3.13	107.68	113.47
2	A	603[A]	GLA	C2-C3-C4	-2.53	106.74	111.04
2	A	602[A]	GLA	O3-C3-C2	-2.10	105.62	110.34
2	B	602[A]	GLA	C6-C5-C4	-2.09	107.85	113.02
2	B	601[B]	FUC	O5-C5-C6	2.01	109.45	106.13
2	A	601[B]	FUC	O5-C5-C6	2.02	109.47	106.13
2	A	602[A]	GLA	C1-C2-C3	2.13	113.60	110.43
2	B	601[A]	FUC	C1-O5-C5	2.20	115.78	112.38
2	A	601[A]	FUC	C1-O5-C5	2.32	115.95	112.38

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	603[B]	GLA	1	0

5.6 Ligand geometry (i)

2 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and

the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
3	ACT	A	604	-	1,3,3	1.82	0	0,3,3	0.00	-
3	ACT	B	604	-	1,3,3	1.85	0	0,3,3	0.00	-

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	ACT	A	604	-	-	0/0/0/0	0/0/0/0
3	ACT	B	604	-	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers (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, 95^{th} 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	309/309 (100%)	-0.26	13 (4%)	40	35	8, 15, 27, 36	0
1	B	309/309 (100%)	-0.16	13 (4%)	40	35	7, 14, 28, 44	0
All	All	618/618 (100%)	-0.21	26 (4%)	40	35	7, 14, 27, 44	0

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	394[A]	THR	8.1
1	B	393[A]	SER	7.9
1	B	306	LEU	6.3
1	A	393[A]	SER	5.9
1	A	394[A]	THR	5.0
1	A	524	PHE	4.1
1	B	392[A]	GLY	4.1
1	B	222	THR	3.8
1	B	293	ILE	3.5
1	A	425	THR	3.5
1	B	395	THR	3.4
1	A	306	LEU	3.2
1	A	426	PHE	3.1
1	B	257	PHE	3.0
1	A	256	ALA	2.8
1	A	257	PHE	2.8
1	B	352[A]	TYR	2.7
1	B	256	ALA	2.5
1	A	427	PRO	2.5
1	A	395	THR	2.4
1	B	294	ALA	2.3
1	A	392[A]	GLY	2.3
1	A	523	GLN	2.3
1	B	391[A]	ASP	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	296	SER	2.1
1	A	294	ALA	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, 95^{th} 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	FUC	A	601[A]	10/11	0.95	0.09	7.31	14,16,17,28	10
2	FUC	A	601[B]	10/11	0.95	0.09	7.31	13,14,15,16	10
2	FUC	B	601[B]	10/11	0.90	0.10	4.13	17,18,20,23	10
2	FUC	B	601[A]	10/11	0.90	0.10	4.13	20,23,26,29	10
2	GLA	A	602[B]	12/12	0.90	0.14	-0.35	22,26,38,44	12
2	GLA	A	602[A]	12/12	0.90	0.14	-0.47	17,26,39,46	12
2	GLA	A	603[B]	11/12	0.87	0.14	-	23,25,29,33	11
2	GLA	B	603[A]	11/12	0.80	0.15	-	20,34,38,41	11
2	GLA	A	603[A]	11/12	0.87	0.14	-	17,21,28,33	11
2	GLA	B	602[B]	12/12	0.84	0.14	-	35,39,44,47	12
2	GLA	B	602[A]	12/12	0.84	0.14	-	23,39,44,50	12
2	GLA	B	603[B]	11/12	0.80	0.15	-	28,33,40,41	11

6.4 Ligands (i)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95^{th} 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
3	ACT	A	604	4/4	0.90	0.13	6.98	23,27,28,31	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å^2)	Q<0.9
3	ACT	B	604	4/4	0.89	0.10	4.04	24,26,30,32	0

6.5 Other polymers (i)

There are no such residues in this entry.