



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 1, 2016 – 01:34 PM GMT

PDB ID : 3U2S
Title : Crystal Structure of PG9 Fab in Complex with V1V2 Region from HIV-1 strain ZM109
Authors : McLellan, J.S.; Pancera, M.; Kwong, P.D.
Deposited on : 2011-10-04
Resolution : 1.80 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
<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

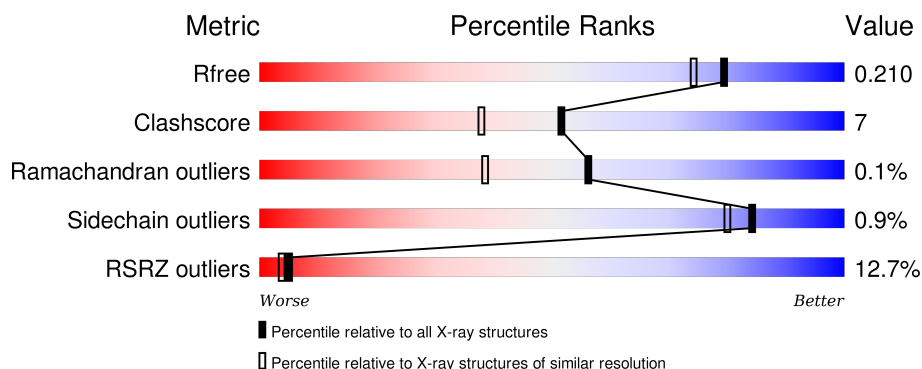
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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	4533 (1.80-1.80)
Clashscore	102246	5383 (1.80-1.80)
Ramachandran outliers	100387	5320 (1.80-1.80)
Sidechain outliers	100360	5319 (1.80-1.80)
RSRZ outliers	91569	4547 (1.80-1.80)

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	248	<div> <div>19%</div> <div> <div></div> <div>83%</div> <div>8%</div> <div>8%</div> </div> </div>
1	H	248	<div> <div>2%</div> <div> <div></div> <div>84%</div> <div>8%</div> <div>7%</div> </div> </div>
2	B	216	<div> <div>4%</div> <div> <div></div> <div>87%</div> <div>11%</div> <div>•</div> </div> </div>
2	L	216	<div> <div>3%</div> <div> <div></div> <div>88%</div> <div>9%</div> <div>•</div> </div> </div>
3	C	124	<div> <div>29%</div> <div> <div></div> <div>56%</div> <div>16%</div> <div>•</div> <div>27%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
3	G	124	

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	SO4	B	215	-	-	-	X
4	SO4	B	216	-	-	-	X
4	SO4	L	214	-	-	-	X

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 16931 atoms, of which 7771 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 PG9 heavy chain.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	H	230	Total	C	H	N	O	S	0	0	0
			3447	1122	1667	298	351	9			
1	A	227	Total	C	H	N	O	S	0	0	0
			3415	1113	1652	295	346	9			

- Molecule 2 is a protein called PG9 light chain.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	L	211	Total	C	H	N	O	S	0	0	0
			3103	974	1537	267	321	4			
2	B	211	Total	C	H	N	O	S	0	0	0
			3103	974	1537	267	321	4			

- Molecule 3 is a protein called Envelope glycoprotein gp120.

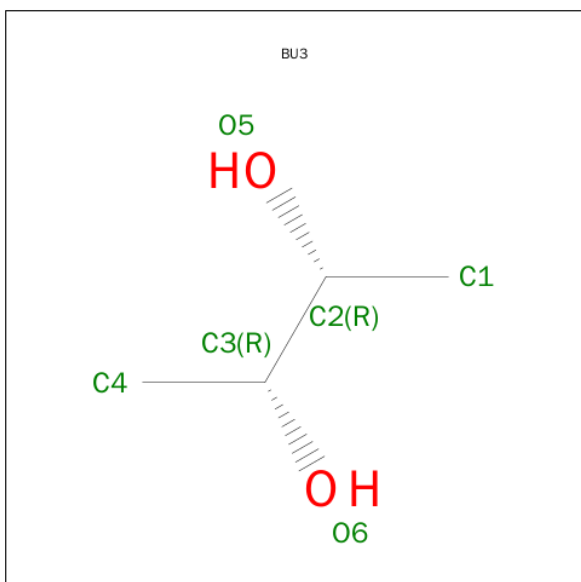
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
3	G	90	Total	C	H	N	O	S	0	0	0
			1379	440	677	117	141	4			
3	C	90	Total	C	H	N	O	S	0	0	0
			1386	442	681	117	142	4			

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	H	1	Total	O	S	0	0
			5	4	1		
4	L	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 5 is (R,R)-2,3-BUTANEDIOL (three-letter code: BU3) (formula: C₄H₁₀O₂).

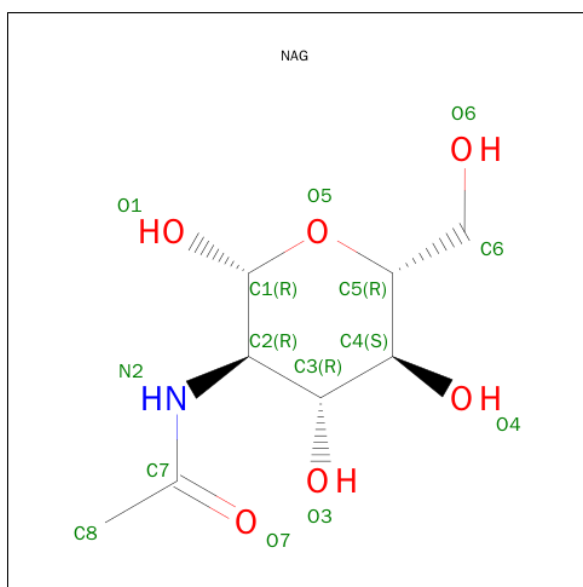


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	L	1	Total	C	H	O	0	0
			16	4	10	2		
5	B	1	Total	C	H	O	0	0
			16	4	10	2		

- Molecule 6 is a polymer of unknown type called SUGAR (7-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	G	7	Total	C	N	O	0	0
			83	46	2	35		
6	C	7	Total	C	N	O	0	0
			83	46	2	35		

- Molecule 7 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	G	1	Total	C	N	O	0	0
			14	8	1	5		
7	C	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	H	232	Total	O	0	0
			232	232		

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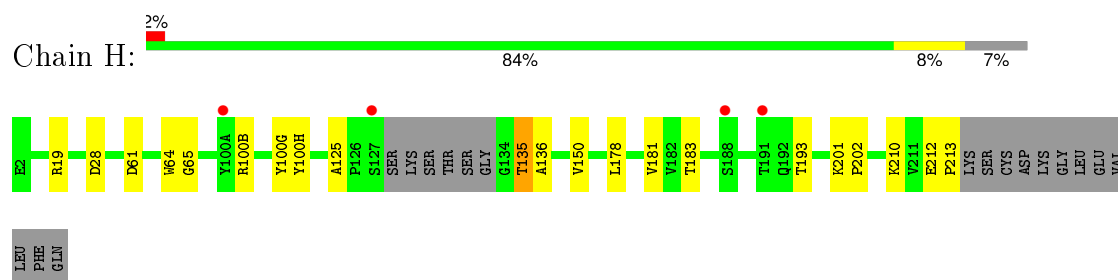
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	L	186	Total 186	O 186	0	0
8	G	38	Total 38	O 38	0	0
8	A	203	Total 203	O 203	0	0
8	B	154	Total 154	O 154	0	0
8	C	34	Total 34	O 34	0	0

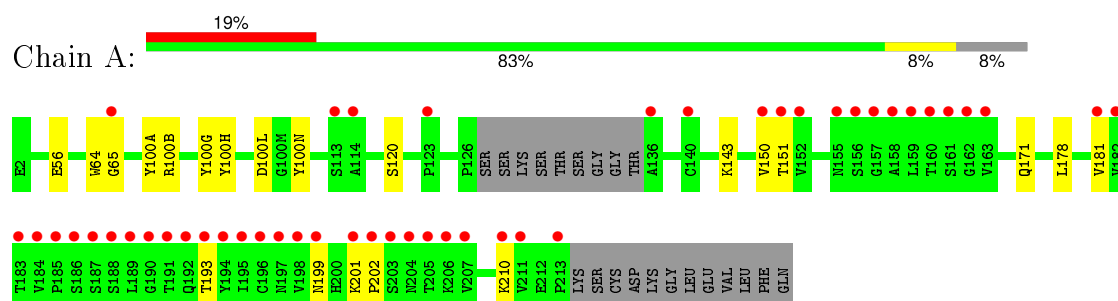
3 Residue-property plots [i](#)

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.

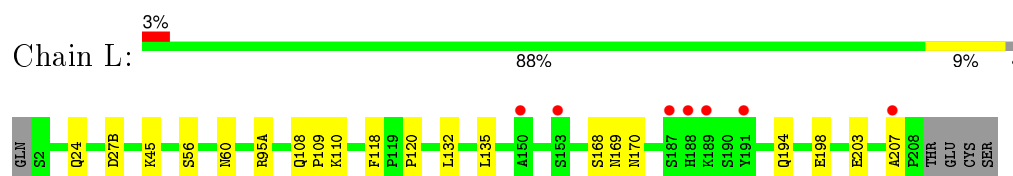
- Molecule 1: PG9 heavy chain



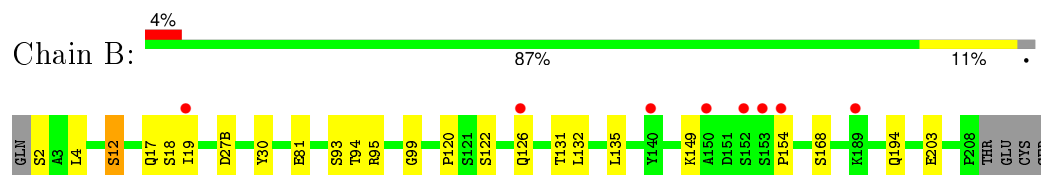
- Molecule 1: PG9 heavy chain



- Molecule 2: PG9 light chain

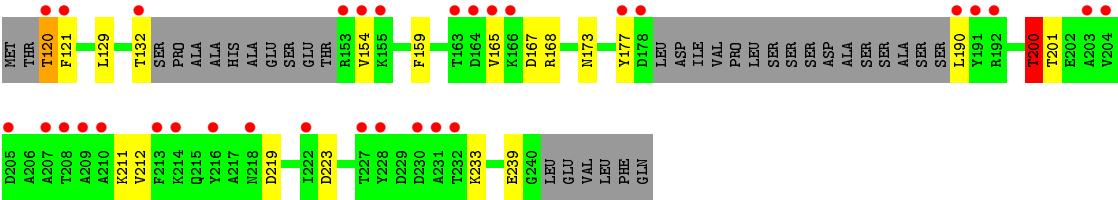


- Molecule 2: PG9 light chain

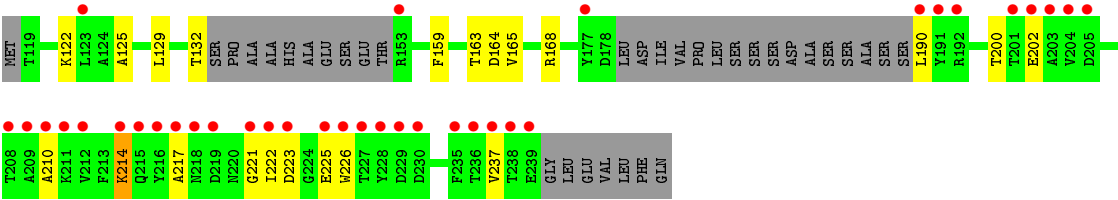


- Molecule 3: Envelope glycoprotein gp120





● Molecule 3: Envelope glycoprotein gp120



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	89.47Å 86.56Å 94.88Å 90.00° 92.08° 90.00°	Depositor
Resolution (Å)	41.58 – 1.80 41.58 – 1.80	Depositor EDS
% Data completeness (in resolution range)	91.0 (41.58-1.80) 90.9 (41.58-1.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.11 (at 1.79Å)	Xtriage
Refinement program	PHENIX (phenix.refine: dev_755)	Depositor
R, R_{free}	0.178 , 0.205 0.182 , 0.210	Depositor DCC
R_{free} test set	6169 reflections (5.31%)	DCC
Wilson B-factor (Å ²)	23.3	Xtriage
Anisotropy	0.533	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 49.6	EDS
Estimated twinning fraction	0.011 for k,h,-l 0.007 for -k,-h,-l 0.025 for h,-k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 122349 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	16931	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.33% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: BMA, NAG, BU3, TYS, SO4, PCA, MAN

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	0/1770	0.53	0/2409
1	H	0.39	0/1787	0.54	0/2432
2	B	0.35	0/1601	0.52	0/2180
2	L	0.37	0/1601	0.53	0/2180
3	C	0.27	0/714	0.44	0/967
3	G	0.28	0/711	0.52	0/962
All	All	0.36	0/8184	0.52	0/11130

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
3	G	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	G	200	THR	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	1763	1652	1662	21	0
1	H	1780	1667	1677	21	0
2	B	1566	1537	1533	18	0
2	L	1566	1537	1533	22	0
3	C	705	681	680	20	0
3	G	702	677	676	17	0
4	B	15	0	0	0	0
4	H	5	0	0	0	0
4	L	5	0	0	1	0
5	B	6	10	10	0	0
5	L	6	10	10	0	0
6	C	83	0	70	0	0
6	G	83	0	70	0	0
7	C	14	0	13	0	0
7	G	14	0	13	0	0
8	A	203	0	0	7	0
8	B	154	0	0	5	0
8	C	34	0	0	2	0
8	G	38	0	0	6	0
8	H	232	0	0	5	1
8	L	186	0	0	10	1
All	All	9160	7771	7947	110	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 7.

The worst 5 of 110 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:H:181:VAL:HG11	2:L:135:LEU:HD13	1.42	1.02
3:G:173:ASN:OD1	8:G:456:HOH:O	1.82	0.96
2:L:56:SER:HG	1:A:100(A):TYR:HH	1.20	0.89
1:A:171:GLN:NE2	8:A:377:HOH:O	2.10	0.84
2:L:207:ALA:O	8:L:668:HOH:O	1.97	0.82

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)
8:H:338:HOH:O	8:L:277:HOH:O[2_645]	2.19	0.01

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	221/248 (89%)	217 (98%)	4 (2%)	0	100	100
1	H	224/248 (90%)	221 (99%)	3 (1%)	0	100	100
2	B	209/216 (97%)	204 (98%)	5 (2%)	0	100	100
2	L	209/216 (97%)	203 (97%)	6 (3%)	0	100	100
3	C	84/124 (68%)	82 (98%)	1 (1%)	1 (1%)	16	4
3	G	84/124 (68%)	82 (98%)	2 (2%)	0	100	100
All	All	1031/1176 (88%)	1009 (98%)	21 (2%)	1 (0%)	56	38

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	C	163	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	189/207 (91%)	188 (100%)	1 (0%)	92	91
1	H	191/207 (92%)	190 (100%)	1 (0%)	92	91

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	B	178/183 (97%)	176 (99%)	2 (1%)	80	74
2	L	178/183 (97%)	177 (99%)	1 (1%)	90	88
3	C	77/105 (73%)	76 (99%)	1 (1%)	76	68
3	G	76/105 (72%)	74 (97%)	2 (3%)	54	37
All	All	889/990 (90%)	881 (99%)	8 (1%)	84	80

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

Mol	Chain	Res	Type
3	G	200	THR
3	C	214	LYS
2	B	12	SER
3	G	120	THR
1	A	56	GLU

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

Mol	Chain	Res	Type
2	L	188	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

6 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
1	TYS	A	100(G)	-	15,16,17	1.12	1 (6%)	16,22,24	0.90	2 (12%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	TYS	A	100(H)	1	15,16,17	0.95	1 (6%)	16,22,24	0.69	0
1	PCA	A	2	1	7,8,9	1.82	2 (28%)	9,10,12	1.67	3 (33%)
1	TYS	H	100(G)	-	15,16,17	1.05	1 (6%)	16,22,24	0.92	1 (6%)
1	TYS	H	100(H)	1	15,16,17	1.02	2 (13%)	16,22,24	0.63	0
1	PCA	H	2	1	7,8,9	1.80	2 (28%)	9,10,12	1.95	5 (55%)

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
1	TYS	A	100(G)	-	-	0/9/11/13	0/1/1/1
1	TYS	A	100(H)	1	-	0/9/11/13	0/1/1/1
1	PCA	A	2	1	-	0/0/11/13	0/1/1/1
1	TYS	H	100(G)	-	-	0/9/11/13	0/1/1/1
1	TYS	H	100(H)	1	-	0/9/11/13	0/1/1/1
1	PCA	H	2	1	-	0/0/11/13	0/1/1/1

The worst 5 of 9 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	100(G)	TYS	OH-CZ	-2.13	1.39	1.42
1	H	100(H)	TYS	OH-CZ	-2.01	1.39	1.42
1	H	100(H)	TYS	O1-S	2.01	1.52	1.45
1	A	100(H)	TYS	O2-S	2.01	1.52	1.45
1	H	100(G)	TYS	O1-S	2.06	1.52	1.45

The worst 5 of 11 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	2	PCA	CA-N-CD	-3.14	103.27	113.81
1	A	2	PCA	CA-N-CD	-2.76	104.55	113.81
1	H	2	PCA	OE-CD-CG	-2.34	121.58	126.81
1	H	2	PCA	CB-CA-C	-2.23	109.71	112.76
1	A	2	PCA	O-C-CA	-2.22	119.58	125.44

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	100(G)	TYS	1	0
1	A	100(H)	TYS	2	0
1	H	100(G)	TYS	1	0
1	H	100(H)	TYS	2	0

5.5 Carbohydrates [i](#)

14 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	$\# Z > 2$
6	NAG	C	560	3,6	14,14,15	0.49	0	15,19,21	0.67	0
6	NAG	C	561	6	14,14,15	0.73	0	15,19,21	0.72	0
6	BMA	C	562	6	11,11,12	0.58	0	14,15,17	1.07	1 (7%)
6	MAN	C	563	6	11,11,12	0.62	0	14,15,17	0.71	0
6	MAN	C	564	6	11,11,12	0.74	0	14,15,17	0.93	0
6	MAN	C	565	6	11,11,12	0.60	0	14,15,17	1.13	1 (7%)
6	MAN	C	566	6	11,11,12	0.67	0	14,15,17	0.72	0
6	NAG	G	560	3,6	14,14,15	0.59	0	15,19,21	1.03	2 (13%)
6	NAG	G	561	6	14,14,15	0.62	0	15,19,21	0.90	0
6	BMA	G	562	6	11,11,12	0.74	0	14,15,17	1.08	1 (7%)
6	MAN	G	563	6	11,11,12	0.62	0	14,15,17	0.83	0
6	MAN	G	564	6	11,11,12	0.62	0	14,15,17	1.07	0
6	MAN	G	565	6	11,11,12	0.63	0	14,15,17	1.06	1 (7%)
6	MAN	G	566	6	11,11,12	0.70	0	14,15,17	0.55	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
6	NAG	C	560	3,6	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	NAG	C	561	6	-	0/6/23/26	0/1/1/1
6	BMA	C	562	6	-	0/2/19/22	0/1/1/1
6	MAN	C	563	6	-	0/2/19/22	0/1/1/1
6	MAN	C	564	6	-	0/2/19/22	0/1/1/1
6	MAN	C	565	6	-	0/2/19/22	0/1/1/1
6	MAN	C	566	6	-	0/2/19/22	0/1/1/1
6	NAG	G	560	3,6	-	0/6/23/26	0/1/1/1
6	NAG	G	561	6	-	0/6/23/26	0/1/1/1
6	BMA	G	562	6	-	0/2/19/22	0/1/1/1
6	MAN	G	563	6	-	0/2/19/22	0/1/1/1
6	MAN	G	564	6	-	0/2/19/22	0/1/1/1
6	MAN	G	565	6	-	0/2/19/22	0/1/1/1
6	MAN	G	566	6	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

The worst 5 of 6 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	G	560	NAG	C2-N2-C7	-2.45	119.90	123.04
6	C	565	MAN	C2-C3-C4	-2.13	107.42	111.04
6	G	565	MAN	C2-C3-C4	-2.12	107.44	111.04
6	G	560	NAG	O5-C5-C6	2.26	112.23	107.35
6	C	562	BMA	C1-C2-C3	2.58	112.60	109.54

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.6 Ligand geometry

9 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
5	BU3	B	213	-	4,5,5	1.51	1 (25%)	6,6,6	0.58	0
4	SO4	B	214	-	4,4,4	0.20	0	6,6,6	0.13	0
4	SO4	B	215	-	4,4,4	0.16	0	6,6,6	0.12	0
4	SO4	B	216	-	4,4,4	0.18	0	6,6,6	0.22	0
7	NAG	C	573	3	14,14,15	0.55	0	15,19,21	0.71	0
7	NAG	G	573	3	14,14,15	0.52	0	15,19,21	0.61	0
4	SO4	H	226	-	4,4,4	0.17	0	6,6,6	0.09	0
5	BU3	L	213	-	4,5,5	1.51	1 (25%)	6,6,6	0.58	0
4	SO4	L	214	-	4,4,4	0.20	0	6,6,6	0.12	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
5	BU3	B	213	-	-	0/4/4/4	0/0/0/0
4	SO4	B	214	-	-	0/0/0/0	0/0/0/0
4	SO4	B	215	-	-	0/0/0/0	0/0/0/0
4	SO4	B	216	-	-	0/0/0/0	0/0/0/0
7	NAG	C	573	3	-	0/6/23/26	0/1/1/1
7	NAG	G	573	3	-	0/6/23/26	0/1/1/1
4	SO4	H	226	-	-	0/0/0/0	0/0/0/0
5	BU3	L	213	-	-	0/4/4/4	0/0/0/0
4	SO4	L	214	-	-	0/0/0/0	0/0/0/0

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	L	213	BU3	O6-C3	-2.57	1.37	1.43
5	B	213	BU3	O6-C3	-2.57	1.37	1.43

There are no bond angle outliers.

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
4	L	214	SO4	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

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	224/248 (90%)	1.08	47 (20%) 1 1	16, 36, 95, 131	0
1	H	227/248 (91%)	0.38	4 (1%) 71 67	16, 30, 63, 85	0
2	B	211/216 (97%)	0.47	8 (3%) 44 38	19, 40, 68, 83	0
2	L	211/216 (97%)	0.35	7 (3%) 50 44	16, 34, 69, 85	0
3	C	90/124 (72%)	1.96	36 (40%) 0 0	30, 68, 109, 126	0
3	G	90/124 (72%)	1.83	32 (35%) 0 0	38, 66, 100, 133	0
All	All	1053/1176 (89%)	0.80	134 (12%) 5 4	16, 39, 90, 133	0

The worst 5 of 134 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	G	190	LEU	8.5
1	A	191	THR	8.2
3	C	221	GLY	8.1
3	C	190	LEU	7.6
1	A	158	ALA	7.6

6.2 Non-standard residues in protein, DNA, RNA chains

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
1	TYS	A	100(H)	16/17	0.93	0.12	-	24,34,64,67	0
1	TYS	H	100(G)	16/17	0.93	0.12	-	30,38,88,90	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
1	PCA	H	2	8/9	0.96	0.09	-	30,39,48,54	0
1	TYS	H	100(H)	16/17	0.89	0.11	-	30,36,91,100	0
1	TYS	A	100(G)	16/17	0.93	0.11	-	27,32,75,76	0
1	PCA	A	2	8/9	0.97	0.11	-	19,24,33,33	0

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(\AA^2)	Q<0.9
6	MAN	G	565	11/12	0.97	0.12	1.37	21,25,38,54	0
6	MAN	C	565	11/12	0.95	0.10	0.95	23,27,40,43	0
6	MAN	C	566	11/12	0.97	0.12	0.64	22,26,32,36	0
6	NAG	G	560	14/15	0.98	0.09	-0.46	26,32,41,43	0
6	MAN	G	566	11/12	0.96	0.09	-0.54	20,23,26,27	0
6	NAG	G	561	14/15	0.93	0.09	-0.81	24,31,43,44	0
6	NAG	C	560	14/15	0.95	0.09	-1.12	28,34,43,50	0
6	MAN	C	563	11/12	0.81	0.23	-	74,80,87,88	0
6	BMA	G	562	11/12	0.93	0.08	-	25,34,42,52	0
6	MAN	C	564	11/12	0.97	0.13	-	25,28,33,37	0
6	NAG	C	561	14/15	0.91	0.10	-	28,36,45,53	0
6	MAN	G	563	11/12	0.86	0.15	-	52,68,75,80	0
6	MAN	G	564	11/12	0.98	0.09	-	21,25,28,28	0
6	BMA	C	562	11/12	0.94	0.08	-	31,43,53,60	0

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, 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(\AA^2)	Q<0.9
4	SO4	L	214	5/5	0.94	0.22	17.45	78,83,90,96	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
4	SO4	B	216	5/5	0.93	0.20	4.59	59,62,77,88	0
4	SO4	B	215	5/5	0.93	0.22	3.42	40,57,71,83	0
4	SO4	H	226	5/5	0.94	0.14	1.24	76,80,89,94	0
7	NAG	G	573	14/15	0.87	0.15	1.01	39,46,53,59	0
5	BU3	L	213	6/6	0.81	0.14	0.94	34,44,58,60	0
7	NAG	C	573	14/15	0.92	0.08	-1.01	31,38,47,51	0
5	BU3	B	213	6/6	0.95	0.11	-1.95	34,44,58,60	0
4	SO4	B	214	5/5	0.88	0.10	-	67,79,95,103	0

6.5 Other polymers

There are no such residues in this entry.