Full wwPDB X-ray Structure Validation Report ⓘ

May 3, 2025 – 02:41 PM EDT

PDB ID	:	3GKR / pdb_00003gkr
Title	:	Crystal Structure of Weissella viridescens FemX:UDP-MurNAc-hexapeptide complex
Authors	:	Delfosse, V.; Piton, J.; Villet, R.; Lecerf, M.; Arthur, M.; Mayer, C.
Deposited on	:	2009-03-11
Resolution	:	1.60 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

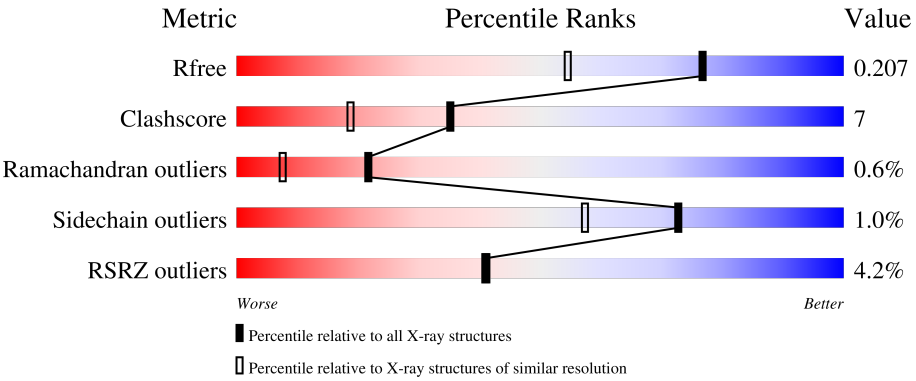
MolProbity	:	4-5-2 with Phenix2.0rc1
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0rc1
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.006 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.43.1

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 1.60 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}	164625	4274 (1.60-1.60)
Clashscore	180529	4682 (1.60-1.60)
Ramachandran outliers	177936	4583 (1.60-1.60)
Sidechain outliers	177891	4582 (1.60-1.60)
RSRZ outliers	164620	4272 (1.60-1.60)

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 of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	336	4% 83% 13% .
2	B	5	60% 40%

2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 3292 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 FemX.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	335	2687	1705	444	526	12	0	0	0

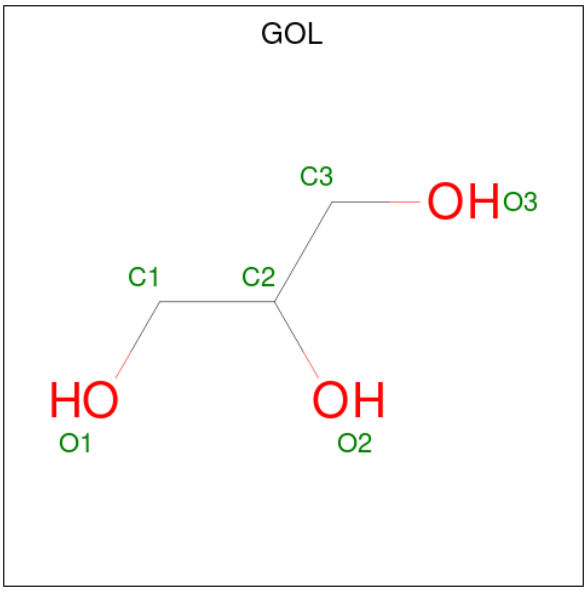
- Molecule 2 is a protein called UDP-MurNAc-peptide.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
2	B	5	77	40	9	26	2	0	0	0

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

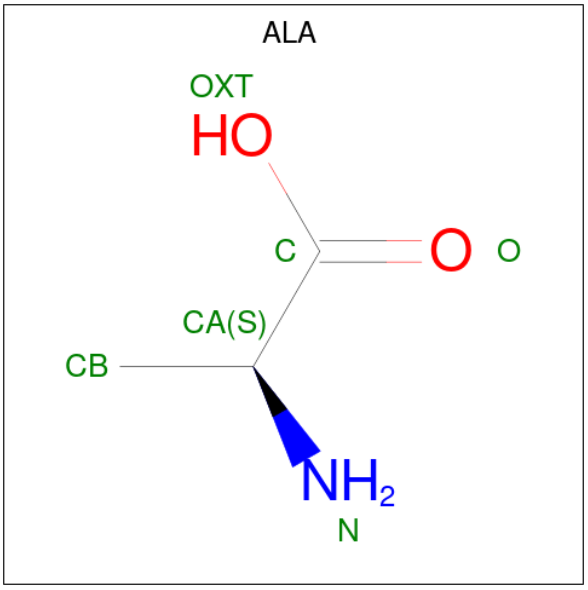
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Mg		
3	A	3	3	3	0	0

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O 6 3 3	0	0

- Molecule 5 is ALANINE (CCD ID: ALA) (formula: $C_3H_7NO_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	B	1	Total C N O 5 3 1 1	0	0

- Molecule 6 is water.

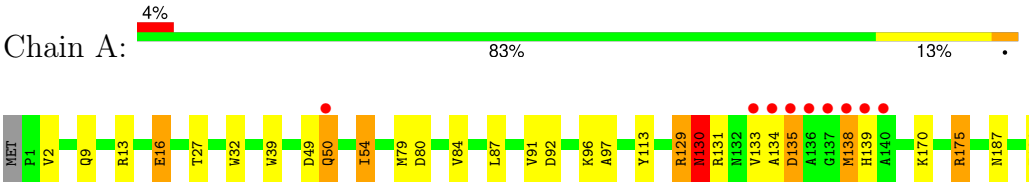
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	491	Total O 491 491	0	0
6	B	23	Total O 23 23	0	0

3 Residue-property plots (i)

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

- Molecule 1: FemX

Chain A:



- Molecule 2: UDP-MurNAc-peptide

Chain B:

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	42.32Å 101.84Å 46.85Å 90.00° 102.85° 90.00°	Depositor
Resolution (Å)	27.20 – 1.60 27.20 – 1.60	Depositor EDS
% Data completeness (in resolution range)	100.0 (27.20-1.60) 99.4 (27.20-1.60)	Depositor EDS
R_{merge}	0.03	Depositor
R_{sym}	0.03	Depositor
$< I/\sigma(I) >$ [1]	3.11 (at 1.60Å)	Xtriage
Refinement program	REFMAC 5.5.0066	Depositor
R, R_{free}	0.173 , 0.216 0.170 , 0.207	Depositor DCC
R_{free} test set	2531 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å^2)	12.4	Xtriage
Anisotropy	0.044	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.41 , 50.1	EDS
L-test for twinning[2]	$< \|L\| > = 0.47$, $< L^2 > = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	3292	wwPDB-VP
Average B, all atoms (Å^2)	15.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.55% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: UMA, GOL, MG, DAL, FGA

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	1.63	21/2748 (0.8%)	1.36	6/3728 (0.2%)
2	B	2.38	0/8	3.05	1/8 (12.5%)
All	All	1.63	21/2756 (0.8%)	1.37	7/3736 (0.2%)

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	16	GLU	C-O	-8.69	1.14	1.24
1	A	129	ARG	C-O	-7.98	1.15	1.24
1	A	80	ASP	C-O	-7.05	1.15	1.23
1	A	331	GLU	C-O	-6.92	1.15	1.24
1	A	130	ASN	C-O	-6.89	1.15	1.24
1	A	218	ARG	C-O	-6.51	1.16	1.24
1	A	84	VAL	CA-CB	6.00	1.62	1.54
1	A	205	ARG	C-N	-5.89	1.25	1.33
1	A	306	HIS	CA-C	5.86	1.60	1.52
1	A	306	HIS	N-CA	-5.80	1.38	1.46
1	A	243	GLY	C-O	-5.80	1.18	1.24
1	A	175	ARG	C-O	-5.70	1.17	1.24
1	A	54	ILE	CG1-CD1	5.56	1.73	1.51
1	A	297	THR	C-O	-5.40	1.17	1.24
1	A	187	ASN	C-O	-5.40	1.18	1.24
1	A	332	LEU	C-O	-5.28	1.17	1.24
1	A	222	ALA	N-CA	5.16	1.52	1.46
1	A	310	LYS	C-O	-5.12	1.17	1.24
1	A	231	PHE	N-CA	5.11	1.52	1.46
1	A	87	LEU	N-CA	5.10	1.52	1.46
1	A	39	TRP	N-CA	5.04	1.51	1.45

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	3	LYS	N-CA-CB	-8.02	96.86	110.50
1	A	13	ARG	N-CA-C	-7.51	103.17	111.36
1	A	50	GLN	N-CA-C	-6.95	104.42	113.17
1	A	170	LYS	N-CA-C	-6.46	104.24	111.28
1	A	205	ARG	O-C-N	-6.01	114.38	122.43
1	A	91	VAL	N-CA-C	5.45	117.86	111.05
1	A	54	ILE	CB-CG1-CD1	5.19	124.69	113.80

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2687	0	2599	40	1
2	B	77	0	57	0	0
3	A	3	0	0	0	0
4	A	6	0	8	0	0
5	B	5	0	4	0	0
6	A	491	0	0	20	1
6	B	23	0	0	0	0
All	All	3292	0	2668	40	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.

All (40) 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:16:GLU:HG3	6:A:967:HOH:O	1.12	1.27
1:A:263:ASN:HB2	6:A:864:HOH:O	1.46	1.14
1:A:224:ASP:HB2	6:A:788:HOH:O	1.56	1.06
1:A:202:MET:SD	6:A:852:HOH:O	2.23	0.95
1:A:190:THR:HG23	6:A:714:HOH:O	1.77	0.85
1:A:49:ASP:HB2	1:A:50:GLN:OE1	1.78	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:298:ASP:C	6:A:904:HOH:O	2.23	0.81
1:A:130:ASN:HD22	1:A:131:ARG:H	1.29	0.79
1:A:134:ALA:O	1:A:135:ASP:HB3	1.90	0.71
1:A:298:ASP:HA	6:A:904:HOH:O	1.90	0.71
1:A:310:LYS:HG3	6:A:860:HOH:O	1.94	0.66
1:A:297:THR:O	6:A:904:HOH:O	2.14	0.64
1:A:298:ASP:CA	6:A:904:HOH:O	2.44	0.62
1:A:16:GLU:OE1	6:A:856:HOH:O	2.16	0.61
1:A:2:VAL:CG2	6:A:678:HOH:O	2.49	0.61
1:A:134:ALA:O	1:A:135:ASP:CB	2.50	0.60
1:A:96:LYS:HE3	6:A:809:HOH:O	2.02	0.60
1:A:96:LYS:CE	6:A:809:HOH:O	2.50	0.59
1:A:138:MET:SD	1:A:139:HIS:CE1	2.97	0.57
1:A:130:ASN:HD22	1:A:131:ARG:N	1.99	0.57
1:A:310:LYS:CG	6:A:860:HOH:O	2.51	0.54
1:A:175:ARG:NH1	6:A:740:HOH:O	2.41	0.53
1:A:50:GLN:OE1	1:A:50:GLN:N	2.43	0.51
1:A:130:ASN:ND2	1:A:131:ARG:H	2.05	0.49
1:A:96:LYS:NZ	6:A:809:HOH:O	2.46	0.48
1:A:175:ARG:NH2	6:A:740:HOH:O	2.47	0.47
1:A:271:VAL:CG1	1:A:275:MET:HE3	2.45	0.45
1:A:199:TYR:CE2	1:A:210:HIS:HB2	2.52	0.45
1:A:54:ILE:O	1:A:54:ILE:HG12	2.17	0.44
1:A:27:THR:HG22	1:A:32:TRP:CE2	2.53	0.43
1:A:2:VAL:HG23	6:A:678:HOH:O	2.16	0.43
1:A:92:ASP:O	1:A:96:LYS:HD2	2.18	0.43
1:A:16:GLU:CG	6:A:967:HOH:O	1.97	0.43
1:A:200:THR:O	1:A:204:GLU:HG3	2.19	0.42
1:A:2:VAL:HG22	1:A:97:ALA:HB2	2.01	0.42
1:A:113:TYR:CD2	1:A:113:TYR:C	2.98	0.42
1:A:129:ARG:HB2	1:A:319:GLU:HB2	2.02	0.42
1:A:133:VAL:O	1:A:133:VAL:HG23	2.20	0.41
1:A:224:ASP:OD1	1:A:224:ASP:C	2.64	0.41
1:A:246:LEU:HD23	1:A:255:MET:HE3	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:A:9:GLN:NE2	6:A:861:HOH:O[1_554]	2.14	0.06

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	333/336 (99%)	324 (97%)	7 (2%)	2 (1%)	22 7

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	135	ASP
1	A	79	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	288/289 (100%)	285 (99%)	3 (1%)	73 57
2	B	1/1 (100%)	1 (100%)	0	100 100
All	All	289/290 (100%)	286 (99%)	3 (1%)	73 57

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	130	ASN
1	A	138	MET
1	A	296	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	9	GLN
1	A	20	GLN
1	A	130	ASN
1	A	139	HIS
1	A	187	ASN
1	A	220	GLN
1	A	263	ASN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

4 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	$\#\|Z\| > 2$
2	FGA	B	2	2	6,8,9	1.06	0	6,9,11	0.95	0
2	UMA	B	1	2	48,50,51	1.38	5 (10%)	66,74,76	1.94	18 (27%)

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	FGA	B	2	2	-	2/8/8/9	-
2	UMA	B	1	2	-	6/39/77/79	0/3/3/3

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1	UMA	O3'-C18	3.86	1.52	1.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1	UMA	O4B-C1B	2.74	1.48	1.42
2	B	1	UMA	C8'-C7'	2.72	1.56	1.50
2	B	1	UMA	C6-C5	2.59	1.41	1.35
2	B	1	UMA	O5'-C1'	2.22	1.47	1.41

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	1	UMA	O5'-C1'-O1'	-5.54	104.12	111.36
2	B	1	UMA	C1'-C2'-N2'	-5.11	102.35	110.92
2	B	1	UMA	C4-N3-C2	-4.38	121.18	126.61
2	B	1	UMA	O4B-C1B-C2B	-3.95	98.15	106.62
2	B	1	UMA	N3-C2-N1	3.83	119.88	114.89
2	B	1	UMA	C5-C4-N3	3.81	120.14	114.80
2	B	1	UMA	C21-N4-C19	-3.36	118.25	122.93
2	B	1	UMA	C23-C21-N4	-3.26	106.00	109.68
2	B	1	UMA	C2B-C1B-N1	-3.23	104.25	113.25
2	B	1	UMA	O4B-C1B-N1	-3.15	101.21	108.36
2	B	1	UMA	O1'-C1'-C2'	3.02	113.86	108.40
2	B	1	UMA	C1B-N1-C2	-2.57	112.99	117.59
2	B	1	UMA	O4-C4-C5	-2.52	120.81	125.16
2	B	1	UMA	C1B-N1-C6	2.46	126.03	120.78
2	B	1	UMA	O3'-C18-C19	-2.34	106.53	111.33
2	B	1	UMA	C4B-O4B-C1B	-2.18	104.66	109.47
2	B	1	UMA	O5'-C1'-C2'	-2.11	106.60	110.59
2	B	1	UMA	C5-C6-N1	-2.03	118.54	121.84

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	1	UMA	C20-C18-C19-O18
2	B	1	UMA	C23-C21-C22-O19
2	B	2	FGA	C-CA-CB-CG
2	B	2	FGA	OE1-CD-CG-CB
2	B	1	UMA	C20-C18-C19-N4
2	B	1	UMA	C20-C18-O3'-C3'
2	B	1	UMA	O3'-C18-C19-O18
2	B	1	UMA	C19-C18-O3'-C3'

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates (i)

There are no oligosaccharides in this entry.

5.6 Ligand geometry (i)

Of 5 ligands modelled in this entry, 3 are monoatomic - leaving 2 for Mogul analysis.

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	$\#\|Z\| > 2$
4	GOL	A	404	-	5,5,5	0.61	0	5,5,5	1.37	1 (20%)
5	ALA	B	101	2	3,4,5	1.08	0	2,4,6	0.68	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
4	GOL	A	404	-	-	2/4/4/4	-
5	ALA	B	101	2	-	0/1/2/4	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	A	404	GOL	O3-C3-C2	-2.09	100.99	110.38

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	404	GOL	C1-C2-C3-O3
4	A	404	GOL	O2-C2-C3-O3

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	335/336 (99%)	-0.44	14 (4%) 41 41	5, 11, 31, 48	0
2	B	1/5 (20%)	-1.14	0 100 100	16, 16, 16, 16	0
All	All	336/341 (98%)	-0.44	14 (4%) 41 41	5, 11, 31, 48	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	136	ALA	10.7
1	A	134	ALA	8.4
1	A	133	VAL	6.1
1	A	140	ALA	4.7
1	A	135	ASP	4.6
1	A	139	HIS	3.7
1	A	138	MET	3.6
1	A	137	GLY	3.0
1	A	297	THR	2.9
1	A	263	ASN	2.6
1	A	298	ASP	2.3
1	A	202	MET	2.3
1	A	300	SER	2.2
1	A	50	GLN	2.0

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. 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	B-factors(Å^2)	Q<0.9
2	DAL	B	4	5/6	0.96	0.05	11,12,13,14	0
2	DAL	B	5	6/6	0.96	0.05	8,12,13,13	0
2	FGA	B	2	9/10	0.97	0.05	12,15,16,18	0
2	UMA	B	1	48/49	0.98	0.04	5,11,14,21	0

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

6.4 Ligands (i)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 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	B-factors(Å^2)	Q<0.9
4	GOL	A	404	6/6	0.83	0.13	7,11,27,29	0
5	ALA	B	101	5/6	0.90	0.12	20,22,23,24	0
3	MG	A	402	1/1	0.96	0.16	23,23,23,23	0
3	MG	A	403	1/1	0.97	0.05	23,23,23,23	0
3	MG	A	401	1/1	0.98	0.11	20,20,20,20	0

6.5 Other polymers (i)

There are no such residues in this entry.