Full wwPDB X-ray Structure Validation Report (i)

Oct 18, 2023 – 04:57 PM EDT

PDB ID	:	2GEY
Title	:	Crystal Structure of AclR a putative hydroxylase from Streptomyces galilaeus
Authors	:	Beinker, P.; Lohkamp, B.; Schneider, G.
Deposited on	:	2006-03-21
Resolution	:	1.80 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 (i) 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 (i)) were used in the production of this report:

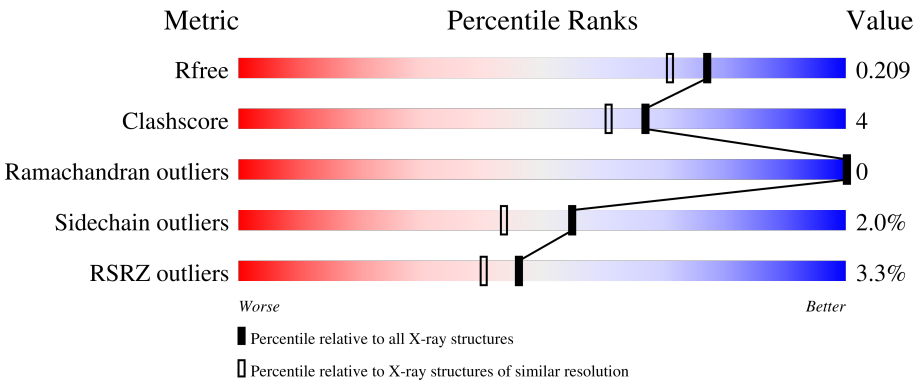
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 1.80 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}	130704	5950 (1.80-1.80)
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)
RSRZ outliers	127900	5850 (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 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	158	2%, 92%, 7%
1	B	158	6%, 77%, 7%, 16%
1	C	158	3%, 90%, 5%
1	D	158	%, 80%, 16%

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 5413 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 AclR protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	157	1308	838	226	235	9	6	11	0
			Total	C	N	O	S			
1	B	133	1117	716	189	202	10	0	10	0
			Total	C	N	O	S			
1	C	150	1224	786	207	222	9	11	8	0
			Total	C	N	O	S			
1	D	133	1127	725	191	201	10	0	12	0

There are 52 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	146	LYS	-	expression tag	UNP Q1XDX7
A	147	LEU	-	expression tag	UNP Q1XDX7
A	148	ALA	-	expression tag	UNP Q1XDX7
A	149	ALA	-	expression tag	UNP Q1XDX7
A	150	ALA	-	expression tag	UNP Q1XDX7
A	151	LEU	-	expression tag	UNP Q1XDX7
A	152	GLU	-	expression tag	UNP Q1XDX7
A	153	HIS	-	expression tag	UNP Q1XDX7
A	154	HIS	-	expression tag	UNP Q1XDX7
A	155	HIS	-	expression tag	UNP Q1XDX7
A	156	HIS	-	expression tag	UNP Q1XDX7
A	157	HIS	-	expression tag	UNP Q1XDX7
A	158	HIS	-	expression tag	UNP Q1XDX7
B	146	LYS	-	expression tag	UNP Q1XDX7
B	147	LEU	-	expression tag	UNP Q1XDX7
B	148	ALA	-	expression tag	UNP Q1XDX7
B	149	ALA	-	expression tag	UNP Q1XDX7
B	150	ALA	-	expression tag	UNP Q1XDX7
B	151	LEU	-	expression tag	UNP Q1XDX7
B	152	GLU	-	expression tag	UNP Q1XDX7
B	153	HIS	-	expression tag	UNP Q1XDX7

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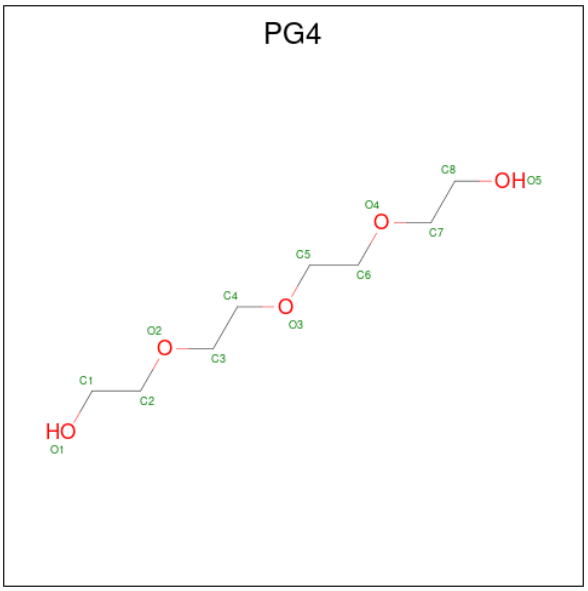
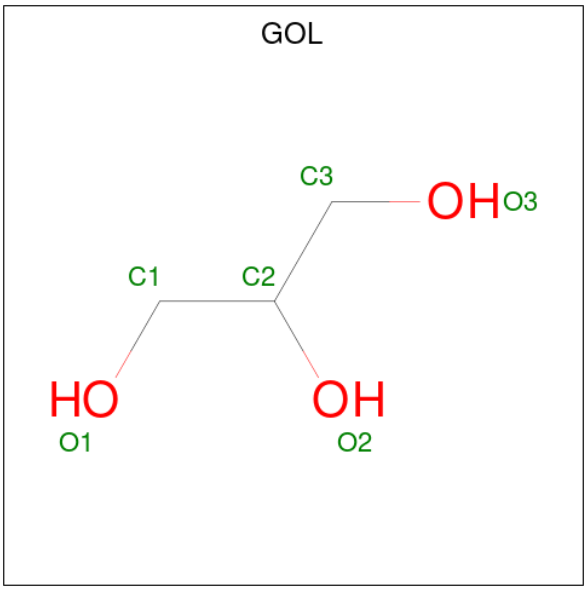
Chain	Residue	Modelled	Actual	Comment	Reference
B	154	HIS	-	expression tag	UNP Q1XDX7
B	155	HIS	-	expression tag	UNP Q1XDX7
B	156	HIS	-	expression tag	UNP Q1XDX7
B	157	HIS	-	expression tag	UNP Q1XDX7
B	158	HIS	-	expression tag	UNP Q1XDX7
C	146	LYS	-	expression tag	UNP Q1XDX7
C	147	LEU	-	expression tag	UNP Q1XDX7
C	148	ALA	-	expression tag	UNP Q1XDX7
C	149	ALA	-	expression tag	UNP Q1XDX7
C	150	ALA	-	expression tag	UNP Q1XDX7
C	151	LEU	-	expression tag	UNP Q1XDX7
C	152	GLU	-	expression tag	UNP Q1XDX7
C	153	HIS	-	expression tag	UNP Q1XDX7
C	154	HIS	-	expression tag	UNP Q1XDX7
C	155	HIS	-	expression tag	UNP Q1XDX7
C	156	HIS	-	expression tag	UNP Q1XDX7
C	157	HIS	-	expression tag	UNP Q1XDX7
C	158	HIS	-	expression tag	UNP Q1XDX7
D	146	LYS	-	expression tag	UNP Q1XDX7
D	147	LEU	-	expression tag	UNP Q1XDX7
D	148	ALA	-	expression tag	UNP Q1XDX7
D	149	ALA	-	expression tag	UNP Q1XDX7
D	150	ALA	-	expression tag	UNP Q1XDX7
D	151	LEU	-	expression tag	UNP Q1XDX7
D	152	GLU	-	expression tag	UNP Q1XDX7
D	153	HIS	-	expression tag	UNP Q1XDX7
D	154	HIS	-	expression tag	UNP Q1XDX7
D	155	HIS	-	expression tag	UNP Q1XDX7
D	156	HIS	-	expression tag	UNP Q1XDX7
D	157	HIS	-	expression tag	UNP Q1XDX7
D	158	HIS	-	expression tag	UNP Q1XDX7

- Molecule 2 is TETRAETHYLENE GLYCOL (three-letter code: PG4) (formula: $C_8H_{18}O_5$).



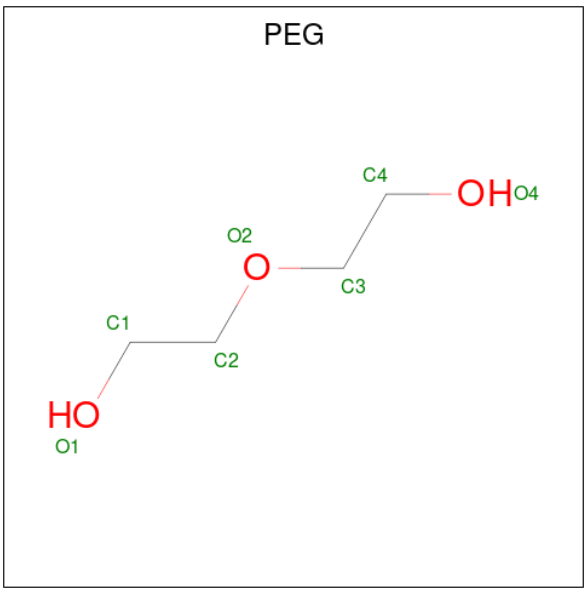
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
2	A	1	13	8	5	0	0
2	B	1	13	8	5	0	0
2	C	1	13	8	5	0	0
2	D	1	13	8	5	0	0

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



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

- Molecule 4 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: $C_4H_{10}O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
4	A	1	7	4	3	0	0

- Molecule 5 is water.

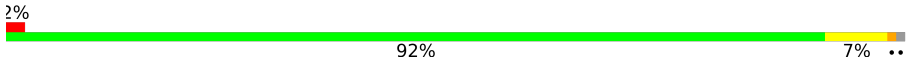
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
5	A	163	163	163	0	0
5	B	95	95	95	0	0
5	C	136	136	136	0	0
5	D	154	154	154	0	0

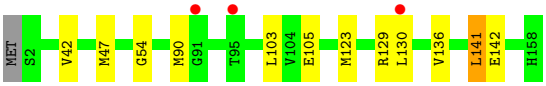
3 Residue-property plots (i)

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: AclR protein

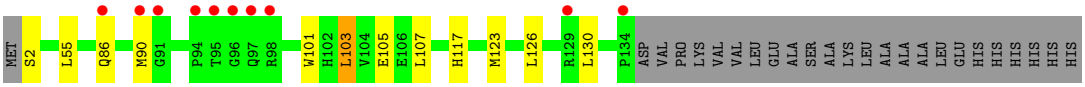
Chain A:





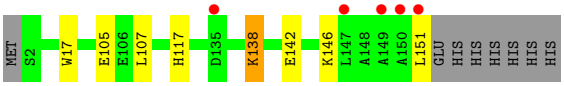
- Molecule 1: AclR protein

Chain B:



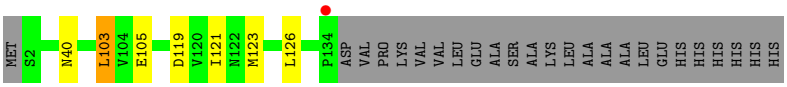
- Molecule 1: AclR protein

Chain C:



- Molecule 1: AclR protein

Chain D:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	80.91Å 81.72Å 136.64Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.98 – 1.80 40.45 – 1.80	Depositor EDS
% Data completeness (in resolution range)	98.1 (43.98-1.80) 98.1 (40.45-1.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$< I/\sigma(I) >$ [1]	2.09 (at 1.79Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.175 , 0.197 0.185 , 0.209	Depositor DCC
R_{free} test set	4133 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å^2)	22.9	Xtriage
Anisotropy	0.047	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 44.5	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	0.014 for k,h,-l	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	5413	wwPDB-VP
Average B, all atoms (Å^2)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.00% 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: GOL, PG4, PEG

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	0.45	2/1372 (0.1%)	0.58	0/1857
1	B	0.34	0/1168	0.55	0/1579
1	C	0.50	1/1273 (0.1%)	0.59	0/1724
1	D	0.46	0/1187	0.61	0/1605
All	All	0.44	3/5000 (0.1%)	0.58	0/6765

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	142	GLU	CG-CD	11.03	1.68	1.51
1	A	142[A]	GLU	CG-CD	-6.75	1.41	1.51
1	A	142[B]	GLU	CG-CD	-6.75	1.41	1.51

There are no bond angle outliers.

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	1308	0	1328	9	0
1	B	1117	0	1138	13	0
1	C	1224	0	1258	7	2
1	D	1127	0	1163	8	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	13	0	18	3	0
2	B	13	0	18	1	0
2	C	13	0	18	2	0
2	D	13	0	18	1	0
3	A	6	0	8	0	0
3	B	12	0	16	1	0
3	C	6	0	8	0	0
3	D	6	0	8	0	0
4	A	7	0	9	0	0
5	A	163	0	0	0	0
5	B	95	0	0	0	2
5	C	136	0	0	0	0
5	D	154	0	0	1	0
All	All	5413	0	5008	38	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 4.

All (38) 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:B:107[B]:LEU:CD2	1:B:117[B]:HIS:CD2	2.52	0.92
1:B:107[B]:LEU:HD21	1:B:117[B]:HIS:CD2	2.05	0.92
1:B:107[B]:LEU:CD2	1:B:117[B]:HIS:HD2	1.85	0.89
1:B:107[B]:LEU:HD21	1:B:117[B]:HIS:HD2	1.36	0.86
1:C:105[B]:GLU:CG	1:C:117[B]:HIS:NE2	2.52	0.73
1:C:105[B]:GLU:HG3	1:C:117[B]:HIS:NE2	2.05	0.72
1:B:103[B]:LEU:HD23	1:B:105:GLU:HG3	1.75	0.68
1:D:103[B]:LEU:HB2	1:D:121[B]:ILE:HD13	1.76	0.68
1:B:103[B]:LEU:CD2	1:B:105:GLU:HG3	2.28	0.63
1:D:103[B]:LEU:HD23	1:D:105:GLU:HG2	1.82	0.60
1:C:17:TRP:CH2	2:C:1001:PG4:H41	2.41	0.55
1:D:103[B]:LEU:HB2	1:D:121[B]:ILE:CD1	2.38	0.53
1:B:123[B]:MET:CE	1:B:126:LEU:HD23	2.40	0.52
1:A:123[A]:MET:CE	2:A:1002:PG4:H22	2.39	0.52
1:B:101:TRP:HZ2	2:B:1004:PG4:H21	1.77	0.49
1:C:105[B]:GLU:CD	1:C:117[B]:HIS:CE1	2.86	0.49
1:B:90:MET:HE2	1:B:130:LEU:HD12	1.94	0.48
1:B:107[B]:LEU:HD23	1:B:117[B]:HIS:CD2	2.46	0.47
1:C:105[B]:GLU:HG3	1:C:117[B]:HIS:CE1	2.50	0.47
1:C:105[A]:GLU:OE1	2:C:1001:PG4:H12	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:90:MET:CE	1:B:130:LEU:HD12	2.44	0.46
1:D:123[B]:MET:CE	1:D:126:LEU:HD23	2.45	0.46
1:D:103[A]:LEU:HD12	1:D:105:GLU:HG2	1.97	0.45
1:A:103[A]:LEU:CD1	1:A:105:GLU:HG3	2.46	0.45
1:A:54:GLY:HA3	2:A:1002:PG4:H42	1.97	0.45
1:A:90:MET:HE1	1:A:130:LEU:HD12	1.99	0.45
1:A:42:VAL:HG11	1:A:47:MET:HE2	1.99	0.44
1:A:136:VAL:CG1	1:A:141:LEU:HD13	2.48	0.44
1:A:123[A]:MET:SD	2:A:1002:PG4:H22	2.57	0.44
1:D:119:ASP:HB2	1:D:121[B]:ILE:HD11	1.99	0.44
1:B:123[B]:MET:HE2	1:B:126:LEU:HD23	2.00	0.43
1:D:40:ASN:ND2	5:D:2024:HOH:O	2.51	0.43
1:A:103[A]:LEU:HD13	1:A:105:GLU:HG3	2.01	0.43
1:B:2:SER:N	3:B:2003:GOL:HO3	2.18	0.42
2:D:1003:PG4:H41	2:D:1003:PG4:H21	1.82	0.41
1:A:90:MET:HB2	1:A:90:MET:HE2	1.92	0.41
1:C:105[B]:GLU:CG	1:C:117[B]:HIS:CE1	3.04	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:C:138:LYS:NZ	5:B:2045:HOH:O[2_554]	0.52	1.68
1:C:138:LYS:CE	5:B:2045:HOH:O[2_554]	2.03	0.17

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	166/158 (105%)	164 (99%)	2 (1%)	0	100 100
1	B	141/158 (89%)	139 (99%)	2 (1%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	156/158 (99%)	154 (99%)	2 (1%)	0	100	100
1	D	143/158 (90%)	143 (100%)	0	0	100	100
All	All	606/632 (96%)	600 (99%)	6 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	146/136 (107%)	144 (99%)	2 (1%)	67	59
1	B	126/136 (93%)	122 (97%)	4 (3%)	39	25
1	C	136/136 (100%)	131 (96%)	5 (4%)	34	19
1	D	128/136 (94%)	126 (98%)	2 (2%)	62	54
All	All	536/544 (98%)	523 (98%)	13 (2%)	55	36

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	129	ARG
1	A	141	LEU
1	B	55	LEU
1	B	86	GLN
1	B	103[A]	LEU
1	B	103[B]	LEU
1	C	107[A]	LEU
1	C	107[B]	LEU
1	C	138	LYS
1	C	146	LYS
1	C	151	LEU
1	D	103[A]	LEU
1	D	103[B]	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	86	GLN
1	A	157	HIS
1	B	86	GLN
1	D	40	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)

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

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

10 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	GOL	B	2005	-	5,5,5	0.36	0	5,5,5	0.21	0
3	GOL	D	2002	-	5,5,5	0.35	0	5,5,5	0.35	0
2	PG4	A	1002	-	12,12,12	0.60	0	11,11,11	0.65	0
2	PG4	D	1003	-	12,12,12	0.52	0	11,11,11	0.50	0
4	PEG	A	3001	-	6,6,6	0.99	0	5,5,5	1.67	2 (40%)
2	PG4	B	1004	-	12,12,12	0.71	0	11,11,11	0.88	1 (9%)
3	GOL	A	2001	-	5,5,5	0.32	0	5,5,5	0.33	0
3	GOL	B	2003	-	5,5,5	0.29	0	5,5,5	0.28	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
2	PG4	C	1001	-	12,12,12	0.43	0	11,11,11	0.72	0
3	GOL	C	2004	-	5,5,5	0.34	0	5,5,5	0.34	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
3	GOL	B	2005	-	-	2/4/4/4	-
3	GOL	D	2002	-	-	2/4/4/4	-
2	PG4	A	1002	-	-	5/10/10/10	-
2	PG4	D	1003	-	-	5/10/10/10	-
4	PEG	A	3001	-	-	1/4/4/4	-
2	PG4	B	1004	-	-	6/10/10/10	-
3	GOL	A	2001	-	-	4/4/4/4	-
3	GOL	B	2003	-	-	2/4/4/4	-
2	PG4	C	1001	-	-	4/10/10/10	-
3	GOL	C	2004	-	-	2/4/4/4	-

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	A	3001	PEG	O2-C3-C4	2.33	120.32	110.07
4	A	3001	PEG	O2-C2-C1	2.23	119.88	110.07
2	B	1004	PG4	O3-C5-C6	2.17	120.18	110.39

There are no chirality outliers.

All (33) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	2001	GOL	O1-C1-C2-C3
3	B	2005	GOL	C1-C2-C3-O3
3	B	2005	GOL	O2-C2-C3-O3
3	C	2004	GOL	O1-C1-C2-C3
2	B	1004	PG4	C6-C5-O3-C4
2	D	1003	PG4	C4-C3-O2-C2
2	A	1002	PG4	O2-C3-C4-O3

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Mol	Chain	Res	Type	Atoms
3	A	2001	GOL	C1-C2-C3-O3
3	B	2003	GOL	O1-C1-C2-C3
3	D	2002	GOL	C1-C2-C3-O3
2	A	1002	PG4	O1-C1-C2-O2
2	B	1004	PG4	O4-C7-C8-O5
2	C	1001	PG4	O4-C7-C8-O5
3	A	2001	GOL	O1-C1-C2-O2
3	C	2004	GOL	O1-C1-C2-O2
3	D	2002	GOL	O2-C2-C3-O3
2	B	1004	PG4	O3-C5-C6-O4
2	C	1001	PG4	C6-C5-O3-C4
2	B	1004	PG4	C4-C3-O2-C2
2	A	1002	PG4	C4-C3-O2-C2
2	D	1003	PG4	O4-C7-C8-O5
2	B	1004	PG4	C1-C2-O2-C3
2	A	1002	PG4	C6-C5-O3-C4
2	B	1004	PG4	O2-C3-C4-O3
3	A	2001	GOL	O2-C2-C3-O3
3	B	2003	GOL	O1-C1-C2-O2
2	D	1003	PG4	C8-C7-O4-C6
2	A	1002	PG4	C1-C2-O2-C3
4	A	3001	PEG	C4-C3-O2-C2
2	C	1001	PG4	C4-C3-O2-C2
2	D	1003	PG4	O2-C3-C4-O3
2	D	1003	PG4	C6-C5-O3-C4
2	C	1001	PG4	C1-C2-O2-C3

There are no ring outliers.

5 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1002	PG4	3	0
2	D	1003	PG4	1	0
2	B	1004	PG4	1	0
3	B	2003	GOL	1	0
2	C	1001	PG4	2	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 (i)

6.1 Protein, DNA and RNA chains (i)

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	157/158 (99%)	0.21	3 (1%)	66	63	23, 27, 33, 36	3 (1%)
1	B	133/158 (84%)	0.46	10 (7%)	14	11	24, 28, 34, 38	2 (1%)
1	C	150/158 (94%)	0.18	5 (3%)	46	40	21, 27, 38, 48	12 (8%)
1	D	133/158 (84%)	0.18	1 (0%)	86	84	21, 27, 33, 37	2 (1%)
All	All	573/632 (90%)	0.26	19 (3%)	46	40	21, 27, 35, 48	19 (3%)

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	134	PRO	4.2
1	D	134	PRO	4.2
1	B	91	GLY	4.0
1	B	97	GLN	3.2
1	C	150	ALA	3.2
1	C	151	LEU	3.2
1	A	95	THR	3.2
1	A	91	GLY	3.1
1	B	86	GLN	2.9
1	B	90	MET	2.9
1	C	135	ASP	2.7
1	B	98	ARG	2.7
1	B	94	PRO	2.6
1	B	129	ARG	2.5
1	B	95	THR	2.4
1	C	147	LEU	2.3
1	B	96	GLY	2.2
1	C	149	ALA	2.1
1	A	130	LEU	2.1

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)

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, 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	B-factors(Å^2)	Q<0.9
2	PG4	B	1004	13/13	0.59	0.24	56,57,58,58	0
4	PEG	A	3001	7/7	0.70	0.21	43,45,46,46	0
3	GOL	B	2003	6/6	0.74	0.23	55,56,56,56	0
3	GOL	C	2004	6/6	0.76	0.31	45,46,47,48	0
2	PG4	A	1002	13/13	0.77	0.20	47,51,57,57	0
2	PG4	C	1001	13/13	0.80	0.23	56,57,59,60	0
2	PG4	D	1003	13/13	0.83	0.17	53,54,58,59	0
3	GOL	B	2005	6/6	0.88	0.13	54,54,55,55	0
3	GOL	D	2002	6/6	0.89	0.14	45,46,46,47	0
3	GOL	A	2001	6/6	0.90	0.19	42,42,43,43	0

6.5 Other polymers (i)

There are no such residues in this entry.