



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 11:22 AM GMT

PDB ID : 3OSV
Title : The crystal structure of FLGD from P. Aeruginosa
Authors : Wang, D.; Luo, M.; Niu, S.
Deposited on : 2010-09-10
Resolution : 2.35 Å(reported)

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

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

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

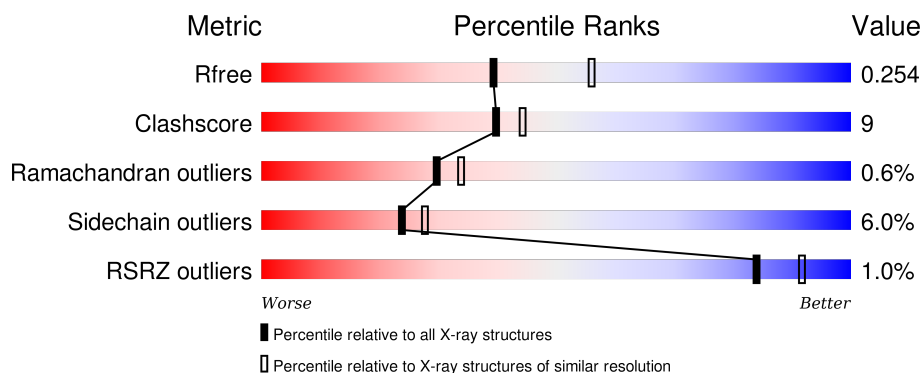
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.35 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	1352 (2.38-2.34)
Clashscore	102246	1456 (2.38-2.34)
Ramachandran outliers	100387	1435 (2.38-2.34)
Sidechain outliers	100360	1436 (2.38-2.34)
RSRZ outliers	91569	1358 (2.38-2.34)

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	138	<div> <div>%</div> <div> <div></div> <div>70%</div> <div>22%</div> <div>• 5%</div> </div> </div>
1	B	138	<div> <div>%</div> <div> <div></div> <div>75%</div> <div>14%</div> <div>5%</div> <div>5%</div> </div> </div>
1	C	138	<div> <div>%</div> <div> <div></div> <div>80%</div> <div>14%</div> <div>• •</div> </div> </div>
1	D	138	<div> <div></div> <div> <div></div> <div>74%</div> <div>16%</div> <div>• • 7%</div> </div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	GOL	A	139	-	-	-	X
2	GOL	A	140	-	-	X	X
2	GOL	A	141	-	-	-	X
2	GOL	B	138	-	-	-	X
2	GOL	B	139	-	-	-	X
2	GOL	B	140	-	-	-	X
2	GOL	C	138	-	-	-	X
2	GOL	D	138	-	-	-	X
2	GOL	D	139	-	-	X	X

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 4112 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 Flagellar basal-body rod modification protein FlgD.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	131	Total	C	N	O	S	11	0	0
			970	612	161	194	3			
1	B	131	Total	C	N	O	S	10	0	0
			963	607	161	192	3			
1	C	132	Total	C	N	O	S	12	0	0
			970	612	162	193	3			
1	D	129	Total	C	N	O	S	11	0	0
			957	605	159	190	3			

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



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

Continued on next page...

Continued from previous page...

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

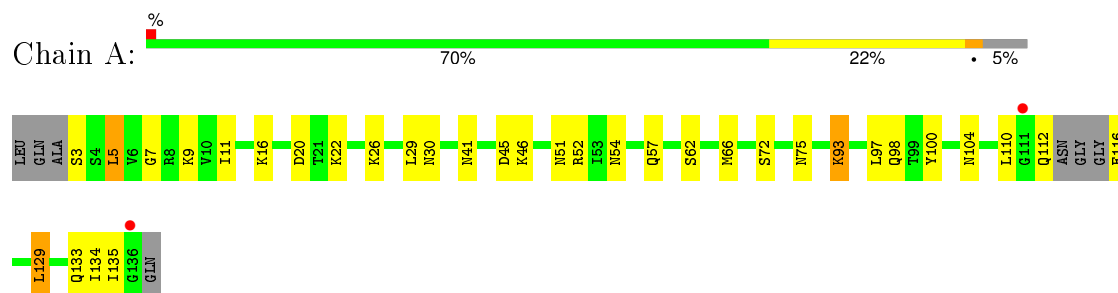
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	41	Total	O	0	0
			41	41		
3	B	48	Total	O	0	0
			48	48		
3	C	56	Total	O	0	0
			56	56		
3	D	47	Total	O	0	0
			47	47		

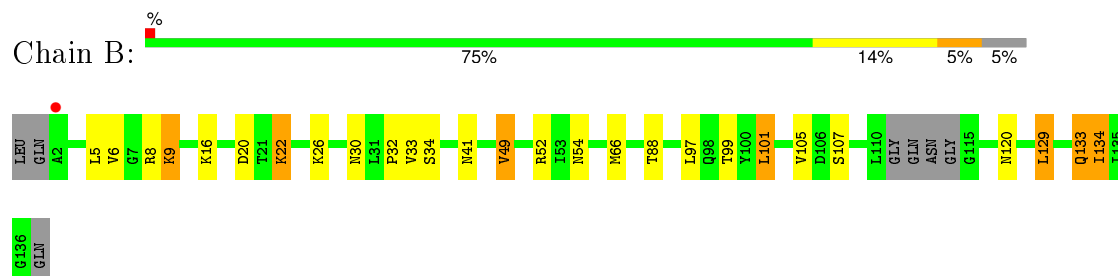
3 Residue-property plots [i](#)

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.

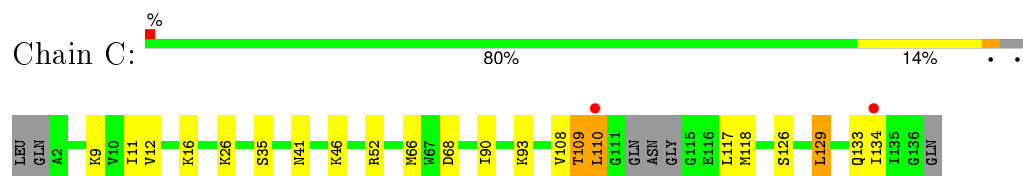
- Molecule 1: Flagellar basal-body rod modification protein FlgD



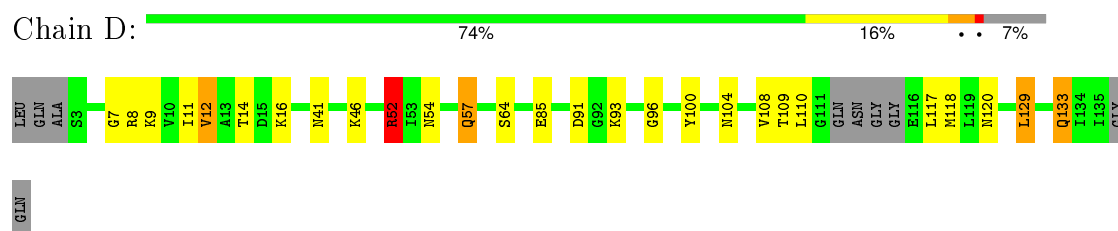
- Molecule 1: Flagellar basal-body rod modification protein FlgD



- Molecule 1: Flagellar basal-body rod modification protein FlgD



- Molecule 1: Flagellar basal-body rod modification protein FlgD



4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	114.83Å 118.38Å 118.30Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.07 – 2.35 48.07 – 2.34	Depositor EDS
% Data completeness (in resolution range)	99.7 (48.07-2.35) 99.7 (48.07-2.34)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.41 (at 2.34Å)	Xtriage
Refinement program	REFMAC 5.5.0072	Depositor
R, R_{free}	0.211 , 0.268 0.203 , 0.254	Depositor DCC
R_{free} test set	1715 reflections (5.33%)	DCC
Wilson B-factor (Å ²)	40.8	Xtriage
Anisotropy	0.192	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 34.6	EDS
Estimated twinning fraction	0.049 for -h,-l,-k 0.000 for k,h,-l 0.000 for -l,-k,-h 0.000 for l,-h,-k 0.000 for -k,-l,h	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.36$	Xtriage
Outliers	1 of 34027 reflections (0.003%)	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4112	wwPDB-VP
Average B, all atoms (Å ²)	33.0	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 47.35 % 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 1.0043e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GOL

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	1.10	5/983 (0.5%)	1.04	7/1333 (0.5%)
1	B	1.30	5/976 (0.5%)	1.06	6/1324 (0.5%)
1	C	1.21	4/983 (0.4%)	1.00	3/1333 (0.2%)
1	D	1.27	3/970 (0.3%)	1.13	8/1316 (0.6%)
All	All	1.22	17/3912 (0.4%)	1.06	24/5306 (0.5%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	46	LYS	CD-CE	-18.77	1.04	1.51
1	B	133	GLN	CA-CB	-17.17	1.16	1.53
1	C	46	LYS	CG-CD	-13.14	1.07	1.52
1	B	133	GLN	CG-CD	-12.37	1.22	1.51
1	C	16	LYS	CD-CE	-11.25	1.23	1.51
1	B	9	LYS	CE-NZ	10.41	1.75	1.49
1	A	9	LYS	CB-CG	-9.07	1.28	1.52
1	D	9	LYS	CG-CD	-9.04	1.21	1.52
1	D	93	LYS	CG-CD	-7.14	1.28	1.52
1	B	22	LYS	CG-CD	-6.39	1.30	1.52
1	C	9	LYS	CG-CD	-6.26	1.31	1.52
1	C	93	LYS	CG-CD	-5.54	1.33	1.52
1	A	16	LYS	CD-CE	5.48	1.65	1.51
1	A	112	GLN	CA-CB	-5.44	1.42	1.53

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	49	VAL	CB-CG2	-5.24	1.41	1.52
1	A	93	LYS	CG-CD	-5.16	1.34	1.52
1	A	110	LEU	CG-CD2	-5.06	1.33	1.51

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	133	GLN	CG-CD-OE1	-10.15	101.29	121.60
1	A	110	LEU	CD1-CG-CD2	8.70	136.58	110.50
1	D	52	ARG	NE-CZ-NH1	8.63	124.61	120.30
1	B	129	LEU	CA-CB-CG	-7.76	97.46	115.30
1	A	112	GLN	CB-CA-C	-7.66	95.09	110.40
1	D	129	LEU	CA-CB-CG	-7.58	97.86	115.30
1	D	133	GLN	CG-CD-OE1	7.52	136.65	121.60
1	D	133	GLN	CG-CD-NE2	-7.34	99.08	116.70
1	C	46	LYS	CB-CG-CD	7.22	130.38	111.60
1	A	129	LEU	CA-CB-CG	-7.14	98.87	115.30
1	D	91	ASP	C-N-CA	-6.90	107.81	122.30
1	C	129	LEU	CA-CB-CG	-6.57	100.19	115.30
1	D	9	LYS	CG-CD-CE	6.55	131.55	111.90
1	B	49	VAL	CB-CA-C	-6.32	99.40	111.40
1	A	16	LYS	CG-CD-CE	-6.10	93.60	111.90
1	B	133	GLN	N-CA-CB	6.09	121.56	110.60
1	C	68	ASP	CB-CG-OD1	5.71	123.44	118.30
1	A	110	LEU	CB-CG-CD2	-5.50	101.64	111.00
1	D	46	LYS	CG-CD-CE	5.39	128.08	111.90
1	B	22	LYS	CB-CG-CD	5.36	125.53	111.60
1	A	45	ASP	CB-CG-OD1	5.26	123.03	118.30
1	A	45	ASP	CB-CG-OD2	-5.10	113.71	118.30
1	B	9	LYS	CB-CG-CD	5.09	124.83	111.60
1	D	16	LYS	CG-CD-CE	-5.01	96.88	111.90

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	133	GLN	Sidechain

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	970	0	978	23	0
1	B	963	0	968	21	0
1	C	970	0	980	13	0
1	D	957	0	969	20	0
2	A	24	0	32	8	0
2	B	18	0	24	0	0
2	C	6	0	8	0	0
2	D	12	0	16	4	0
3	A	41	0	0	0	0
3	B	48	0	0	2	0
3	C	56	0	0	0	0
3	D	47	0	0	0	0
All	All	4112	0	3975	68	0

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 9.

All (68) 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:D:109:THR:HG22	1:D:110:LEU:H	1.24	1.02
1:A:51:ASN:HD21	2:A:139:GOL:H31	1.28	0.95
1:D:96:GLY:CA	2:D:139:GOL:H31	2.04	0.88
1:A:133:GLN:NE2	1:B:134:ILE:H	1.76	0.84
1:C:109:THR:HG22	1:C:110:LEU:H	1.44	0.81
1:D:41:ASN:HD21	1:D:52:ARG:HH11	1.24	0.81
1:D:109:THR:HG22	1:D:110:LEU:N	1.97	0.80
1:D:52:ARG:HE	1:D:54:ASN:HD21	1.32	0.77
1:C:109:THR:CG2	1:C:110:LEU:H	1.97	0.77
1:D:96:GLY:HA3	2:D:139:GOL:H31	1.66	0.76
1:A:52:ARG:HE	1:A:54:ASN:HD21	1.34	0.76
1:B:52:ARG:HE	1:B:54:ASN:HD21	1.34	0.75
2:A:140:GOL:H12	1:C:52:ARG:HH12	1.54	0.73
1:B:41:ASN:HD21	1:B:52:ARG:HH11	1.33	0.73
1:D:96:GLY:HA2	2:D:139:GOL:H31	1.72	0.72

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:11:ILE:HB	1:A:133:GLN:HB2	1.72	0.72
1:A:26:LYS:HG2	1:A:66:MET:HE2	1.74	0.70
1:D:109:THR:CG2	1:D:110:LEU:H	1.92	0.69
1:C:109:THR:HG22	1:C:110:LEU:N	2.08	0.68
1:C:35:SER:HB2	1:C:90:ILE:CD1	2.24	0.67
1:D:109:THR:OG1	1:D:120:ASN:ND2	2.27	0.67
1:B:16:LYS:HE3	3:B:183:HOH:O	1.94	0.67
1:C:117:LEU:O	1:C:118:MET:HG2	1.95	0.66
1:C:26:LYS:HG2	1:C:66:MET:HG2	1.79	0.65
1:A:41:ASN:HD21	1:A:52:ARG:HH11	1.44	0.64
1:C:41:ASN:HD21	1:C:52:ARG:HH11	1.45	0.62
1:A:20:ASP:OD2	1:A:22:LYS:HE3	2.00	0.61
1:B:52:ARG:HE	1:B:54:ASN:ND2	2.01	0.59
1:A:52:ARG:HE	1:A:54:ASN:ND2	2.02	0.58
1:A:135:ILE:H	1:B:30:ASN:HD21	1.52	0.57
1:D:117:LEU:O	1:D:118:MET:HG2	2.05	0.57
1:B:26:LYS:HG2	1:B:66:MET:HG2	1.87	0.56
1:C:35:SER:HB2	1:C:90:ILE:HD11	1.89	0.54
1:A:133:GLN:HE22	1:B:134:ILE:H	1.50	0.54
2:A:140:GOL:H12	1:C:52:ARG:NH1	2.23	0.52
1:B:20:ASP:OD1	1:B:22:LYS:HG2	2.10	0.52
1:D:41:ASN:HD21	1:D:52:ARG:NH1	2.01	0.51
1:D:117:LEU:C	1:D:118:MET:HG2	2.29	0.51
1:A:133:GLN:NE2	1:B:134:ILE:N	2.53	0.51
1:B:41:ASN:HD21	1:B:52:ARG:NH1	2.06	0.50
1:B:99:THR:HB	1:B:101:LEU:HD13	1.93	0.50
1:A:98:GLN:OE1	2:A:141:GOL:H32	2.13	0.49
1:D:7:GLY:HA2	1:D:104:ASN:HD21	1.78	0.49
1:C:133:GLN:O	1:C:134:ILE:C	2.51	0.49
1:A:54:ASN:HB2	2:A:138:GOL:H11	1.95	0.48
1:C:11:ILE:HG22	1:C:133:GLN:HG3	1.96	0.47
1:A:30:ASN:ND2	1:A:100:TYR:OH	2.47	0.47
1:D:52:ARG:HE	1:D:54:ASN:ND2	2.06	0.46
1:A:51:ASN:ND2	2:A:139:GOL:H31	2.12	0.46
1:A:133:GLN:HE21	1:B:134:ILE:H	1.56	0.46
1:D:11:ILE:HG22	1:D:100:TYR:HB3	1.98	0.46
1:B:5:LEU:O	1:B:8:ARG:HB2	2.16	0.45
1:D:109:THR:CG2	1:D:110:LEU:N	2.62	0.44
1:A:5:LEU:HD22	1:B:32:PRO:HA	2.00	0.43
1:B:16:LYS:CE	3:B:183:HOH:O	2.60	0.43
1:A:29:LEU:HD13	1:A:97:LEU:HD13	2.01	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:33:VAL:HG12	1:B:34:SER:N	2.34	0.42
1:D:11:ILE:CG2	1:D:100:TYR:HB3	2.49	0.42
2:A:140:GOL:C1	1:C:52:ARG:HH22	2.31	0.42
1:B:88:THR:HB	1:B:97:LEU:HD11	2.02	0.41
1:B:6:VAL:HG22	1:B:107:SER:HA	2.01	0.41
1:D:85:GLU:HG3	2:D:139:GOL:H2	2.02	0.41
1:A:133:GLN:HE21	1:B:134:ILE:N	2.16	0.41
1:A:62:SER:OG	1:B:134:ILE:HG23	2.21	0.41
1:A:72:SER:OG	2:A:140:GOL:H31	2.21	0.41
1:A:7:GLY:HA2	1:A:104:ASN:HD21	1.86	0.41
1:A:75:ASN:HD21	1:D:57:GLN:H	1.68	0.40
1:D:12:VAL:HG22	1:D:14:THR:HB	2.02	0.40

There are no symmetry-related clashes.

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	127/138 (92%)	123 (97%)	2 (2%)	2 (2%)	12	10
1	B	127/138 (92%)	123 (97%)	3 (2%)	1 (1%)	24	26
1	C	128/138 (93%)	124 (97%)	4 (3%)	0	100	100
1	D	125/138 (91%)	118 (94%)	7 (6%)	0	100	100
All	All	507/552 (92%)	488 (96%)	16 (3%)	3 (1%)	30	34

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	134	ILE
1	A	5	LEU
1	A	134	ILE

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	109/114 (96%)	103 (94%)	6 (6%)	27	32
1	B	107/114 (94%)	101 (94%)	6 (6%)	26	31
1	C	108/114 (95%)	102 (94%)	6 (6%)	26	31
1	D	108/114 (95%)	100 (93%)	8 (7%)	17	19
All	All	432/456 (95%)	406 (94%)	26 (6%)	24	28

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	SER
1	A	46	LYS
1	A	57	GLN
1	A	93	LYS
1	A	116	GLU
1	A	129	LEU
1	B	9	LYS
1	B	49	VAL
1	B	101	LEU
1	B	105	VAL
1	B	120	ASN
1	B	129	LEU
1	C	12	VAL
1	C	108	VAL
1	C	109	THR
1	C	110	LEU
1	C	126	SER
1	C	129	LEU
1	D	8	ARG
1	D	12	VAL
1	D	52	ARG
1	D	57	GLN
1	D	64	SER
1	D	108	VAL

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	D	129	LEU
1	D	133	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	41	ASN
1	A	54	ASN
1	A	75	ASN
1	A	104	ASN
1	A	133	GLN
1	B	30	ASN
1	B	41	ASN
1	B	54	ASN
1	B	120	ASN
1	C	30	ASN
1	C	41	ASN
1	C	104	ASN
1	D	41	ASN
1	D	54	ASN
1	D	104	ASN
1	D	120	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 carbohydrates 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	$\# Z > 2$
2	GOL	A	138	-	5,5,5	0.16	0	5,5,5	0.76	0
2	GOL	A	139	-	5,5,5	0.44	0	5,5,5	0.83	0
2	GOL	A	140	-	5,5,5	0.28	0	5,5,5	0.45	0
2	GOL	A	141	-	5,5,5	0.43	0	5,5,5	0.72	0
2	GOL	B	138	-	5,5,5	0.49	0	5,5,5	0.49	0
2	GOL	B	139	-	5,5,5	0.53	0	5,5,5	0.22	0
2	GOL	B	140	-	5,5,5	0.45	0	5,5,5	1.08	0
2	GOL	C	138	-	5,5,5	0.32	0	5,5,5	1.18	0
2	GOL	D	138	-	5,5,5	0.51	0	5,5,5	0.66	0
2	GOL	D	139	-	5,5,5	0.66	0	5,5,5	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
2	GOL	A	138	-	-	0/4/4/4	0/0/0/0
2	GOL	A	139	-	-	0/4/4/4	0/0/0/0
2	GOL	A	140	-	-	0/4/4/4	0/0/0/0
2	GOL	A	141	-	-	0/4/4/4	0/0/0/0
2	GOL	B	138	-	-	0/4/4/4	0/0/0/0
2	GOL	B	139	-	-	0/4/4/4	0/0/0/0
2	GOL	B	140	-	-	0/4/4/4	0/0/0/0
2	GOL	C	138	-	-	0/4/4/4	0/0/0/0
2	GOL	D	138	-	-	0/4/4/4	0/0/0/0
2	GOL	D	139	-	-	0/4/4/4	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.

5 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	138	GOL	1	0
2	A	139	GOL	2	0
2	A	140	GOL	4	0
2	A	141	GOL	1	0
2	D	139	GOL	4	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	131/138 (94%)	0.06	2 (1%) 76 85	22, 31, 54, 78	5 (3%)
1	B	131/138 (94%)	-0.11	1 (0%) 87 93	22, 31, 53, 63	4 (3%)
1	C	132/138 (95%)	-0.02	2 (1%) 76 85	21, 30, 52, 72	5 (3%)
1	D	129/138 (93%)	-0.10	0 100 100	21, 30, 47, 71	5 (3%)
All	All	523/552 (94%)	-0.04	5 (0%) 84 92	21, 30, 52, 78	19 (3%)

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	136	GLY	3.3
1	C	110	LEU	2.7
1	A	111	GLY	2.3
1	C	134	ILE	2.3
1	B	2	ALA	2.3

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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains.

The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	GOL	D	138	6/6	0.94	0.28	14.97	43,51,54,56	0
2	GOL	D	139	6/6	0.88	0.28	14.48	41,46,53,53	0
2	GOL	A	141	6/6	0.74	0.33	13.24	56,62,64,65	0
2	GOL	A	139	6/6	0.83	0.27	6.58	47,53,54,57	0
2	GOL	B	140	6/6	0.89	0.25	6.56	47,50,52,53	0
2	GOL	C	138	6/6	0.88	0.24	6.42	44,46,50,54	0
2	GOL	A	140	6/6	0.92	0.23	6.39	52,55,57,61	0
2	GOL	B	138	6/6	0.96	0.20	4.81	41,49,51,54	0
2	GOL	B	139	6/6	0.94	0.18	3.93	47,49,51,53	0
2	GOL	A	138	6/6	0.94	0.15	0.53	36,45,48,52	0

6.5 Other polymers ⓘ

There are no such residues in this entry.