



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 28, 2024 – 07:06 am BST

PDB ID : 1UZZ
Title : Erythrina cristagalli bound to N-linked oligosaccharide and lactose
Authors : Turton, K.; Natesh, R.; Thiyagarajan, N.; Chaddock, J.A.; Acharya, K.R.
Deposited on : 2004-03-20
Resolution : 2.13 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 : 4.02b-467
Mogul : 1.8.4, CSD as541be (2020)
Xtrriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36.2

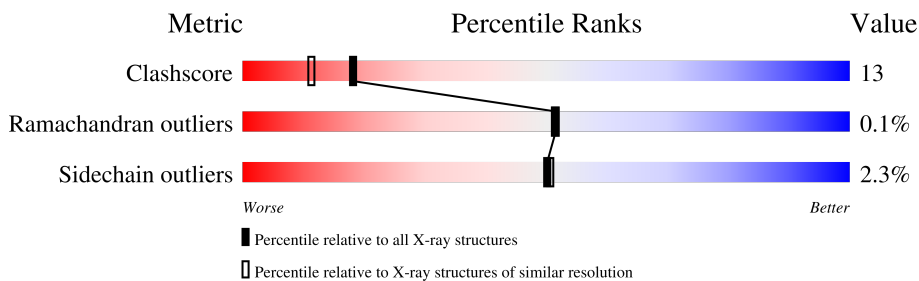
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.13 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))
Clashscore	141614	2653 (2.16-2.12)
Ramachandran outliers	138981	2618 (2.16-2.12)
Sidechain outliers	138945	2617 (2.16-2.12)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	242	
1	B	242	
1	C	242	
1	D	242	

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	1240	-	X	-	-
2	GOL	B	1240	-	X	-	-
2	GOL	D	1241	-	X	-	-

2 Entry composition [i](#)

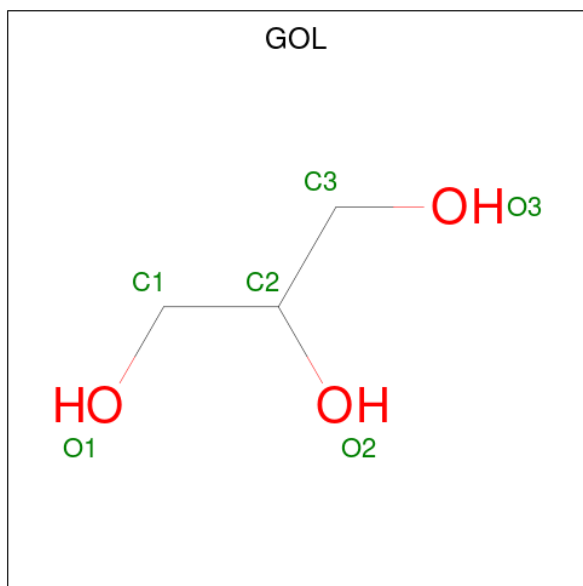
There are 5 unique types of molecules in this entry. The entry contains 8343 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Lectin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	240	Total 1856	C 1189	N 302	O 362	S 3	0	0	1
1	B	240	Total 1856	C 1189	N 302	O 362	S 3	0	0	1
1	C	241	Total 1863	C 1193	N 303	O 364	S 3	0	0	1
1	D	241	Total 1863	C 1193	N 303	O 364	S 3	0	0	1

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



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

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

- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

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

- Molecule 4 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Mn	0	0
			1	1		
4	B	1	Total	Mn	0	0
			1	1		
4	C	1	Total	Mn	0	0
			1	1		
4	D	1	Total	Mn	0	0
			1	1		

- Molecule 5 is water.

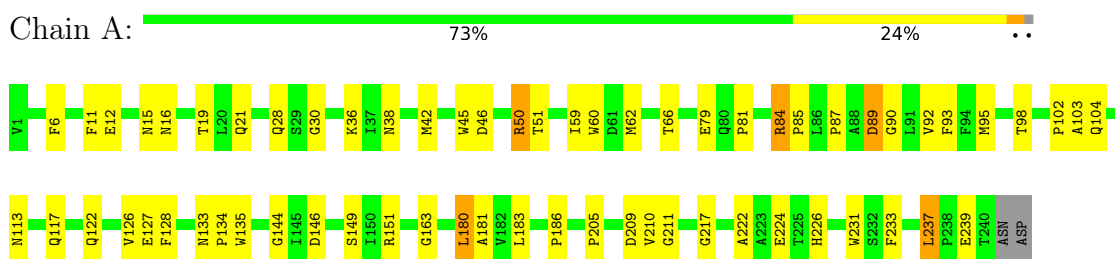
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	250	Total	O	0	0
			250	250		
5	B	242	Total	O	0	0
			242	242		
5	C	188	Total	O	0	0
			188	188		
5	D	199	Total	O	0	0
			199	199		

3 Residue-property plots [i](#)

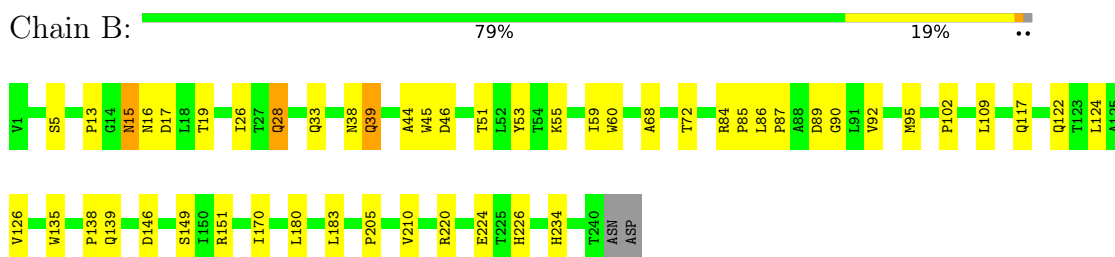
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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. 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.

Note EDS was not executed.

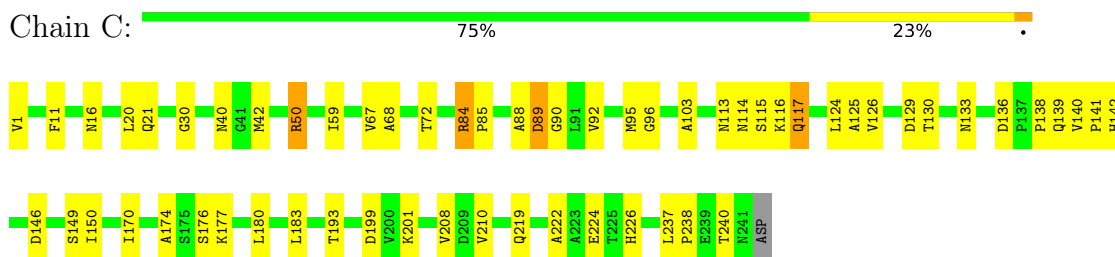
- Molecule 1: Lectin



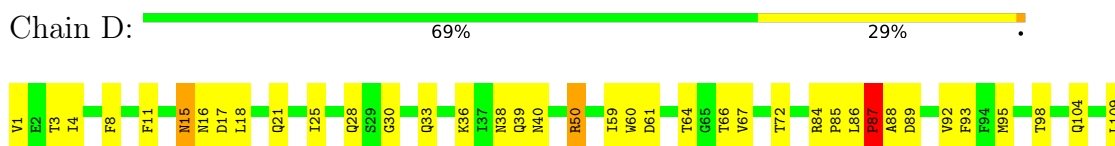
- Molecule 1: Lectin

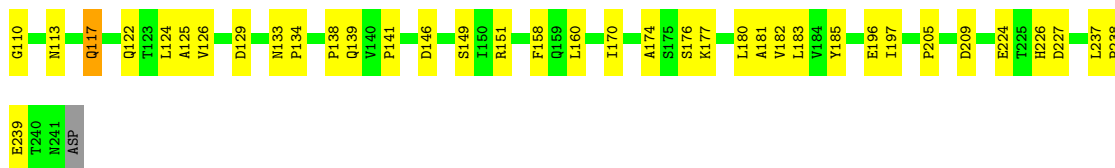


- Molecule 1: Lectin



- Molecule 1: Lectin





4 Data and refinement statistics

Xtrriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	55.28Å 55.37Å 86.93Å 86.23° 75.37° 82.13°	Depositor
Resolution (Å)	40.00 – 2.13	Depositor
% Data completeness (in resolution range)	90.3 (40.00-2.13)	Depositor
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.191 , 0.241	Depositor
Estimated twinning fraction	No twinning to report.	Xtrriage
Total number of atoms	8343	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

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: CA, GOL, MN

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	0.56	3/1907 (0.2%)	0.69	1/2609 (0.0%)
1	B	0.57	4/1907 (0.2%)	0.67	0/2609
1	C	0.51	3/1914 (0.2%)	0.70	2/2619 (0.1%)
1	D	0.50	3/1914 (0.2%)	0.69	2/2619 (0.1%)
All	All	0.54	13/7642 (0.2%)	0.69	5/10456 (0.0%)

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	89	ASP	CG-OD1	12.49	1.54	1.25
1	B	86	LEU	C-N	-11.44	1.12	1.34
1	A	87	PRO	N-CA	-10.76	1.28	1.47
1	B	90	GLY	CA-C	7.77	1.64	1.51
1	D	87	PRO	CA-CB	7.23	1.68	1.53
1	A	90	GLY	C-O	6.97	1.34	1.23
1	D	87	PRO	N-CA	-6.93	1.35	1.47
1	C	89	ASP	CB-CG	-6.49	1.38	1.51
1	B	89	ASP	C-N	6.47	1.44	1.33
1	D	87	PRO	CA-C	-6.03	1.40	1.52
1	B	87	PRO	CA-CB	5.26	1.64	1.53
1	C	90	GLY	CA-C	5.22	1.60	1.51
1	A	90	GLY	N-CA	5.05	1.53	1.46

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	89	ASP	CB-CG-OD2	11.56	128.71	118.30
1	D	87	PRO	N-CA-C	7.00	130.29	112.10
1	D	87	PRO	CA-CB-CG	-6.83	91.01	104.00
1	A	89	ASP	CB-CG-OD2	6.35	124.01	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	89	ASP	OD1-CG-OD2	-5.31	113.21	123.30

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	1856	0	1793	49	0
1	B	1856	0	1793	37	0
1	C	1863	0	1800	47	0
1	D	1863	0	1800	67	0
2	A	6	0	4	2	0
2	B	6	0	4	0	0
2	D	6	0	4	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
5	A	250	0	0	8	0
5	B	242	0	0	5	0
5	C	188	0	0	4	0
5	D	199	0	0	7	0
All	All	8343	0	7198	197	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 13.

All (197) 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:50:ARG:HD3	1:A:102:PRO:HB3	1.45	0.98
1:D:67:VAL:HG13	1:D:174:ALA:HB1	1.54	0.87
1:A:95:MET:HG2	1:A:210:VAL:HG12	1.60	0.81
1:C:138:PRO:HG2	1:C:139:GLN:NE2	1.97	0.79
1:C:114:ASN:ND2	1:C:116:LYS:HE2	1.98	0.78
1:D:59:ILE:HD13	1:D:95:MET:HE1	1.71	0.72
1:C:180:LEU:HD21	5:D:2171:HOH:O	1.88	0.72
1:C:59:ILE:CD1	1:C:208:VAL:HG22	2.23	0.69
1:C:67:VAL:HG13	1:C:174:ALA:HB1	1.75	0.68
1:C:84:ARG:HB2	1:C:84:ARG:NH1	2.07	0.68
1:A:135:TRP:HB2	1:A:151:ARG:HG2	1.73	0.68
1:B:95:MET:HG2	1:B:210:VAL:HG12	1.74	0.68
1:D:59:ILE:HG23	1:D:60:TRP:HD1	1.59	0.67
1:B:220:ARG:HB3	1:B:220:ARG:HH11	1.60	0.67
1:B:117:GLN:HG2	1:B:149:SER:HB2	1.76	0.66
1:C:84:ARG:HB2	1:C:84:ARG:HH11	1.59	0.66
1:C:16:ASN:HB2	5:C:2031:HOH:O	1.94	0.66
1:B:220:ARG:HB3	1:B:220:ARG:NH1	2.12	0.65
1:B:28:GLN:HE21	1:B:28:GLN:HA	1.62	0.65
1:C:95:MET:HG2	1:C:210:VAL:HG12	1.80	0.64
1:B:92:VAL:HG23	1:B:126:VAL:O	1.97	0.64
1:D:117:GLN:HE22	1:D:151:ARG:HD2	1.62	0.64
1:B:13:PRO:HG3	1:B:26:ILE:HG22	1.78	0.64
1:B:72:THR:HG22	1:B:170:ILE:HB	1.81	0.63
1:B:117:GLN:HG2	1:B:149:SER:CB	2.30	0.62
1:A:103:ALA:HB1	1:A:113:ASN:ND2	2.15	0.61
1:B:38:ASN:HB3	1:B:39:GLN:OE1	2.01	0.61
1:C:59:ILE:HD13	1:C:208:VAL:HG22	1.83	0.61
1:D:92:VAL:HG23	1:D:126:VAL:O	2.00	0.61
1:D:84:ARG:HA	1:D:85:PRO:C	2.21	0.61
1:C:103:ALA:HB1	1:C:113:ASN:ND2	2.15	0.61
1:A:12:GLU:HG3	1:A:15:ASN:HB2	1.84	0.60
1:D:138:PRO:HG2	1:D:139:GLN:NE2	2.16	0.60
1:D:8:PHE:CE2	1:D:18:LEU:HD21	2.37	0.60
1:A:21:GLN:OE1	1:A:50:ARG:NH1	2.35	0.59
1:C:224:GLU:HG3	1:C:226:HIS:CE1	2.37	0.59
1:A:60:TRP:HA	5:A:2100:HOH:O	2.03	0.59
1:D:39:GLN:H	1:D:39:GLN:CD	2.06	0.58
1:A:146:ASP:OD1	5:A:2172:HOH:O	2.17	0.58
1:A:117:GLN:HA	1:A:149:SER:HB2	1.85	0.58
1:A:127:GLU:HG2	5:A:2172:HOH:O	2.03	0.58
1:D:67:VAL:HG11	5:D:2163:HOH:O	2.03	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:28:GLN:HB2	5:D:2040:HOH:O	2.04	0.57
1:A:104:GLN:HG2	1:A:113:ASN:HD22	1.70	0.57
1:D:86:LEU:N	1:D:86:LEU:HD12	2.20	0.57
1:D:180:LEU:HD12	1:D:181:ALA:N	2.19	0.57
1:C:59:ILE:HD11	1:C:208:VAL:HG22	1.87	0.57
1:C:146:ASP:HB3	1:C:149:SER:O	2.05	0.57
1:C:193:THR:HB	1:D:180:LEU:HD21	1.86	0.56
1:D:86:LEU:H	1:D:86:LEU:CD1	2.18	0.56
1:D:59:ILE:CG2	1:D:60:TRP:HD1	2.19	0.56
1:C:130:THR:O	1:C:140:VAL:HG11	2.06	0.56
1:A:36:LYS:HE3	1:A:38:ASN:ND2	2.21	0.55
1:B:117:GLN:HA	1:B:149:SER:HB2	1.87	0.55
1:D:86:LEU:HD12	1:D:86:LEU:H	1.71	0.55
1:A:180:LEU:CD1	1:B:180:LEU:HD22	2.35	0.55
1:C:11:PHE:O	1:C:30:GLY:HA2	2.06	0.55
1:D:237:LEU:C	1:D:237:LEU:HD13	2.27	0.55
1:D:224:GLU:HG3	1:D:226:HIS:CE1	2.42	0.55
1:A:237:LEU:HG	5:A:2100:HOH:O	2.06	0.55
1:A:98:THR:HG22	1:A:209:ASP:OD1	2.08	0.54
1:B:39:GLN:HG3	5:B:2065:HOH:O	2.07	0.54
1:C:84:ARG:HH11	1:C:84:ARG:CB	2.19	0.54
1:D:66:THR:HB	1:D:238:PRO:HG2	1.89	0.54
1:C:59:ILE:N	1:C:59:ILE:HD12	2.23	0.54
1:C:138:PRO:HG2	1:C:139:GLN:HE22	1.71	0.53
1:C:1:VAL:HG12	1:C:238:PRO:HA	1.90	0.53
1:C:96:GLY:O	1:C:208:VAL:HB	2.08	0.53
1:D:67:VAL:CG1	1:D:174:ALA:HB1	2.33	0.53
1:A:6:PHE:CZ	1:A:233:PHE:HB3	2.45	0.52
1:C:133:ASN:HB2	1:C:136:ASP:OD2	2.09	0.52
1:D:237:LEU:HD22	1:D:238:PRO:HD2	1.91	0.52
1:C:115:SER:HA	1:C:150:ILE:HG22	1.89	0.52
1:D:21:GLN:OE1	1:D:50:ARG:HD3	2.10	0.52
1:A:144:GLY:C	5:A:2172:HOH:O	2.48	0.51
1:B:59:ILE:HG23	1:B:60:TRP:HD1	1.75	0.51
1:B:33:GLN:NE2	5:B:2061:HOH:O	2.43	0.51
1:C:117:GLN:HG3	5:C:2122:HOH:O	2.11	0.51
1:D:88:ALA:HB1	1:D:89:ASP:CG	2.31	0.51
1:D:129:ASP:O	1:D:141:PRO:HA	2.10	0.51
1:A:28:GLN:HE21	1:A:28:GLN:HA	1.75	0.51
1:D:117:GLN:HA	1:D:149:SER:HB2	1.93	0.51
1:A:180:LEU:HD12	1:B:180:LEU:HD22	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:237:LEU:HD11	5:C:2182:HOH:O	2.10	0.51
1:D:11:PHE:O	1:D:30:GLY:HA2	2.11	0.51
1:A:42:MET:CE	1:A:81:PRO:HB2	2.41	0.51
1:B:15:ASN:C	1:B:15:ASN:HD22	2.11	0.51
1:A:222:ALA:HB1	5:A:2248:HOH:O	2.10	0.50
1:B:5:SER:HB3	1:B:234:HIS:HD2	1.77	0.50
1:A:92:VAL:HG23	1:A:126:VAL:O	2.12	0.50
1:D:122:GLN:HB3	1:D:205:PRO:HD3	1.94	0.50
1:D:180:LEU:HD11	1:D:182:VAL:HG23	1.94	0.50
1:D:61:ASP:HB3	1:D:64:THR:OG1	2.12	0.50
1:D:92:VAL:HG22	1:D:93:PHE:N	2.26	0.50
1:D:38:ASN:HB2	1:D:40:ASN:OD1	2.11	0.50
1:C:115:SER:HA	1:C:150:ILE:CG2	2.41	0.49
1:D:196:GLU:HG3	1:D:197:ILE:N	2.27	0.49
1:C:140:VAL:HG13	1:C:141:PRO:HA	1.94	0.49
1:B:15:ASN:HD22	1:B:16:ASN:N	2.09	0.49
1:D:146:ASP:HB3	1:D:149:SER:O	2.12	0.49
1:D:124:LEU:HD23	1:D:124:LEU:C	2.33	0.49
1:C:40:ASN:ND2	1:C:42:MET:HG2	2.28	0.49
1:B:122:GLN:HB3	1:B:205:PRO:HD3	1.95	0.48
1:A:59:ILE:CG2	1:A:60:TRP:HD1	2.26	0.48
1:A:84:ARG:HA	1:A:85:PRO:C	2.32	0.48
1:A:79:GLU:HG3	1:A:163:GLY:HA2	1.95	0.48
1:A:180:LEU:HD22	5:B:2210:HOH:O	2.12	0.48
1:D:113:ASN:ND2	5:D:2108:HOH:O	2.46	0.48
1:A:128:PHE:CE1	1:A:183:LEU:HD12	2.49	0.48
1:D:36:LYS:HZ2	1:D:38:ASN:HD22	1.61	0.48
1:B:5:SER:HB3	1:B:234:HIS:CD2	2.48	0.48
1:C:72:THR:HG22	1:C:170:ILE:HB	1.94	0.48
1:C:21:GLN:OE1	1:C:50:ARG:HD2	2.14	0.47
1:C:124:LEU:HD13	1:C:124:LEU:C	2.35	0.47
1:A:89:ASP:OD1	2:A:1240:GOL:O2	2.33	0.47
1:B:15:ASN:ND2	1:B:17:ASP:H	2.12	0.47
1:D:36:LYS:NZ	1:D:38:ASN:HD22	2.12	0.47
1:B:53:TYR:CE2	1:B:55:LYS:HB2	2.50	0.47
1:A:66:THR:OG1	1:A:237:LEU:HD21	2.15	0.47
1:A:186:PRO:HD2	5:A:2192:HOH:O	2.15	0.47
1:C:84:ARG:HA	1:C:85:PRO:C	2.36	0.46
1:D:25:ILE:HD11	1:D:33:GLN:NE2	2.29	0.46
1:B:224:GLU:HG3	1:B:226:HIS:CE1	2.51	0.46
1:C:240:THR:HG21	5:C:2184:HOH:O	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:103:ALA:HB1	1:A:113:ASN:HD21	1.81	0.46
1:A:239:GLU:OE2	1:A:239:GLU:HA	2.16	0.46
1:C:68:ALA:HB2	1:C:237:LEU:HB2	1.96	0.46
1:D:4:ILE:HG13	5:D:2079:HOH:O	2.15	0.46
1:D:92:VAL:CG2	1:D:93:PHE:N	2.79	0.46
1:A:95:MET:HE2	1:A:95:MET:HB3	1.61	0.46
1:A:180:LEU:HG	1:A:181:ALA:N	2.29	0.46
1:B:19:THR:O	1:B:51:THR:HA	2.16	0.46
1:B:102:PRO:HG2	5:B:2133:HOH:O	2.16	0.46
1:C:199:ASP:OD1	1:C:201:LYS:HB2	2.16	0.46
1:D:17:ASP:O	1:D:18:LEU:HD23	2.15	0.46
1:D:183:LEU:C	1:D:183:LEU:HD23	2.36	0.45
1:A:122:GLN:HB3	1:A:205:PRO:HD3	1.99	0.45
1:B:183:LEU:C	1:B:183:LEU:HD23	2.37	0.45
1:B:39:GLN:NE2	5:B:2067:HOH:O	2.47	0.45
1:D:15:ASN:ND2	1:D:17:ASP:H	2.14	0.45
1:B:38:ASN:ND2	1:B:44:ALA:HB2	2.31	0.45
1:B:138:PRO:HG2	1:B:139:GLN:NE2	2.31	0.45
1:C:92:VAL:HG23	1:C:126:VAL:O	2.16	0.45
1:D:180:LEU:HD12	1:D:181:ALA:H	1.81	0.45
1:B:180:LEU:C	1:B:180:LEU:HD12	2.37	0.45
1:A:84:ARG:HH11	1:A:84:ARG:HB3	1.82	0.45
1:D:86:LEU:N	1:D:86:LEU:CD1	2.80	0.44
1:C:124:LEU:HD13	1:C:125:ALA:N	2.32	0.44
1:D:3:THR:HG22	1:D:4:ILE:N	2.32	0.44
1:B:59:ILE:HD12	1:B:68:ALA:HB3	1.98	0.44
1:C:21:GLN:OE1	1:C:50:ARG:CD	2.66	0.44
1:B:28:GLN:HE21	1:B:28:GLN:CA	2.28	0.44
1:C:20:LEU:N	1:C:20:LEU:HD22	2.33	0.43
1:B:135:TRP:HB2	1:B:151:ARG:HG2	1.99	0.43
1:C:129:ASP:HB3	1:C:142:HIS:CE1	2.54	0.43
1:D:33:GLN:NE2	5:D:2046:HOH:O	2.51	0.43
1:D:104:GLN:O	1:D:110:GLY:HA2	2.19	0.43
1:D:15:ASN:HD22	1:D:16:ASN:N	2.16	0.43
1:C:88:ALA:HB1	1:C:89:ASP:CG	2.40	0.43
1:A:211:GLY:HA2	1:A:231:TRP:HH2	1.84	0.42
1:A:224:GLU:HG3	1:A:226:HIS:CE1	2.54	0.42
1:D:39:GLN:CD	1:D:39:GLN:N	2.70	0.42
1:D:133:ASN:HB3	1:D:134:PRO:CD	2.48	0.42
1:A:211:GLY:HA2	1:A:231:TRP:CH2	2.54	0.42
1:D:33:GLN:HG2	1:D:227:ASP:OD1	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:98:THR:HA	1:D:209:ASP:OD2	2.19	0.42
1:D:117:GLN:HG2	5:D:2111:HOH:O	2.18	0.42
1:D:160:LEU:HD12	1:D:185:TYR:OH	2.19	0.42
1:A:217:GLY:HA2	2:A:1240:GOL:O2	2.19	0.42
1:B:45:TRP:CE3	1:B:46:ASP:HB2	2.54	0.42
1:D:88:ALA:HA	1:D:89:ASP:HA	1.80	0.42
1:A:19:THR:O	1:A:51:THR:HA	2.19	0.42
1:D:72:THR:HG22	1:D:170:ILE:HB	2.01	0.42
1:D:88:ALA:O	1:D:224:GLU:HG2	2.20	0.42
1:C:183:LEU:C	1:C:183:LEU:HD23	2.40	0.41
1:A:84:ARG:HB3	1:A:84:ARG:NH1	2.35	0.41
1:A:92:VAL:HG23	1:A:126:VAL:C	2.40	0.41
1:D:3:THR:CG2	1:D:4:ILE:N	2.82	0.41
1:D:1:VAL:HA	1:D:239:GLU:HG3	2.02	0.41
1:D:67:VAL:HG13	1:D:174:ALA:CB	2.39	0.41
1:D:124:LEU:HD23	1:D:125:ALA:N	2.36	0.41
1:B:84:ARG:HA	1:B:85:PRO:C	2.41	0.41
1:D:59:ILE:CD1	1:D:95:MET:HE1	2.47	0.41
1:D:176:SER:O	1:D:177:LYS:HB2	2.21	0.41
1:C:176:SER:O	1:C:177:LYS:HB2	2.19	0.41
1:A:11:PHE:O	1:A:30:GLY:HA2	2.21	0.41
1:A:92:VAL:HG22	1:A:93:PHE:N	2.35	0.41
1:A:222:ALA:CB	5:A:2248:HOH:O	2.68	0.41
1:C:114:ASN:CG	1:C:116:LYS:HE2	2.40	0.41
1:D:141:PRO:HG2	1:D:158:PHE:O	2.20	0.41
1:A:133:ASN:HB3	1:A:134:PRO:CD	2.51	0.40
1:A:45:TRP:CE3	1:A:46:ASP:HB2	2.55	0.40
1:A:28:GLN:HA	1:A:28:GLN:NE2	2.36	0.40
1:B:146:ASP:HB3	1:B:149:SER:O	2.22	0.40
1:C:219:GLN:HG3	1:C:222:ALA:HB2	2.03	0.40
1:C:117:GLN:HA	1:C:149:SER:HB2	2.04	0.40
1:D:21:GLN:OE1	1:D:50:ARG:CD	2.70	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	238/242 (98%)	230 (97%)	8 (3%)	0	100	100
1	B	238/242 (98%)	231 (97%)	7 (3%)	0	100	100
1	C	239/242 (99%)	226 (95%)	13 (5%)	0	100	100
1	D	239/242 (99%)	222 (93%)	16 (7%)	1 (0%)	34	29
All	All	954/968 (99%)	909 (95%)	44 (5%)	1 (0%)	51	51

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	87	PRO

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	205/208 (99%)	199 (97%)	6 (3%)	42	40
1	B	205/208 (99%)	200 (98%)	5 (2%)	49	49
1	C	206/208 (99%)	203 (98%)	3 (2%)	65	68
1	D	206/208 (99%)	201 (98%)	5 (2%)	49	49
All	All	822/832 (99%)	803 (98%)	19 (2%)	50	51

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	ASN
1	A	50	ARG
1	A	62	MET
1	A	84	ARG
1	A	180	LEU

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Mol	Chain	Res	Type
1	A	237	LEU
1	B	15	ASN
1	B	28	GLN
1	B	39	GLN
1	B	109	LEU
1	B	124	LEU
1	C	50	ARG
1	C	84	ARG
1	C	117	GLN
1	D	15	ASN
1	D	50	ARG
1	D	87	PRO
1	D	109	LEU
1	D	117	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	16	ASN
1	A	28	GLN
1	A	33	GLN
1	A	38	ASN
1	A	113	ASN
1	A	159	GLN
1	B	15	ASN
1	B	28	GLN
1	B	33	GLN
1	B	38	ASN
1	B	80	GLN
1	C	33	GLN
1	C	38	ASN
1	C	80	GLN
1	C	113	ASN
1	C	117	GLN
1	C	139	GLN
1	C	156	GLN
1	D	15	ASN
1	D	28	GLN
1	D	33	GLN
1	D	38	ASN
1	D	39	GLN
1	D	80	GLN

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Mol	Chain	Res	Type
1	D	113	ASN
1	D	119	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](#)

Of 11 ligands modelled in this entry, 8 are monoatomic - leaving 3 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
2	GOL	B	1240	-	5,5,5	4.59	5 (100%)	5,5,5	5.74	3 (60%)
2	GOL	A	1240	-	5,5,5	4.56	5 (100%)	5,5,5	5.81	3 (60%)
2	GOL	D	1241	-	5,5,5	4.51	5 (100%)	5,5,5	5.76	3 (60%)

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	B	1240	-	-	2/4/4/4	-
2	GOL	A	1240	-	-	3/4/4/4	-
2	GOL	D	1241	-	-	2/4/4/4	-

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1240	GOL	C3-C2	-7.69	1.20	1.51
2	A	1240	GOL	C3-C2	-7.43	1.21	1.51
2	D	1241	GOL	C3-C2	-7.36	1.21	1.51
2	A	1240	GOL	O1-C1	4.55	1.61	1.42
2	B	1240	GOL	O1-C1	4.54	1.61	1.42
2	D	1241	GOL	O1-C1	4.51	1.61	1.42
2	A	1240	GOL	O3-C3	3.47	1.57	1.42
2	D	1241	GOL	O3-C3	3.36	1.56	1.42
2	B	1240	GOL	O3-C3	3.12	1.55	1.42
2	D	1241	GOL	C1-C2	-2.87	1.39	1.51
2	A	1240	GOL	O2-C2	-2.85	1.34	1.43
2	B	1240	GOL	O2-C2	-2.84	1.34	1.43
2	A	1240	GOL	C1-C2	-2.83	1.40	1.51
2	D	1241	GOL	O2-C2	-2.80	1.35	1.43
2	B	1240	GOL	C1-C2	-2.78	1.40	1.51

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1240	GOL	O3-C3-C2	10.65	161.25	110.20
2	D	1241	GOL	O3-C3-C2	10.47	160.40	110.20
2	B	1240	GOL	O3-C3-C2	10.44	160.25	110.20
2	D	1241	GOL	O2-C2-C3	6.78	138.99	109.12
2	A	1240	GOL	O2-C2-C3	6.67	138.50	109.12
2	B	1240	GOL	O2-C2-C3	6.67	138.49	109.12
2	B	1240	GOL	O1-C1-C2	3.30	126.02	110.20
2	A	1240	GOL	O1-C1-C2	3.25	125.80	110.20
2	D	1241	GOL	O1-C1-C2	3.18	125.43	110.20

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1240	GOL	O1-C1-C2-C3

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Mol	Chain	Res	Type	Atoms
2	A	1240	GOL	C1-C2-C3-O3
2	B	1240	GOL	C1-C2-C3-O3
2	D	1241	GOL	O1-C1-C2-C3
2	D	1241	GOL	C1-C2-C3-O3
2	B	1240	GOL	O1-C1-C2-O2
2	A	1240	GOL	O1-C1-C2-O2

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1240	GOL	2	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	B	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	B	86:LEU	C	87:PRO	N	1.12

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.