



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 27, 2014 – 03:46 PM GMT

PDB ID : 2P2K
Title : Crystal structure of a lectin from Canavalia gladiata seeds (CGL) in complex with man1-4man-OMe
Authors : Moreno, F.B.M.B.; Bezerra, G.A.; Oliveira, T.M.; Souza, E.P.; Rocha, B.A.M.; Benevides, G.B.; Delatorre, P.; Cavada, B.S.; de Azevedo Jr., W.F.
Deposited on : 2007-03-07
Resolution : 1.98 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

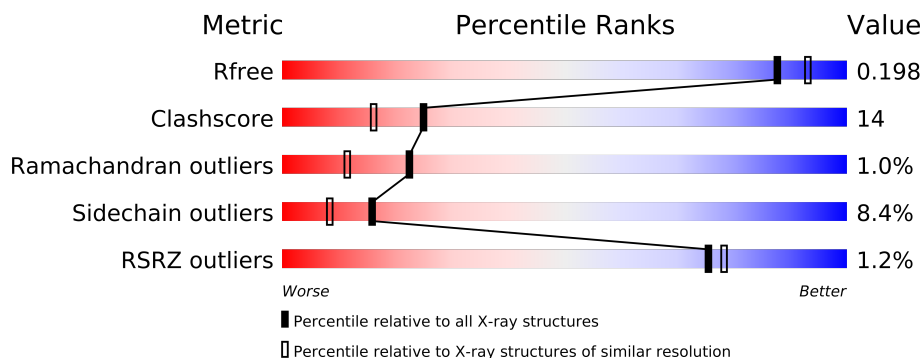
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 1.98 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}	66092	6577 (2.00-1.96)
Clashscore	79885	8091 (2.00-1.96)
Ramachandran outliers	78287	7989 (2.00-1.96)
Sidechain outliers	78261	7987 (2.00-1.96)
RSRZ outliers	66119	6578 (2.00-1.96)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	237	
1	B	237	
1	C	237	
1	D	237	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 7697 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Canavalia gladiata lectin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	234	Total	C	N	O	S	0	0	0
			1779	1123	298	356	2			
1	B	234	Total	C	N	O	S	0	0	0
			1782	1125	298	357	2			
1	C	234	Total	C	N	O	S	0	0	0
			1779	1123	298	356	2			
1	D	234	Total	C	N	O	S	0	0	0
			1782	1125	298	357	2			

- Molecule 2 is a polymer of unknown type called SUGAR (2-MER).

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	O	2	Total	C	O	0	0
			24	13	11		
2	P	2	Total	C	O	0	0
			24	13	11		
2	Q	2	Total	C	O	0	0
			24	13	11		
2	R	2	Total	C	O	0	0
			24	13	11		

- Molecule 3 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

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

- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

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

- Molecule 5 is water.

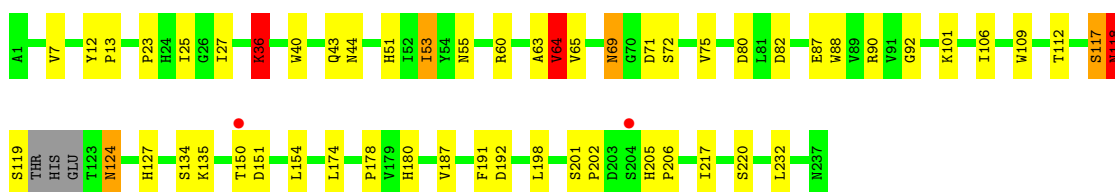
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	149	Total 149	O 149	0	0
5	B	142	Total 142	O 142	0	0
5	C	82	Total 82	O 82	0	0
5	D	89	Total 89	O 89	0	0
5	O	2	Total 2	O 2	0	0
5	P	3	Total 3	O 3	0	0
5	Q	3	Total 3	O 3	0	0
5	R	1	Total 1	O 1	0	0

3 Residue-property plots

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: *Canavalia gladiata* lectin

Chain A: 



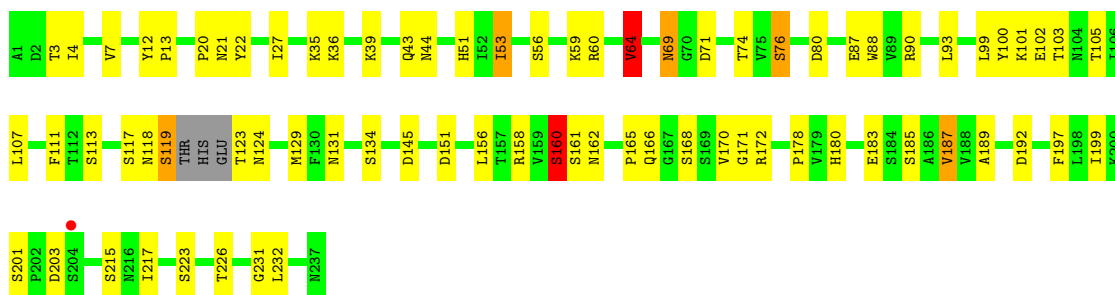
- Molecule 1: *Canavalia gladiata* lectin

Chain B: 



- Molecule 1: *Canavalia gladiata* lectin

Chain C: 



- Molecule 1: *Canavalia gladiata* lectin

Chain D: 





4 Data and refinement statistics

Property	Value	Source
Space group	P 32	Depositor
Cell constants a, b, c, α , β , γ	68.93Å 68.93Å 160.44Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	39.84 – 1.98 39.83 – 1.98	Depositor EDS
% Data completeness (in resolution range)	99.6 (39.84-1.98) 99.5 (39.83-1.98)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	6.80	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.31 (at 1.98Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.182 , 0.236 0.172 , 0.198	Depositor DCC
R_{free} test set	2983 reflections (5.32%)	DCC
Wilson B-factor (Å ²)	19.2	Xtriage
Anisotropy	0.012	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 15.8	EDS
Estimated twinning fraction	0.027 for -h,-k,l 0.486 for h,-h-k,-l 0.028 for -k,-h,-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 59097 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7697	wwPDB-VP
Average B, all atoms (Å ²)	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.16% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CA, MN, MMA, MAN

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.92	2/1819 (0.1%)	0.97	4/2476 (0.2%)
1	B	0.95	2/1822 (0.1%)	1.02	6/2480 (0.2%)
1	C	0.92	1/1820 (0.1%)	1.12	6/2479 (0.2%)
1	D	0.86	2/1822 (0.1%)	1.01	10/2480 (0.4%)
All	All	0.91	7/7283 (0.1%)	1.03	26/9915 (0.3%)

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	2
1	C	0	1
1	D	0	1
2	O	1	0
2	P	1	0
2	Q	1	0
2	R	1	0
All	All	4	4

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	119	SER	C-N	18.07	1.75	1.34
1	B	101	LYS	CE-NZ	7.33	1.67	1.49
1	D	36	LYS	CE-NZ	6.54	1.65	1.49
1	D	36	LYS	CB-CG	6.53	1.70	1.52
1	B	89	VAL	CB-CG1	5.69	1.64	1.52
1	A	36	LYS	CE-NZ	5.19	1.62	1.49
1	A	36	LYS	CD-CE	5.10	1.64	1.51

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	119	SER	O-C-N	-24.89	82.87	122.70
1	C	119	SER	CA-C-N	16.08	152.57	117.20
1	C	119	SER	C-N-CA	10.68	148.39	121.70
1	D	230	LEU	N-CA-C	-10.22	83.40	111.00
1	D	36	LYS	CD-CE-NZ	9.30	133.09	111.70
1	C	64	VAL	CB-CA-C	-8.54	95.17	111.40
1	D	229	LEU	C-N-CA	8.32	142.50	121.70
1	A	64	VAL	CB-CA-C	-8.29	95.65	111.40
1	D	64	VAL	CB-CA-C	-8.01	96.18	111.40
1	B	64	VAL	CB-CA-C	-7.82	96.55	111.40
1	D	174	LEU	CA-CB-CG	7.67	132.95	115.30
1	A	64	VAL	CG1-CB-CG2	7.26	122.52	110.90
1	B	151	ASP	N-CA-C	7.19	130.40	111.00
1	B	174	LEU	CA-CB-CG	7.12	131.68	115.30
1	D	229	LEU	O-C-N	-6.78	111.86	122.70
1	C	64	VAL	CG1-CB-CG2	6.15	120.74	110.90
1	D	64	VAL	CG1-CB-CG2	5.99	120.48	110.90
1	D	229	LEU	CA-C-N	5.95	130.29	117.20
1	D	229	LEU	N-CA-C	5.91	126.94	111.00
1	B	151	ASP	N-CA-CB	-5.80	100.16	110.60
1	D	81	LEU	CB-CG-CD1	5.68	120.65	111.00
1	A	60	ARG	NE-CZ-NH2	-5.46	117.57	120.30
1	A	60	ARG	NE-CZ-NH1	5.46	123.03	120.30
1	C	21	ASN	CB-CA-C	5.33	121.05	110.40
1	B	64	VAL	CG1-CB-CG2	5.10	119.05	110.90
1	B	101	LYS	CD-CE-NZ	-5.03	100.13	111.70

All (4) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	O	240	MMA	C1
2	P	242	MMA	C1
2	Q	244	MMA	C1
2	R	246	MMA	C1

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	116	LYS	Peptide
1	B	149	GLY	Peptide
1	C	160	SER	Peptide

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Group
1	D	116	LYS	Peptide

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1779	0	1730	42	0
1	B	1782	0	1731	48	0
1	C	1779	0	1730	56	0
1	D	1782	0	1730	62	1
2	O	24	0	23	1	0
2	P	24	0	23	0	0
2	Q	24	0	23	0	0
2	R	24	0	23	2	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	149	0	0	5	0
5	B	142	0	0	6	0
5	C	82	0	0	8	0
5	D	89	0	0	6	0
5	O	2	0	0	1	0
5	P	3	0	0	0	1
5	Q	3	0	0	0	0
5	R	1	0	0	0	0
All	All	7697	0	7013	196	1

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 14.

All (196) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:119:SER:C	1:C:123:THR:N	1.75	1.36
1:D:150:THR:HA	1:D:152:GLY:H	1.00	1.12
1:B:150:THR:HA	1:B:152:GLY:H	0.96	1.10
1:B:64:VAL:HG13	1:B:74:THR:HG22	1.33	1.09
1:A:53:ILE:CD1	5:C:300:HOH:O	2.02	1.08
1:D:64:VAL:HG13	1:D:74:THR:HG22	1.34	1.06
1:D:117:SER:HB3	1:D:118:ASN:HA	1.05	1.02
1:B:150:THR:HA	1:B:152:GLY:N	1.76	1.00
1:B:203:ASP:O	1:B:205:HIS:N	1.96	0.96
1:D:117:SER:HB3	1:D:118:ASN:CA	1.95	0.95
1:B:74:THR:HG23	5:B:366:HOH:O	1.62	0.95
1:D:117:SER:CB	1:D:118:ASN:HA	1.95	0.94
1:A:53:ILE:HD11	5:C:300:HOH:O	1.67	0.93
1:C:119:SER:O	1:C:123:THR:N	2.02	0.91
1:D:150:THR:HA	1:D:152:GLY:N	1.85	0.90
1:A:205:HIS:HD2	1:A:206:PRO:O	1.55	0.89
1:C:44:ASN:HD21	1:C:201:SER:H	1.20	0.88
1:A:118:ASN:HB3	5:A:360:HOH:O	1.73	0.88
1:B:134:SER:H	1:B:137:GLN:NE2	1.76	0.83
1:B:117:SER:CB	1:B:118:ASN:HA	2.08	0.82
1:A:118:ASN:OD1	1:A:119:SER:N	2.13	0.81
1:D:150:THR:CA	1:D:152:GLY:H	1.90	0.81
1:A:205:HIS:CD2	1:A:206:PRO:O	2.34	0.81
1:C:88:TRP:HB3	1:C:217:ILE:HD11	1.62	0.81
1:A:44:ASN:HD21	1:A:201:SER:H	1.27	0.81
1:D:99:LEU:HD12	2:O:240:MMA:H4	1.61	0.81
1:B:194:THR:HG22	5:B:316:HOH:O	1.83	0.78
1:A:51:HIS:HB2	1:A:64:VAL:HG23	1.64	0.77
1:C:166:GLN:HG3	5:C:315:HOH:O	1.84	0.77
1:D:44:ASN:HD21	1:D:201:SER:H	1.36	0.74
1:A:36:LYS:HD2	1:A:75:VAL:HG23	1.70	0.74
1:A:124:ASN:HD21	1:B:132:GLN:H	1.35	0.73
1:B:117:SER:HB3	1:B:118:ASN:HA	1.69	0.73
1:B:194:THR:CG2	5:B:316:HOH:O	2.35	0.73
1:B:36:LYS:HE3	5:B:272:HOH:O	1.90	0.72
1:B:203:ASP:C	1:B:205:HIS:H	1.92	0.72
1:D:150:THR:HB	1:D:151:ASP:HA	1.70	0.72
1:C:53:ILE:C	1:C:53:ILE:HD13	2.10	0.71
1:B:117:SER:HB2	1:B:187:VAL:HG12	1.70	0.71
1:A:53:ILE:HD12	5:C:300:HOH:O	1.78	0.71
1:C:107:LEU:O	5:C:321:HOH:O	2.11	0.68
1:A:44:ASN:HD21	1:A:201:SER:N	1.91	0.68
1:D:205:HIS:HD2	1:D:206:PRO:O	1.77	0.68

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:80:ASP:OD1	1:B:82:ASP:HB2	1.95	0.66
1:A:117:SER:HA	1:A:187:VAL:HG22	1.78	0.66
1:B:64:VAL:HG13	1:B:74:THR:CG2	2.21	0.66
1:A:90:ARG:NH1	1:A:217:ILE:HG23	2.11	0.65
1:C:124:ASN:ND2	1:D:131:ASN:H	1.94	0.65
1:B:99:LEU:HD13	1:B:100:TYR:CZ	2.32	0.65
1:A:178:PRO:HB3	1:A:217:ILE:CD1	2.27	0.64
1:D:122:GLU:HG2	5:D:310:HOH:O	1.98	0.64
1:C:131:ASN:O	5:C:321:HOH:O	2.15	0.64
1:C:156:LEU:O	1:C:171:GLY:HA3	1.98	0.63
1:C:51:HIS:HB2	1:C:64:VAL:HG23	1.81	0.63
1:D:44:ASN:ND2	1:D:201:SER:H	1.96	0.63
5:B:352:HOH:O	1:D:194:THR:HG21	1.97	0.63
1:C:160:SER:O	1:C:162:ASN:N	2.31	0.63
1:B:134:SER:H	1:B:137:GLN:HE21	1.46	0.63
1:D:122:GLU:CG	5:D:310:HOH:O	2.48	0.62
1:B:51:HIS:HB2	1:B:64:VAL:HG23	1.82	0.62
1:D:36:LYS:HE3	5:D:243:HOH:O	1.98	0.62
1:D:51:HIS:HB2	1:D:64:VAL:HG23	1.80	0.61
1:C:88:TRP:HB3	1:C:217:ILE:CD1	2.28	0.61
1:D:178:PRO:HB3	1:D:217:ILE:HD11	1.82	0.61
1:C:7:VAL:HG22	1:C:27:ILE:CD1	2.32	0.60
1:C:56:SER:CB	1:C:189:ALA:H	2.15	0.59
1:B:117:SER:HB3	1:B:118:ASN:CA	2.32	0.59
1:B:53:ILE:HD13	5:D:292:HOH:O	2.01	0.59
1:D:201:SER:OG	1:D:203:ASP:HB2	2.03	0.59
1:D:20:PRO:HB2	1:D:22:TYR:CZ	2.38	0.58
1:B:194:THR:HG21	5:D:306:HOH:O	2.02	0.58
1:A:25:ILE:HD12	1:A:75:VAL:HG12	1.86	0.58
1:D:88:TRP:HB3	1:D:217:ILE:CD1	2.34	0.57
1:A:88:TRP:HB3	1:A:217:ILE:HD11	1.87	0.57
1:A:151:ASP:HB2	5:A:361:HOH:O	2.05	0.57
1:C:53:ILE:O	1:C:53:ILE:HD13	2.04	0.56
1:D:99:LEU:HD12	5:O:242:HOH:O	2.05	0.56
1:A:36:LYS:HD2	1:A:75:VAL:CG2	2.34	0.56
1:D:64:VAL:HG13	1:D:74:THR:CG2	2.22	0.56
1:B:88:TRP:HB3	1:B:217:ILE:HD11	1.87	0.56
1:B:25:ILE:HD12	1:B:75:VAL:HG12	1.86	0.56
1:C:20:PRO:HB2	1:C:22:TYR:CZ	2.41	0.56
1:B:137:GLN:HG2	1:B:140:LEU:HD12	1.88	0.56
1:C:69:ASN:HD21	1:C:71:ASP:HB2	1.71	0.56
1:B:134:SER:N	1:B:137:GLN:NE2	2.52	0.55

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:44:ASN:HD21	1:B:201:SER:H	1.52	0.55
1:B:44:ASN:HD21	1:B:201:SER:N	2.04	0.55
1:C:145:ASP:HB3	1:C:158:ARG:HG2	1.88	0.55
1:D:56:SER:CB	1:D:189:ALA:H	2.20	0.55
1:D:111:PHE:CE2	1:D:113:SER:HB2	2.42	0.55
1:D:88:TRP:HB3	1:D:217:ILE:HD11	1.89	0.55
1:C:60:ARG:HD3	1:C:76:SER:OG	2.07	0.55
1:C:100:TYR:HD2	2:R:247:MAN:H62	1.71	0.55
1:D:60:ARG:HD3	1:D:76:SER:OG	2.07	0.55
1:A:178:PRO:HB3	1:A:217:ILE:HD11	1.88	0.55
1:B:90:ARG:NH1	1:B:217:ILE:CG2	2.70	0.54
1:A:44:ASN:ND2	1:A:201:SER:H	2.02	0.54
1:A:151:ASP:CB	5:A:361:HOH:O	2.56	0.54
1:C:44:ASN:ND2	1:C:201:SER:H	1.97	0.53
1:D:90:ARG:NH1	1:D:217:ILE:HG23	2.24	0.52
1:D:145:ASP:HB3	1:D:158:ARG:HG2	1.91	0.52
1:A:124:ASN:ND2	1:B:131:ASN:H	2.07	0.52
1:D:178:PRO:HB3	1:D:217:ILE:CD1	2.39	0.52
1:B:178:PRO:HB3	1:B:217:ILE:CD1	2.40	0.51
1:A:43:GLN:HE22	1:A:69:ASN:HD21	1.57	0.51
1:D:156:LEU:O	1:D:171:GLY:HA3	2.10	0.51
1:C:99:LEU:HD22	2:R:246:MMA:H4	1.92	0.51
1:C:59:LYS:HD3	1:C:80:ASP:HB2	1.92	0.51
1:C:35:LYS:O	1:C:36:LYS:HG2	2.10	0.51
1:D:116:LYS:O	1:D:187:VAL:N	2.44	0.50
1:C:90:ARG:NH1	1:C:217:ILE:HG23	2.26	0.50
1:A:127:HIS:HB3	1:B:127:HIS:HB3	1.93	0.50
1:C:4:ILE:HD13	1:C:215:SER:HB3	1.93	0.50
1:C:43:GLN:HE22	1:C:69:ASN:ND2	2.09	0.50
1:B:99:LEU:HD13	1:B:100:TYR:CE2	2.46	0.49
1:A:112:THR:O	1:A:191:PHE:HA	2.11	0.49
1:B:105:THR:O	1:B:197:PHE:HA	2.12	0.49
1:B:90:ARG:NH1	1:B:217:ILE:HG22	2.27	0.49
1:B:122:GLU:OE2	1:B:123:THR:OG1	2.31	0.49
1:C:102:GLU:C	1:C:165:PRO:HB3	2.33	0.49
1:C:119:SER:HB3	1:D:131:ASN:HB3	1.95	0.48
1:D:54:TYR:CE1	1:D:81:LEU:HD22	2.48	0.48
1:C:124:ASN:HD21	1:D:132:GLN:H	1.61	0.48
5:A:327:HOH:O	1:C:74:THR:HG21	2.14	0.48
1:C:76:SER:HB3	5:C:295:HOH:O	2.12	0.48
1:D:91:VAL:HG12	1:D:214:ILE:HG12	1.95	0.48
1:B:96:SER:OG	1:B:230:LEU:HA	2.13	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:187:VAL:HG11	5:C:292:HOH:O	2.13	0.48
1:D:170:VAL:HG21	1:D:231:GLY:HA2	1.94	0.48
1:D:170:VAL:HB	1:D:226:THR:HG22	1.96	0.48
1:A:51:HIS:O	1:A:63:ALA:HA	2.14	0.47
1:A:69:ASN:HD21	1:A:71:ASP:HB3	1.79	0.47
1:C:170:VAL:HB	1:C:226:THR:HG22	1.97	0.47
1:B:7:VAL:HG21	1:B:52:ILE:HG12	1.96	0.47
1:B:57:VAL:HG21	1:D:64:VAL:HG22	1.96	0.46
1:D:54:TYR:CD1	1:D:81:LEU:HD22	2.51	0.46
1:C:12:TYR:HA	1:C:13:PRO:HD3	1.77	0.46
1:B:118:ASN:OD1	5:B:367:HOH:O	2.21	0.46
1:A:7:VAL:HG22	1:A:27:ILE:HD12	1.97	0.46
1:D:110:SER:HB3	1:D:129:MET:HG3	1.98	0.46
1:C:43:GLN:HE22	1:C:69:ASN:HD21	1.63	0.46
1:A:92:GLY:HA2	1:A:109:TRP:CH2	2.50	0.46
1:C:158:ARG:HB2	1:C:166:GLN:HB3	1.98	0.46
1:C:178:PRO:HB3	1:C:217:ILE:CD1	2.46	0.45
1:C:217:ILE:CD1	1:D:138:LYS:HD2	2.46	0.45
1:A:12:TYR:HA	1:A:13:PRO:HD3	1.83	0.45
1:A:80:ASP:OD1	1:A:82:ASP:HB2	2.17	0.45
1:B:133:PHE:HA	1:B:137:GLN:NE2	2.31	0.45
1:C:22:TYR:CD2	1:C:39:LYS:HE3	2.52	0.45
1:C:111:PHE:CE2	1:C:113:SER:HB2	2.51	0.45
1:C:107:LEU:HD12	1:C:107:LEU:N	2.31	0.45
1:A:69:ASN:ND2	1:A:71:ASP:H	2.15	0.45
1:D:205:HIS:CD2	1:D:206:PRO:O	2.65	0.44
1:D:102:GLU:C	1:D:165:PRO:HB3	2.38	0.44
1:C:36:LYS:HE3	1:C:76:SER:O	2.18	0.44
1:B:112:THR:O	1:B:191:PHE:HA	2.18	0.43
1:A:106:ILE:HB	1:A:154:LEU:HB3	2.00	0.43
1:A:174:LEU:N	1:A:174:LEU:HD12	2.32	0.43
1:D:208:ASP:OD2	1:D:227:GLY:HA2	2.19	0.43
1:B:51:HIS:O	1:B:63:ALA:HA	2.19	0.43
1:C:87:GLU:HG3	1:C:180:HIS:CD2	2.54	0.43
1:D:106:ILE:HB	1:D:154:LEU:HB3	2.01	0.43
1:D:15:THR:HG21	1:D:21:ASN:HD22	1.84	0.43
1:D:116:LYS:HG3	1:D:188:VAL:HB	2.01	0.43
1:C:105:THR:O	1:C:197:PHE:HA	2.19	0.43
1:C:129:MET:HE3	1:D:123:THR:HB	2.01	0.42
1:A:90:ARG:NH1	1:A:217:ILE:CG2	2.80	0.42
1:D:88:TRP:HB3	1:D:217:ILE:HD12	2.01	0.42
1:D:50:ALA:O	1:D:194:THR:HA	2.20	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
5:A:315:HOH:O	1:C:118:ASN:HA	2.20	0.42
1:A:65:VAL:O	1:A:72:SER:HA	2.20	0.42
1:C:170:VAL:HG21	1:C:231:GLY:HA2	2.00	0.42
1:D:126:LEU:CD2	1:D:179:VAL:HG22	2.49	0.42
1:C:53:ILE:C	1:C:53:ILE:CD1	2.84	0.42
1:C:124:ASN:HA	1:D:129:MET:O	2.20	0.42
1:A:69:ASN:C	1:A:69:ASN:HD22	2.22	0.42
1:A:198:LEU:HD23	1:A:198:LEU:C	2.39	0.41
1:D:135:LYS:HG3	1:D:149:GLY:O	2.20	0.41
1:B:59:LYS:HD3	1:B:80:ASP:HA	2.02	0.41
1:D:150:THR:CB	1:D:151:ASP:HA	2.44	0.41
1:D:149:GLY:HA2	1:D:150:THR:O	2.21	0.41
1:A:55:ASN:HD21	1:C:53:ILE:HD12	1.85	0.41
1:A:87:GLU:HG2	1:A:180:HIS:CD2	2.55	0.41
1:C:93:LEU:O	1:C:172:ARG:HA	2.21	0.41
1:D:93:LEU:O	1:D:172:ARG:HA	2.20	0.41
1:C:103:THR:O	1:C:199:ILE:HA	2.20	0.41
1:C:3:THR:O	1:C:215:SER:HA	2.21	0.41
1:D:15:THR:CG2	1:D:21:ASN:HD22	2.33	0.41
1:B:106:ILE:HB	1:B:154:LEU:HB3	2.02	0.40
1:D:89:VAL:HG22	1:D:181:ILE:HB	2.03	0.40
1:B:187:VAL:HG21	5:D:299:HOH:O	2.21	0.40
1:D:107:LEU:N	1:D:107:LEU:HD12	2.36	0.40
1:B:64:VAL:CG1	1:B:74:THR:HG22	2.25	0.40
1:C:129:MET:O	1:D:124:ASN:HA	2.20	0.40
1:B:111:PHE:HB3	1:B:128:PHE:CZ	2.56	0.40
1:A:23:PRO:HB2	1:A:40:TRP:O	2.22	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	Distance(Å)	Clash(Å)
1:D:76:SER:CB	5:P:246:HOH:O[2.664]	1.29	0.91

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	230/237 (97%)	219 (95%)	9 (4%)	2 (1%)	25	13
1	B	230/237 (97%)	219 (95%)	8 (4%)	3 (1%)	18	7
1	C	232/237 (98%)	219 (94%)	11 (5%)	2 (1%)	25	13
1	D	230/237 (97%)	219 (95%)	9 (4%)	2 (1%)	25	13
All	All	922/948 (97%)	876 (95%)	37 (4%)	9 (1%)	22	11

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	151	ASP
1	B	204	SER
1	C	161	SER
1	C	203	ASP
1	D	203	ASP
1	B	203	ASP
1	A	202	PRO
1	D	151	ASP
1	A	118	ASN

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	199/202 (98%)	185 (93%)	14 (7%)	21	13
1	B	199/202 (98%)	185 (93%)	14 (7%)	21	13
1	C	199/202 (98%)	183 (92%)	16 (8%)	17	9
1	D	199/202 (98%)	176 (88%)	23 (12%)	8	3
All	All	796/808 (98%)	729 (92%)	67 (8%)	16	8

All (67) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	36	LYS
1	A	53	ILE
1	A	64	VAL

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	69	ASN
1	A	101	LYS
1	A	117	SER
1	A	118	ASN
1	A	124	ASN
1	A	134	SER
1	A	135	LYS
1	A	150	THR
1	A	192	ASP
1	A	220	SER
1	A	232	LEU
1	B	64	VAL
1	B	81	LEU
1	B	99	LEU
1	B	101	LYS
1	B	107	LEU
1	B	117	SER
1	B	151	ASP
1	B	174	LEU
1	B	184	SER
1	B	192	ASP
1	B	194	THR
1	B	216	ASN
1	B	230	LEU
1	B	232	LEU
1	C	53	ILE
1	C	64	VAL
1	C	69	ASN
1	C	76	SER
1	C	101	LYS
1	C	117	SER
1	C	134	SER
1	C	151	ASP
1	C	160	SER
1	C	168	SER
1	C	183	GLU
1	C	185	SER
1	C	187	VAL
1	C	192	ASP
1	C	223	SER
1	C	232	LEU
1	D	56	SER

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	D	64	VAL
1	D	71	ASP
1	D	76	SER
1	D	81	LEU
1	D	101	LYS
1	D	117	SER
1	D	122	GLU
1	D	132	GLN
1	D	134	SER
1	D	135	LYS
1	D	161	SER
1	D	164	SER
1	D	168	SER
1	D	174	LEU
1	D	183	GLU
1	D	185	SER
1	D	192	ASP
1	D	194	THR
1	D	208	ASP
1	D	223	SER
1	D	230	LEU
1	D	232	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	21	ASN
1	A	41	ASN
1	A	44	ASN
1	A	69	ASN
1	A	83	ASN
1	A	104	ASN
1	A	124	ASN
1	A	166	GLN
1	A	205	HIS
1	B	41	ASN
1	B	44	ASN
1	B	104	ASN
1	B	118	ASN
1	B	137	GLN
1	B	166	GLN
1	C	41	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	C	44	ASN
1	C	69	ASN
1	C	124	ASN
1	C	131	ASN
1	D	21	ASN
1	D	44	ASN
1	D	104	ASN
1	D	132	GLN
1	D	205	HIS

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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

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

5.5 Carbohydrates ⓘ

8 carbohydrates are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	MMA	O	240	2	13,13,13	0.94	1 (7%)	18,18,18	1.27	3 (16%)
2	MAN	O	241	2	10,11,12	0.87	0	11,15,17	1.45	2 (18%)
2	MMA	P	242	2	13,13,13	0.59	0	18,18,18	3.73	7 (38%)
2	MAN	P	243	2	10,11,12	0.49	0	11,15,17	0.90	0
2	MMA	Q	244	2	13,13,13	0.66	0	18,18,18	3.76	8 (44%)
2	MAN	Q	245	2	10,11,12	0.53	0	11,15,17	1.16	1 (9%)
2	MMA	R	246	2	13,13,13	0.97	1 (7%)	18,18,18	1.38	2 (11%)
2	MAN	R	247	2	10,11,12	0.73	0	11,15,17	1.38	2 (18%)

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	MMA	O	240	2	1/1/5/5	0/4/24/24	0/1/1/1
2	MAN	O	241	2	-	0/2/19/22	0/1/1/1
2	MMA	P	242	2	1/1/5/5	0/4/24/24	0/1/1/1
2	MAN	P	243	2	-	0/2/19/22	0/1/1/1
2	MMA	Q	244	2	1/1/5/5	0/4/24/24	0/1/1/1
2	MAN	Q	245	2	-	0/2/19/22	0/1/1/1
2	MMA	R	246	2	1/1/5/5	0/4/24/24	0/1/1/1
2	MAN	R	247	2	-	0/2/19/22	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	O	240	MMA	O1-C1	2.56	1.44	1.40
2	R	246	MMA	O1-C1	2.47	1.44	1.40

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	Q	244	MMA	O1-C1-C2	9.63	119.98	108.18
2	P	242	MMA	O1-C1-C2	8.59	118.71	108.18
2	P	242	MMA	C1-C2-C3	-7.77	94.88	110.00
2	Q	244	MMA	C1-C2-C3	-7.68	95.05	110.00
2	P	242	MMA	C1-O5-C5	-7.53	99.09	113.73
2	Q	244	MMA	C1-O5-C5	-7.03	100.05	113.73
2	P	242	MMA	O6-C6-C5	-3.89	97.97	111.36
2	R	246	MMA	O5-C5-C4	3.84	116.87	109.76
2	Q	244	MMA	O2-C2-C1	3.57	117.82	110.04
2	P	242	MMA	O2-C2-C1	3.44	117.54	110.04
2	P	242	MMA	O5-C1-O1	3.31	118.75	110.85
2	O	241	MAN	O5-C5-C4	2.78	114.18	110.65
2	Q	244	MMA	O2-C2-C3	-2.74	104.22	110.35
2	O	240	MMA	C3-C4-C5	2.62	114.87	110.20
2	Q	244	MMA	O6-C6-C5	-2.58	102.48	111.36
2	O	241	MAN	O3-C3-C4	2.58	116.13	110.35
2	P	242	MMA	O5-C5-C6	2.57	112.65	106.34
2	Q	244	MMA	O5-C1-O1	2.40	116.58	110.85
2	O	240	MMA	O1-C1-C2	2.33	111.04	108.18
2	O	240	MMA	O5-C1-C2	-2.26	105.67	110.31

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	R	247	MAN	O3-C3-C4	2.19	115.26	110.35
2	Q	245	MAN	O3-C3-C2	-2.18	105.95	109.94
2	R	246	MMA	C1-O5-C5	2.07	117.75	113.73
2	R	247	MAN	O5-C5-C6	2.04	109.13	106.98
2	Q	244	MMA	O5-C5-C6	2.01	111.29	106.34

All (4) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	Q	244	MMA	C1
2	P	242	MMA	C1
2	R	246	MMA	C1
2	O	240	MMA	C1

There are no torsion outliers.

There are no ring outliers.

5.6 Ligand geometry ⓘ

Of 8 ligands modelled in this entry, 8 are monoatomic - leaving 0 for Mogul analysis.

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.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	234/237 (98%)	-0.30	2 (0%) 81 84	8, 14, 29, 46	6 (2%)
1	B	234/237 (98%)	-0.28	2 (0%) 81 84	9, 14, 28, 42	4 (1%)
1	C	234/237 (98%)	-0.08	1 (0%) 90 92	11, 20, 37, 45	6 (2%)
1	D	234/237 (98%)	-0.11	4 (1%) 67 69	9, 20, 36, 45	6 (2%)
All	All	936/948 (98%)	-0.19	9 (0%) 75 82	8, 17, 32, 46	22 (2%)

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	204	SER	5.2
1	B	149	GLY	3.2
1	A	204	SER	2.6
1	D	1	ALA	2.5
1	D	203	ASP	2.3
1	D	204	SER	2.2
1	D	205	HIS	2.1
1	A	150	THR	2.1
1	B	150	THR	2.0

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

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

6.3 Carbohydrates ⓘ

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	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	MMA	O	240	13/13	0.31	6.66	42,53,59,60	0
2	MMA	R	246	13/13	0.37	4.52	40,52,58,59	0
2	MAN	R	247	11/12	0.13	0.73	19,29,33,34	0
2	MAN	O	241	11/12	0.13	0.58	21,27,32,33	0
2	MMA	Q	244	13/13	0.10	0.40	18,24,26,27	0
2	MAN	P	243	11/12	0.09	0.20	8,13,15,16	0
2	MAN	Q	245	11/12	0.08	-0.15	9,14,16,18	0
2	MMA	P	242	13/13	0.10	-0.22	18,21,26,28	0

6.4 Ligands ⓘ

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	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	MN	B	238	1/1	0.08	1.76	16,16,16,16	0
3	MN	A	238	1/1	0.08	0.85	16,16,16,16	0
4	CA	D	239	1/1	0.06	-2.06	17,17,17,17	0
4	CA	B	239	1/1	0.06	-2.34	11,11,11,11	0
4	CA	C	239	1/1	0.06	-2.90	17,17,17,17	0
3	MN	D	238	1/1	0.04	-3.18	21,21,21,21	0
3	MN	C	238	1/1	0.04	-4.32	21,21,21,21	0
4	CA	A	239	1/1	0.04	-6.35	12,12,12,12	0

6.5 Other polymers ⓘ

There are no such residues in this entry.