Full wwPDB X-ray Structure Validation Report (i)

Jun 11, 2024 – 07:23 PM EDT

PDB ID	:	1LEN
Title	:	REFINEMENT OF TWO CRYSTAL FORMS OF LENTIL LECTIN AT 1.8 ANGSTROMS RESOLUTION
Authors	:	Van Overberge, D.; Loris, R.; Wyns, L.
Deposited on	:	1993-11-17
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:

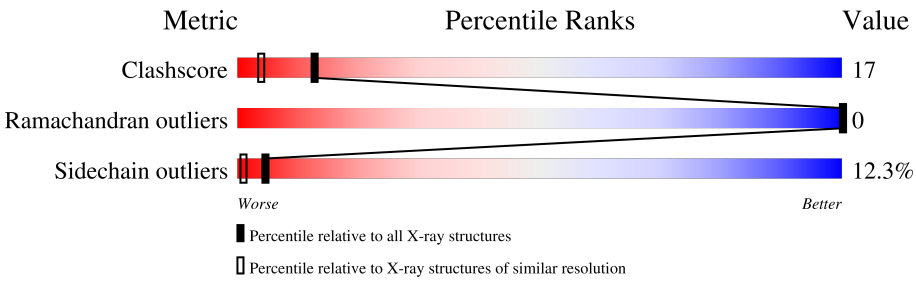
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

1 Overall quality at a glance (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))
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (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%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	181	69% 24% 6% •
1	C	181	59% 32% 8% •
2	B	52	50% 35% 6% 10%
2	D	52	52% 35% • 10%

2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 3820 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.

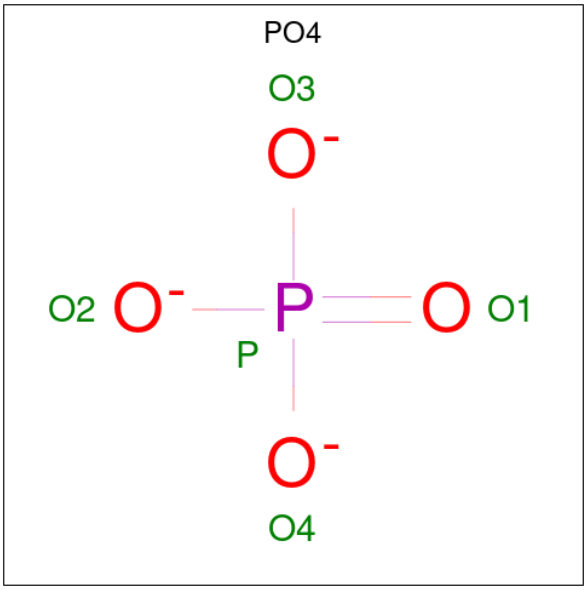
- Molecule 1 is a protein called LECTIN.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
1	A	181	1415	903	231	281	0	3	0
1	C	181	1420	906	232	282	0	2	0

- Molecule 2 is a protein called LECTIN.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	B	47	370	238	60	72	0	1	0
2	D	47	370	238	60	72	0	1	0

- Molecule 3 is PHOSPHATE ION (three-letter code: PO4) (formula: O_4P).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total O P 5 4 1	0	0
3	C	1	Total O P 5 4 1	0	0

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

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

- Molecule 5 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total Ca 1 1	0	0
5	C	1	Total Ca 1 1	0	0

- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	101	Total O 101 101	0	0
6	B	27	Total O 27 27	0	0
6	C	88	Total O 88 88	0	0
6	D	15	Total O 15 15	0	0

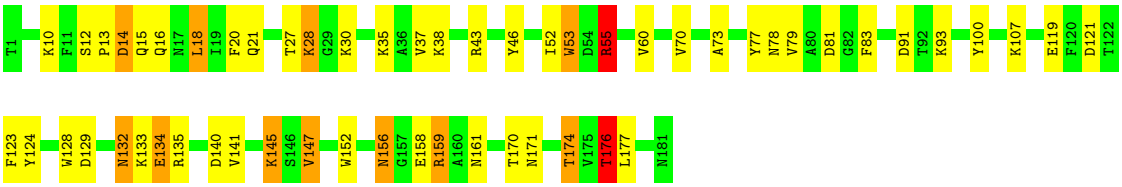
3 Residue-property plots (i)

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. 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.

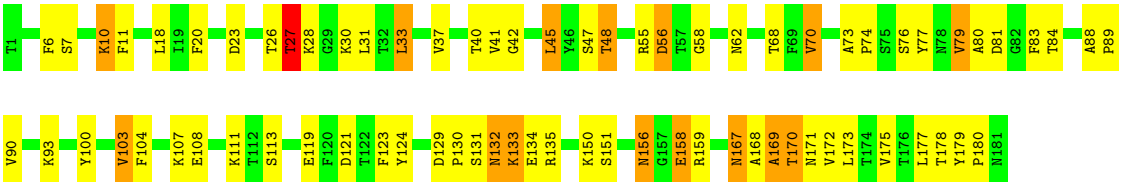
• Molecule 1: LECTIN

Chain A:



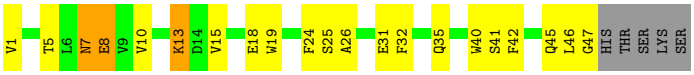
• Molecule 1: LECTIN

Chain C:



• Molecule 2: LECTIN

Chain B:



• Molecule 2: LECTIN

Chain D:

4 Data and refinement statistics (i)

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

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	58.49Å 56.37Å 82.73Å 90.00° 104.40° 90.00°	Depositor
Resolution (Å)	(Not available) – 1.80	Depositor
% Data completeness (in resolution range)	(Not available) ((Not available)-1.80)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	RESTRAIN	Depositor
R, R_{free}	0.175 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3820	wwPDB-VP
Average B, all atoms ($Å^2$)	26.0	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, PO4, 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	1.19	4/1465 (0.3%)	1.82	22/2002 (1.1%)
1	C	1.17	2/1464 (0.1%)	1.78	30/1999 (1.5%)
2	B	1.30	1/385 (0.3%)	2.09	15/527 (2.8%)
2	D	1.26	0/385	1.75	7/527 (1.3%)
All	All	1.20	7/3699 (0.2%)	1.83	74/5055 (1.5%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	158	GLU	CD-OE2	7.52	1.33	1.25
1	A	158	GLU	CD-OE1	7.30	1.33	1.25
1	C	158	GLU	CD-OE1	7.26	1.33	1.25
2	B	19	TRP	NE1-CE2	-6.49	1.29	1.37
1	C	158	GLU	CD-OE2	5.96	1.32	1.25
1	A	53	TRP	NE1-CE2	-5.36	1.30	1.37
1	A	152	TRP	NE1-CE2	-5.36	1.30	1.37

All (74) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	129	ASP	CB-CG-OD2	10.89	128.10	118.30
2	B	31	GLU	OE1-CD-OE2	-10.67	110.49	123.30
1	A	100	TYR	CB-CG-CD2	-10.15	114.91	121.00
1	A	79	VAL	CA-CB-CG2	9.75	125.53	110.90
2	B	8	GLU	OE1-CD-OE2	-8.60	112.98	123.30
1	A	147	VAL	CA-CB-CG2	8.50	123.65	110.90
1	C	119	GLU	OE1-CD-OE2	-8.49	113.12	123.30
1	C	79	VAL	CA-CB-CG2	8.35	123.43	110.90
2	D	15	VAL	CA-CB-CG2	8.25	123.28	110.90
2	B	24	PHE	CG-CD1-CE1	8.22	129.84	120.80

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	47	GLY	CA-C-O	8.13	135.24	120.60
2	B	10	VAL	CA-CB-CG2	7.77	122.56	110.90
1	C	83	PHE	CB-CG-CD1	-7.62	115.46	120.80
2	B	1	VAL	CA-CB-CG1	7.57	122.26	110.90
1	C	100	TYR	CB-CG-CD2	-7.35	116.59	121.00
1	A	18	LEU	CB-CG-CD1	7.30	123.41	111.00
2	B	15	VAL	CA-CB-CG2	7.26	121.79	110.90
1	C	45	LEU	CB-CG-CD2	7.23	123.29	111.00
1	A	20	PHE	CG-CD1-CE1	-7.17	112.92	120.80
1	C	121	ASP	CB-CG-OD2	7.17	124.75	118.30
1	C	169	ALA	C-N-CA	7.08	139.39	121.70
1	A	158	GLU	OE1-CD-OE2	-7.03	114.87	123.30
1	C	130	PRO	C-N-CA	7.00	139.21	121.70
1	C	103	VAL	CA-CB-CG2	6.98	121.37	110.90
1	C	129	ASP	CB-CG-OD2	6.91	124.52	118.30
1	C	11	PHE	CG-CD1-CE1	6.85	128.33	120.80
1	A	55	ARG	C-N-CA	6.79	138.68	121.70
1	A	83	PHE	CG-CD1-CE1	-6.77	113.36	120.80
2	D	1	VAL	CA-CB-CG1	6.76	121.04	110.90
1	A	141	VAL	CA-CB-CG2	6.73	121.00	110.90
2	B	24	PHE	CD1-CE1-CZ	-6.63	112.15	120.10
1	C	90	VAL	O-C-N	-6.47	112.35	122.70
1	C	135	ARG	NE-CZ-NH1	6.43	123.52	120.30
1	A	119	GLU	OE1-CD-OE2	-6.42	115.60	123.30
1	C	173	LEU	O-C-N	-6.28	112.66	122.70
1	A	100	TYR	CG-CD1-CE1	-6.27	116.28	121.30
1	C	58	GLY	O-C-N	-6.19	112.80	122.70
1	C	123	PHE	CG-CD1-CE1	-6.12	114.07	120.80
2	B	24	PHE	CB-CG-CD1	6.03	125.02	120.80
1	A	159	ARG	O-C-N	6.02	132.33	122.70
1	A	140	ASP	CB-CG-OD1	5.98	123.68	118.30
1	C	37	VAL	CA-CB-CG2	5.83	119.65	110.90
1	A	123	PHE	CG-CD1-CE1	-5.78	114.44	120.80
2	B	24	PHE	CZ-CE2-CD2	5.77	127.03	120.10
1	C	23	ASP	C-N-CA	-5.72	110.28	122.30
2	D	8	GLU	OE1-CD-OE2	-5.68	116.48	123.30
1	C	175	VAL	CA-CB-CG2	5.68	119.41	110.90
2	D	24	PHE	CB-CG-CD2	-5.65	116.84	120.80
1	C	47	SER	CA-C-N	5.65	129.63	117.20
1	C	27	THR	O-C-N	5.64	131.72	122.70
1	A	176	THR	CA-CB-CG2	5.61	120.25	112.40
2	B	7	ASN	N-CA-CB	5.60	120.69	110.60

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	18	LEU	CB-CG-CD1	5.56	120.46	111.00
1	C	56	ASP	CB-CG-OD2	-5.54	113.31	118.30
2	B	31	GLU	CG-CD-OE1	5.54	129.37	118.30
2	D	47	GLY	CA-C-O	5.50	130.50	120.60
1	C	56	ASP	CB-CG-OD1	5.46	123.21	118.30
2	B	46	LEU	CA-C-N	5.38	126.96	116.20
2	B	40	TRP	C-N-CA	5.36	135.10	121.70
1	C	70	VAL	CG1-CB-CG2	-5.35	102.34	110.90
1	C	33	LEU	CB-CG-CD1	5.34	120.08	111.00
1	A	46	TYR	O-C-N	-5.31	114.20	122.70
2	D	31	GLU	OE1-CD-OE2	-5.31	116.93	123.30
1	C	84	THR	O-C-N	5.29	131.16	122.70
1	A	14	ASP	N-CA-CB	5.28	120.10	110.60
1	C	11	PHE	CD1-CE1-CZ	-5.26	113.79	120.10
2	B	45	GLN	CG-CD-OE1	-5.19	111.22	121.60
1	A	55	ARG	O-C-N	5.18	130.99	122.70
1	A	121	ASP	CB-CG-OD2	5.17	122.95	118.30
2	D	40	TRP	CH2-CZ2-CE2	5.12	122.52	117.40
1	C	80	ALA	CA-C-O	-5.08	109.44	120.10
1	A	83	PHE	CB-CG-CD1	-5.06	117.26	120.80
1	C	170	THR	OG1-CB-CG2	5.04	121.60	110.00
1	A	176	THR	N-CA-CB	5.04	119.87	110.30

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	1415	0	1363	40	0
1	C	1420	0	1367	61	0
2	B	370	0	354	17	0
2	D	370	0	354	11	0
3	A	5	0	0	0	0
3	C	5	0	0	0	0
4	A	1	0	0	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	C	1	0	0	0	0
5	A	1	0	0	0	0
5	C	1	0	0	0	0
6	A	101	0	0	4	0
6	B	27	0	0	0	0
6	C	88	0	0	2	0
6	D	15	0	0	0	0
All	All	3820	0	3438	117	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 17.

All (117) 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)
2:D:43:ASN:HD22	2:D:43:ASN:C	1.64	0.99
2:D:43:ASN:C	2:D:43:ASN:ND2	2.20	0.91
1:A:55:ARG:HH11	1:A:55:ARG:HG2	1.34	0.91
1:C:73:ALA:H	1:C:156:ASN:HD21	1.10	0.90
1:C:170:THR:HG23	1:C:172:VAL:HG23	1.53	0.88
1:A:132:ASN:HD22	1:A:134:GLU:H	1.20	0.85
1:A:73:ALA:H	1:A:156:ASN:HD21	1.23	0.85
1:C:10:LYS:NZ	1:C:10:LYS:HB2	1.95	0.82
2:D:43:ASN:ND2	2:D:44:SER:N	2.29	0.80
1:C:132:ASN:ND2	1:C:134:GLU:H	1.80	0.77
1:C:132:ASN:HD22	1:C:133:LYS:N	1.83	0.76
2:B:13:LYS:NZ	2:B:13:LYS:CB	2.48	0.76
1:A:21:GLN:HE22	1:A:43:ARG:HH21	1.33	0.76
1:C:167:ASN:HD22	1:C:167:ASN:C	1.91	0.74
1:A:132:ASN:ND2	1:A:134:GLU:H	1.87	0.73
1:C:170:THR:HG23	1:C:172:VAL:CG2	2.19	0.72
1:A:55:ARG:HG2	1:A:55:ARG:NH1	2.02	0.72
1:A:128:TRP:HB2	1:A:145:LYS:HG2	1.71	0.72
1:C:177:LEU:C	1:C:177:LEU:HD23	2.11	0.71
6:A:273:HOH:O	1:C:48:THR:HG23	1.91	0.71
1:C:132:ASN:HD22	1:C:132:ASN:C	1.95	0.70
1:C:27:THR:O	1:C:27:THR:CG2	2.40	0.70
1:C:76:SER:OG	1:C:77[B]:TYR:HD1	1.75	0.70
1:C:158:GLU:HG3	1:C:159:ARG:N	2.07	0.69
1:C:73:ALA:H	1:C:156:ASN:ND2	1.89	0.69
1:A:27:THR:O	1:A:27:THR:HG23	1.93	0.67

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:55:ARG:HB2	2:D:18:GLU:OE2	1.93	0.67
2:B:26:ALA:HB1	2:B:35[B]:GLN:HG3	1.77	0.66
1:A:77[B]:TYR:HD1	1:A:78:ASN:N	1.92	0.66
1:C:76:SER:OG	1:C:77[B]:TYR:CD1	2.48	0.65
1:A:77[B]:TYR:CD1	1:A:78:ASN:N	2.66	0.64
1:A:132:ASN:ND2	1:A:134:GLU:HB2	2.13	0.64
2:B:13:LYS:NZ	2:B:13:LYS:HB2	2.13	0.62
1:C:159:ARG:HD3	2:D:38:HIS:ND1	2.16	0.61
1:A:73:ALA:H	1:A:156:ASN:ND2	1.97	0.61
1:C:27:THR:O	1:C:27:THR:HG22	2.00	0.60
2:D:43:ASN:HD22	2:D:44:SER:N	1.95	0.60
1:A:174:THR:HG23	6:A:230:HOH:O	2.01	0.59
1:C:68:THR:HG21	1:C:159:ARG:HD2	1.84	0.59
2:B:13:LYS:HB2	2:B:13:LYS:HZ1	1.67	0.59
1:C:170:THR:CG2	1:C:172:VAL:HB	2.34	0.58
1:C:167:ASN:ND2	1:C:169:ALA:H	2.02	0.57
1:A:147:VAL:HG23	2:B:8:GLU:HG2	1.86	0.57
1:C:180:PRO:HD2	6:C:238:HOH:O	2.03	0.57
1:C:73:ALA:HB2	1:C:79:VAL:HG13	1.86	0.56
1:A:176:THR:HG23	6:A:246:HOH:O	2.05	0.56
1:C:20:PHE:HE1	1:C:31:LEU:CD1	2.19	0.56
1:C:167:ASN:C	1:C:167:ASN:ND2	2.60	0.55
1:A:38:LYS:HG2	2:B:32:PHE:CD2	2.42	0.54
1:C:170:THR:HG22	1:C:172:VAL:H	1.72	0.54
1:C:177:LEU:HD23	1:C:178:THR:N	2.22	0.54
1:C:132:ASN:ND2	1:C:132:ASN:C	2.60	0.53
1:C:40:THR:HG22	2:D:28:THR:OG1	2.09	0.53
1:C:70:VAL:CG2	1:C:159:ARG:HG3	2.40	0.52
1:A:147:VAL:CG2	2:B:8:GLU:HG2	2.40	0.52
1:C:158:GLU:HG3	1:C:159:ARG:H	1.72	0.52
1:C:70:VAL:HG23	1:C:159:ARG:HG2	1.92	0.51
1:C:73:ALA:N	1:C:156:ASN:HD21	1.93	0.51
1:A:27:THR:O	1:A:27:THR:CG2	2.58	0.51
1:C:10:LYS:HB2	1:C:10:LYS:HZ2	1.76	0.51
1:A:12:SER:HB2	1:A:15:GLN:OE1	2.12	0.50
1:A:21:GLN:HE21	1:A:43:ARG:HE	1.60	0.50
2:D:26:ALA:HB1	2:D:35[B]:GLN:HG3	1.94	0.49
1:C:88:ALA:HB1	1:C:89:PRO:HD2	1.95	0.48
1:C:33:LEU:O	1:C:42:GLY:HA3	2.13	0.48
2:D:43:ASN:HD21	2:D:45:GLN:HG3	1.79	0.48
2:B:13:LYS:NZ	2:B:13:LYS:HB3	2.26	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:10:LYS:HB3	1:C:10:LYS:HE2	1.64	0.48
1:A:13:PRO:HD2	1:A:14:ASP:H	1.79	0.47
2:B:13:LYS:HB3	2:B:13:LYS:HZ2	1.80	0.47
1:A:38:LYS:HG2	2:B:32:PHE:HD2	1.80	0.47
1:C:167:ASN:HD21	1:C:169:ALA:HB3	1.79	0.46
1:A:55:ARG:NH1	1:A:55:ARG:CG	2.76	0.46
1:A:161:ASN:HB3	6:A:253:HOH:O	2.16	0.46
1:A:124:TYR:CD1	1:A:133:LYS:HA	2.50	0.46
1:C:6:PHE:C	1:C:6:PHE:CD1	2.88	0.45
1:C:28:LYS:HB3	1:C:28:LYS:HE2	1.59	0.45
1:C:103:VAL:HG23	1:C:104:PHE:CD2	2.52	0.45
1:C:170:THR:CG2	1:C:172:VAL:H	2.29	0.45
1:C:76:SER:C	1:C:77[B]:TYR:HD1	2.19	0.45
1:A:70:VAL:HG22	1:A:159:ARG:HA	1.99	0.45
1:C:62:ASN:HD22	1:C:168:ALA:H	1.65	0.45
1:A:177:LEU:C	1:A:177:LEU:HD23	2.38	0.45
1:C:132:ASN:HD22	1:C:134:GLU:H	1.63	0.45
1:C:170:THR:O	1:C:171:ASN:HB2	2.17	0.45
1:A:28:LYS:HB2	1:A:28:LYS:HE3	1.39	0.44
1:C:70:VAL:HG23	1:C:159:ARG:CG	2.48	0.44
1:C:177:LEU:C	1:C:177:LEU:CD2	2.81	0.44
1:A:176:THR:HG22	2:B:5:THR:HG23	2.00	0.43
2:D:28:THR:HB	2:D:32:PHE:HB3	2.00	0.43
1:C:73:ALA:HB1	1:C:74:PRO:HD2	2.00	0.43
1:C:41:VAL:HB	2:D:27:THR:HG22	2.01	0.43
1:C:88:ALA:HB1	1:C:89:PRO:CD	2.47	0.43
1:C:179:TYR:HA	1:C:180:PRO:HD3	1.62	0.43
2:B:13:LYS:CB	2:B:13:LYS:HZ2	2.27	0.43
1:C:124:TYR:CD1	1:C:133:LYS:HA	2.53	0.43
1:A:135:ARG:HH11	1:A:135:ARG:HD2	1.66	0.43
1:C:70:VAL:CG2	1:C:159:ARG:CG	2.97	0.42
1:A:53:TRP:CD2	2:B:13:LYS:HG3	2.53	0.42
1:A:55:ARG:HB2	2:B:18:GLU:OE2	2.20	0.42
1:C:167:ASN:HD22	1:C:169:ALA:H	1.65	0.42
2:B:13:LYS:CB	2:B:13:LYS:HZ1	2.24	0.42
1:C:124:TYR:CE2	1:C:133:LYS:HD3	2.55	0.42
1:C:170:THR:CG2	1:C:172:VAL:CB	2.98	0.42
1:C:124:TYR:HD1	6:C:195:HOH:O	2.03	0.42
1:A:60:VAL:CG2	2:B:13:LYS:HE3	2.50	0.41
1:A:170:THR:O	1:A:171:ASN:HB2	2.21	0.41
1:A:21:GLN:NE2	1:A:43:ARG:HH21	2.09	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:21:GLN:NE2	1:A:43:ARG:HE	2.18	0.41
1:C:20:PHE:CZ	1:C:26:THR:HG23	2.56	0.41

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	182/181 (101%)	179 (98%)	3 (2%)	0	100	100
1	C	181/181 (100%)	176 (97%)	5 (3%)	0	100	100
2	B	46/52 (88%)	46 (100%)	0	0	100	100
2	D	46/52 (88%)	45 (98%)	1 (2%)	0	100	100
All	All	455/466 (98%)	446 (98%)	9 (2%)	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	158/156 (101%)	138 (87%)	20 (13%)	4	1
1	C	158/156 (101%)	138 (87%)	20 (13%)	4	1
2	B	41/45 (91%)	37 (90%)	4 (10%)	8	2

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
2	D	41/45 (91%)	36 (88%)	5 (12%)	5 1
All	All	398/402 (99%)	349 (88%)	49 (12%)	4 1

All (49) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	LYS
1	A	16	GLN
1	A	18	LEU
1	A	28	LYS
1	A	30	LYS
1	A	35	LYS
1	A	37[A]	VAL
1	A	37[B]	VAL
1	A	52	ILE
1	A	55	ARG
1	A	81	ASP
1	A	91	ASP
1	A	93	LYS
1	A	107	LYS
1	A	132	ASN
1	A	134	GLU
1	A	145	LYS
1	A	156	ASN
1	A	174	THR
1	A	176	THR
2	B	7	ASN
2	B	13	LYS
2	B	25	SER
2	B	41	SER
1	C	7	SER
1	C	10	LYS
1	C	27	THR
1	C	30	LYS
1	C	45	LEU
1	C	48	THR
1	C	56	ASP
1	C	81	ASP
1	C	93	LYS
1	C	107	LYS
1	C	108	GLU
1	C	111	LYS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	C	113	SER
1	C	131	SER
1	C	132	ASN
1	C	133	LYS
1	C	150	LYS
1	C	151	SER
1	C	156	ASN
1	C	167	ASN
2	D	1	VAL
2	D	5	THR
2	D	13	LYS
2	D	41	SER
2	D	43	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	21	GLN
1	A	59	ASN
1	A	105	ASN
1	A	132	ASN
1	A	142	ASN
1	A	156	ASN
1	A	171	ASN
1	C	59	ASN
1	C	62	ASN
1	C	105	ASN
1	C	132	ASN
1	C	142	ASN
1	C	156	ASN
1	C	167	ASN
2	D	43	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 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
3	PO4	C	182	-	4,4,4	0.91	0	6,6,6	0.23	0
3	PO4	A	182	-	4,4,4	0.84	0	6,6,6	0.32	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.

No monomer is involved in short contacts.

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates (i)

EDS was not executed - this section is therefore empty.

6.4 Ligands (i)

EDS was not executed - this section is therefore empty.

6.5 Other polymers (i)

EDS was not executed - this section is therefore empty.