Full wwPDB X-ray Structure Validation Report (i)

Jun 13, 2024 – 07:25 AM EDT

PDB ID	:	1MCB
Title	:	PRINCIPLES AND PITFALLS IN DESIGNING SITE DIRECTED PEPTIDE LIGANDS
Authors	:	Edmundson, A.B.; Harris, D.L.; Fan, Z.-C.; Guddat, L.W.
Deposited on	:	1993-02-25
Resolution	:	2.70 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

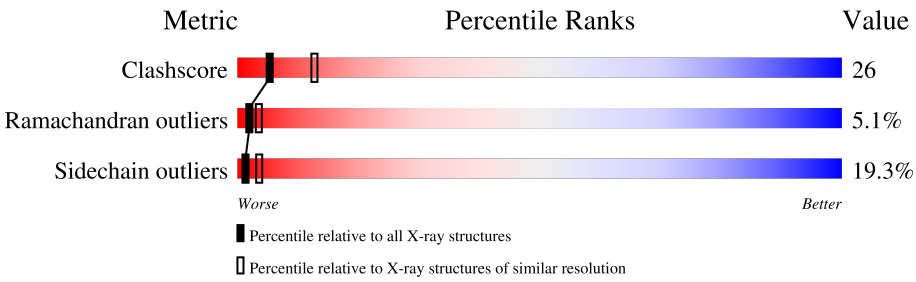
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (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 2.70 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	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)

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	216	49% 38% 11% •
1	B	216	40% 49% 10% •
2	P	5	80% 20%

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	DHI	P	3	X	-	-	-

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 3251 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 IMMUNOGLOBULIN LAMBDA DIMER MCG (LIGHT CHAIN).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	216	1605	1000	266	334	5	0	0	0
1	B	216	1605	1000	266	334	5	0	0	0

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	PRO	-	INSERTION	UNP Q6PIK1
A	14	LEU	PRO	CONFLICT	UNP Q6PIK1
A	42	ALA	PRO	CONFLICT	UNP Q6PIK1
A	48	VAL	LEU	CONFLICT	UNP Q6PIK1
A	49	ILE	MET	CONFLICT	UNP Q6PIK1
A	94	GLU	ALA	CONFLICT	UNP Q6PIK1
A	97	ASP	ASN	CONFLICT	UNP Q6PIK1
A	99	PHE	TYR	CONFLICT	UNP Q6PIK1
B	1	PRO	-	INSERTION	UNP Q6PIK1
B	14	LEU	PRO	CONFLICT	UNP Q6PIK1
B	42	ALA	PRO	CONFLICT	UNP Q6PIK1
B	48	VAL	LEU	CONFLICT	UNP Q6PIK1
B	49	ILE	MET	CONFLICT	UNP Q6PIK1
B	94	GLU	ALA	CONFLICT	UNP Q6PIK1
B	97	ASP	ASN	CONFLICT	UNP Q6PIK1
B	99	PHE	TYR	CONFLICT	UNP Q6PIK1

- Molecule 2 is a protein called PEPTIDE N-ACETYL-L-GLN-D-PHE-L-HIS-D-PRO-OH.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	P	5	41	27	7	7	0	0	0

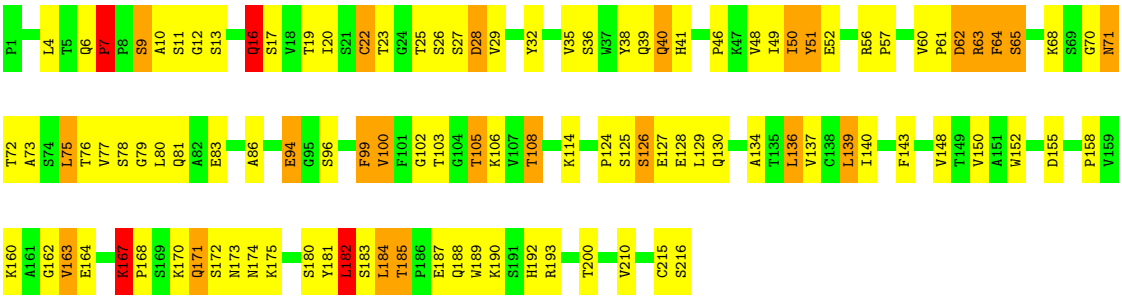
3 Residue-property plots (i)

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.

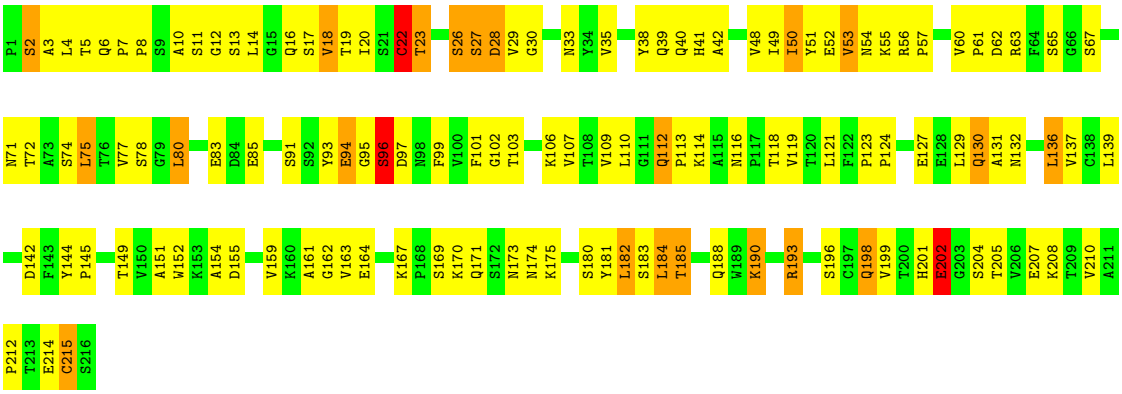
• Molecule 1: IMMUNOGLOBULIN LAMBDA DIMER MCG (LIGHT CHAIN)

Chain A: 49% 38% 11% •



• Molecule 1: IMMUNOGLOBULIN LAMBDA DIMER MCG (LIGHT CHAIN)

Chain B: 40% 49% 10% •



• Molecule 2: PEPTIDE N-ACETYL-L-GLN-D-PHE-L-HIS-D-PRO-OH

Chain P: 80% 20%

4 Data and refinement statistics (i)

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

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α, β, γ	72.30Å 72.30Å 185.90Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	6.00 – 2.70	Depositor
% Data completeness (in resolution range)	(Not available) (6.00-2.70)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	PROLSQ, X-PLOR	Depositor
R, R_{free}	0.204 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3251	wwPDB-VP
Average B, all atoms ($Å^2$)	0.0	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: DPR, ACE, DPN, DHI

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	0.88	0/1644	1.40	10/2241 (0.4%)
1	B	0.86	0/1644	1.39	8/2241 (0.4%)
2	P	0.76	0/9	1.37	0/11
All	All	0.87	0/3297	1.39	18/4493 (0.4%)

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	A	0	1
2	P	1	2
All	All	1	3

There are no bond length outliers.

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	136	LEU	CA-CB-CG	8.37	134.54	115.30
1	A	94	GLU	CA-CB-CG	7.48	129.85	113.40
1	B	33	ASN	N-CA-C	7.17	130.35	111.00
1	A	22	CYS	CA-CB-SG	6.77	126.19	114.00
1	B	22	CYS	CA-CB-SG	6.71	126.09	114.00
1	B	22	CYS	CB-CA-C	6.46	123.32	110.40
1	A	182	LEU	CA-CB-CG	6.14	129.43	115.30
1	B	193	ARG	NE-CZ-NH2	-5.90	117.35	120.30
1	A	136	LEU	CA-CB-CG	5.72	128.45	115.30
1	B	33	ASN	CA-C-N	5.40	129.09	117.20
1	A	7	PRO	N-CA-CB	-5.40	96.66	102.60

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	189	TRP	CA-CB-CG	5.32	123.81	113.70
1	B	129	LEU	CB-CA-C	5.28	120.24	110.20
1	A	128	GLU	CA-CB-CG	5.27	125.00	113.40
1	B	110	LEU	CA-CB-CG	5.25	127.39	115.30
1	A	152	TRP	CA-CB-CG	5.13	123.45	113.70
1	A	167	LYS	CA-CB-CG	5.13	124.69	113.40
1	A	100	VAL	CB-CA-C	5.04	120.97	111.40

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	P	3	DHI	CA

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	6	GLN	Peptide
2	P	3	DHI	Peptide,Mainchain

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	1605	0	1540	73	0
1	B	1605	0	1540	103	0
2	P	41	0	33	9	0
All	All	3251	0	3113	168	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 26.

All (168) 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:57:PRO:HA	1:B:97:ASP:HB2	1.30	1.09
1:A:61:PRO:HB2	1:A:64:PHE:HD2	1.44	0.82

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:106:LYS:N	1:B:106:LYS:HD2	1.98	0.78
1:A:64:PHE:CE1	1:A:77:VAL:HG22	2.19	0.77
1:B:116:ASN:HD22	1:B:201:HIS:HD2	1.35	0.75
1:A:162:GLY:O	1:A:182:LEU:HA	1.87	0.74
1:A:49:ILE:HD11	1:A:75:LEU:HD11	1.70	0.73
1:B:106:LYS:HD2	1:B:106:LYS:H	1.54	0.72
1:B:127:GLU:O	1:B:130:GLN:HB2	1.92	0.70
1:A:57:PRO:HA	1:B:97:ASP:CB	2.17	0.69
1:B:93:TYR:CD1	2:P:1:GLN:HB3	2.27	0.69
1:B:185:THR:HG23	1:B:188:GLN:HG3	1.76	0.67
1:A:167:LYS:HD3	1:B:169:SER:OG	1.94	0.67
1:A:57:PRO:CA	1:B:97:ASP:HB2	2.19	0.67
1:B:51:TYR:HD2	1:B:52:GLU:HG3	1.60	0.65
1:B:93:TYR:HA	1:B:99:PHE:HD1	1.60	0.65
1:A:20:ILE:HG21	1:A:105:THR:OG1	1.97	0.64
1:B:163:VAL:HG22	1:B:182:LEU:HG	1.79	0.64
1:B:26:SER:H	1:B:71:ASN:ND2	1.98	0.61
1:B:185:THR:HG23	1:B:188:GLN:CG	2.31	0.61
1:B:23:THR:HG23	1:B:72:THR:OG1	2.01	0.61
1:B:106:LYS:H	1:B:106:LYS:CD	2.13	0.60
1:B:155:ASP:OD1	1:B:193:ARG:HB2	2.00	0.60
1:B:51:TYR:CD2	1:B:52:GLU:HG3	2.36	0.60
1:B:136:LEU:HD11	1:B:210:VAL:HG21	1.84	0.59
1:B:10:ALA:O	1:B:107:VAL:HA	2.01	0.59
1:B:170:LYS:NZ	1:B:174:ASN:HB3	2.17	0.59
1:A:51:TYR:HD2	2:P:3:DHI:HE2	1.50	0.59
1:B:151:ALA:HB3	1:B:198:GLN:HE21	1.67	0.59
1:A:65:SER:O	1:A:75:LEU:HA	2.04	0.58
1:A:60:VAL:HG13	1:A:61:PRO:HD2	1.86	0.57
1:B:170:LYS:HZ2	1:B:174:ASN:HB3	1.69	0.57
1:A:49:ILE:CD1	1:A:75:LEU:HD11	2.34	0.57
2:P:2:DPN:HE1	2:P:4:DPR:OXT	2.05	0.57
1:B:106:LYS:N	1:B:106:LYS:CD	2.67	0.56
1:A:9:SER:HA	1:A:106:LYS:O	2.05	0.56
1:B:182:LEU:HD22	1:B:184:LEU:HD23	1.87	0.56
1:B:52:GLU:O	1:B:54:ASN:N	2.37	0.56
1:A:163:VAL:HA	1:A:181:TYR:O	2.06	0.56
1:A:38:TYR:CE2	1:A:48:VAL:HG22	2.40	0.56
1:B:154:ALA:HB2	1:B:159:VAL:HG21	1.87	0.55
1:B:93:TYR:CA	1:B:99:PHE:HD1	2.18	0.54
1:A:29:VAL:O	1:A:68:LYS:HD2	2.08	0.54

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:79:GLY:O	1:A:81:GLN:HG3	2.07	0.54
1:A:188:GLN:O	1:A:192:HIS:ND1	2.41	0.54
1:B:164:GLU:O	1:B:180:SER:HA	2.08	0.54
1:B:112:GLN:HB3	1:B:113:PRO:HD2	1.90	0.53
1:A:35:VAL:HG11	1:A:73:ALA:CB	2.37	0.53
1:B:119:VAL:O	1:B:208:LYS:HE2	2.08	0.53
1:A:143:PHE:HZ	1:A:168:PRO:HB3	1.73	0.53
1:B:96:SER:OG	1:B:97:ASP:N	2.41	0.53
1:B:50:ILE:HG13	1:B:56:ARG:HG2	1.91	0.53
1:B:136:LEU:CD1	1:B:210:VAL:HG21	2.39	0.53
1:A:12:GLY:HA3	1:A:80:LEU:CD1	2.39	0.52
1:A:19:THR:HG1	1:A:76:THR:HG1	1.56	0.52
1:B:163:VAL:HA	1:B:181:TYR:O	2.08	0.52
1:B:170:LYS:CG	1:B:174:ASN:HA	2.39	0.52
1:B:35:VAL:HA	1:B:91:SER:O	2.10	0.52
1:A:41:HIS:ND1	1:A:86:ALA:HB2	2.25	0.52
1:A:25:THR:HG22	1:A:26:SER:N	2.24	0.52
1:B:162:GLY:O	1:B:182:LEU:HA	2.10	0.52
1:A:125:SER:O	1:A:129:LEU:HG	2.10	0.51
1:A:215:CYS:O	1:A:216:SER:CB	2.58	0.51
1:B:85:GLU:HG3	1:B:109:VAL:HG23	1.91	0.51
1:A:25:THR:HG22	1:A:26:SER:H	1.76	0.51
1:B:16:GLN:CD	1:B:17:SER:H	2.14	0.51
1:B:124:PRO:HD3	1:B:136:LEU:HD22	1.92	0.51
1:A:155:ASP:OD1	1:A:193:ARG:HB2	2.09	0.51
1:B:51:TYR:O	1:B:55:LYS:HB2	2.11	0.51
1:A:63:ARG:HG2	1:A:78:SER:HB3	1.92	0.50
1:B:26:SER:O	1:B:27:SER:C	2.51	0.50
1:B:4:LEU:H	1:B:102:GLY:HA2	1.77	0.49
1:B:127:GLU:O	1:B:130:GLN:CB	2.60	0.49
1:B:145:PRO:O	1:B:201:HIS:HE1	1.95	0.49
1:A:27:SER:O	1:A:32:TYR:HE1	1.96	0.49
1:A:41:HIS:HE1	1:A:83:GLU:O	1.95	0.49
1:B:12:GLY:O	1:B:109:VAL:HA	2.13	0.49
1:B:152:TRP:CD1	1:B:163:VAL:HG11	2.48	0.49
1:B:95:GLY:O	1:B:96:SER:HB3	2.13	0.49
1:A:139:LEU:HG	1:B:137:VAL:HG21	1.95	0.48
1:B:51:TYR:HD1	1:B:57:PRO:HD3	1.78	0.48
1:B:170:LYS:HG2	1:B:174:ASN:HA	1.95	0.48
1:A:140:ILE:HD11	1:A:150:VAL:HG21	1.96	0.48
1:A:4:LEU:HB2	1:A:102:GLY:HA2	1.95	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:11:SER:OG	1:B:202:GLU:OE2	2.32	0.48
1:B:20:ILE:HD11	1:B:107:VAL:CG2	2.44	0.47
1:B:50:ILE:CG1	1:B:56:ARG:HG2	2.44	0.47
1:B:152:TRP:CE3	1:B:182:LEU:HD12	2.50	0.47
1:A:41:HIS:CE1	1:A:83:GLU:O	2.68	0.47
1:B:99:PHE:HB3	2:P:1:GLN:HE21	1.80	0.47
1:A:40:GLN:HB2	1:A:46:PRO:HA	1.97	0.46
1:A:164:GLU:HG2	1:B:171:GLN:HG2	1.96	0.46
1:B:14:LEU:CD1	1:B:109:VAL:HG11	2.45	0.46
1:B:93:TYR:O	1:B:94:GLU:O	2.32	0.46
1:A:50:ILE:HD13	1:A:75:LEU:HD21	1.98	0.46
1:B:99:PHE:CG	2:P:1:GLN:HG2	2.50	0.46
1:A:124:PRO:HG3	1:A:134:ALA:HB1	1.97	0.46
1:B:57:PRO:O	1:B:60:VAL:HG23	2.15	0.46
1:A:185:THR:OG1	1:A:188:GLN:OE1	2.33	0.46
1:B:185:THR:O	1:B:188:GLN:HB2	2.16	0.46
1:A:99:PHE:CE2	2:P:2:DPN:O	2.69	0.46
1:B:121:LEU:CD2	1:B:210:VAL:HG13	2.46	0.45
1:B:61:PRO:O	1:B:63:ARG:N	2.49	0.45
1:B:149:THR:O	1:B:199:VAL:HA	2.17	0.45
1:B:99:PHE:CB	2:P:1:GLN:HG2	2.46	0.45
1:B:6:GLN:HB3	1:B:7:PRO:HD2	1.99	0.45
1:A:23:THR:HA	1:A:71:ASN:O	2.17	0.45
1:A:40:GLN:HE22	1:B:40:GLN:HE21	1.65	0.45
1:A:80:LEU:HD23	1:A:80:LEU:HA	1.83	0.45
1:B:190:LYS:NZ	1:B:212:PRO:HB2	2.32	0.45
1:B:20:ILE:HD11	1:B:107:VAL:HG21	1.99	0.44
1:A:4:LEU:HG	1:A:100:VAL:HG13	1.98	0.44
1:A:171:GLN:HG3	1:A:175:LYS:O	2.17	0.44
1:A:81:GLN:HB3	1:A:83:GLU:CD	2.37	0.44
1:A:173:ASN:O	1:A:175:LYS:N	2.51	0.43
1:B:14:LEU:HD12	1:B:109:VAL:HG11	2.00	0.43
1:B:173:ASN:C	1:B:175:LYS:H	2.22	0.43
1:A:56:ARG:HE	1:A:56:ARG:HB2	1.58	0.43
1:A:61:PRO:HB2	1:A:64:PHE:CD2	2.36	0.43
1:A:181:TYR:CD2	1:B:139:LEU:HD11	2.53	0.43
1:B:99:PHE:HE2	1:B:101:PHE:CE1	2.36	0.43
1:B:131:ALA:O	1:B:132:ASN:HB3	2.19	0.43
1:A:16:GLN:CG	1:A:17:SER:H	2.30	0.43
1:A:140:ILE:HG22	1:A:143:PHE:CE2	2.54	0.43
1:B:112:GLN:CB	1:B:113:PRO:HD2	2.49	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:P:0:ACE:O	2:P:2:DPN:N	2.52	0.43
1:A:56:ARG:HD3	1:A:64:PHE:O	2.18	0.43
1:B:5:THR:O	1:B:22:CYS:HA	2.19	0.43
1:B:99:PHE:CD1	2:P:1:GLN:HA	2.53	0.43
1:B:26:SER:H	1:B:71:ASN:HD21	1.64	0.42
1:A:126:SER:O	1:A:129:LEU:HB2	2.19	0.42
1:B:39:GLN:OE1	1:B:41:HIS:NE2	2.52	0.42
1:B:154:ALA:CB	1:B:159:VAL:HG21	2.48	0.42
1:A:127:GLU:O	1:A:130:GLN:N	2.52	0.42
1:B:57:PRO:HG2	1:B:60:VAL:HG21	2.02	0.42
1:A:140:ILE:HG22	1:A:143:PHE:CD2	2.55	0.42
1:B:2:SER:OG	1:B:3:ALA:N	2.52	0.42
1:A:11:SER:HB3	1:A:108:THR:OG1	2.19	0.42
1:A:140:ILE:HD11	1:A:150:VAL:CG2	2.49	0.42
1:B:49:ILE:O	1:B:57:PRO:HD2	2.20	0.42
1:B:142:ASP:HA	1:B:175:LYS:HB3	2.00	0.42
1:A:160:LYS:HA	1:A:160:LYS:HD2	1.78	0.42
1:A:167:LYS:HD3	1:B:169:SER:HG	1.83	0.42
1:B:11:SER:OG	1:B:114:LYS:HE2	2.19	0.42
1:B:116:ASN:ND2	1:B:201:HIS:HD2	2.11	0.42
1:A:70:GLY:C	1:A:72:THR:H	2.22	0.41
1:A:137:VAL:HG11	1:B:137:VAL:HG11	2.01	0.41
1:A:11:SER:HA	1:A:108:THR:O	2.20	0.41
1:A:28:ASP:O	1:A:32:TYR:HD1	2.03	0.41
1:A:150:VAL:HG11	1:A:180:SER:CB	2.50	0.41
1:A:65:SER:O	1:A:75:LEU:HD23	2.21	0.41
1:B:190:LYS:HA	1:B:190:LYS:HD2	1.65	0.41
1:B:99:PHE:CE2	1:B:101:PHE:CE1	3.08	0.41
1:B:144:TYR:HA	1:B:145:PRO:C	2.41	0.41
1:A:215:CYS:O	1:A:215:CYS:SG	2.79	0.41
1:B:17:SER:HA	1:B:77:VAL:O	2.21	0.41
1:B:196:SER:HA	1:B:208:LYS:O	2.21	0.41
1:B:18:VAL:HG13	1:B:80:LEU:HD13	2.03	0.41
1:B:65:SER:O	1:B:75:LEU:HA	2.21	0.41
1:B:123:PRO:HA	1:B:136:LEU:HD13	2.03	0.41
1:B:67:SER:OG	1:B:74:SER:HB2	2.21	0.41
1:B:199:VAL:O	1:B:205:THR:HA	2.21	0.41
1:B:215:CYS:O	1:B:215:CYS:SG	2.78	0.41
1:A:61:PRO:O	1:A:63:ARG:N	2.54	0.40
1:A:184:LEU:HD12	1:A:184:LEU:N	2.36	0.40
1:A:124:PRO:HD3	1:A:136:LEU:HD23	2.03	0.40

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:193:ARG:HH11	1:A:193:ARG:HG3	1.86	0.40
1:B:80:LEU:HD21	1:B:109:VAL:HG22	2.04	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	214/216 (99%)	177 (83%)	27 (13%)	10 (5%)	2 4
1	B	214/216 (99%)	177 (83%)	25 (12%)	12 (6%)	2 3
2	P	1/5 (20%)	0	1 (100%)	0	100 100
All	All	429/437 (98%)	354 (82%)	53 (12%)	22 (5%)	2 3

All (22) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	7	PRO
1	A	10	ALA
1	A	96	SER
1	A	174	ASN
1	B	2	SER
1	B	53	VAL
1	B	62	ASP
1	B	94	GLU
1	B	161	ALA
1	A	62	ASP
1	B	28	ASP
1	B	29	VAL
1	B	42	ALA
1	A	16	GLN
1	B	8	PRO

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	96	SER
1	A	172	SER
1	B	202	GLU
1	A	39	GLN
1	B	30	GLY
1	A	40	GLN
1	A	71	ASN

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	181/181 (100%)	144 (80%)	37 (20%)	1 3
1	B	181/181 (100%)	148 (82%)	33 (18%)	1 4
2	P	1/1 (100%)	1 (100%)	0	100 100
All	All	363/363 (100%)	293 (81%)	70 (19%)	1 3

All (70) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	PRO
1	A	9	SER
1	A	13	SER
1	A	16	GLN
1	A	22	CYS
1	A	28	ASP
1	A	36	SER
1	A	50	ILE
1	A	51	TYR
1	A	52	GLU
1	A	62	ASP
1	A	63	ARG
1	A	64	PHE
1	A	65	SER
1	A	75	LEU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	94	GLU
1	A	99	PHE
1	A	103	THR
1	A	105	THR
1	A	108	THR
1	A	114	LYS
1	A	126	SER
1	A	139	LEU
1	A	148	VAL
1	A	158	PRO
1	A	163	VAL
1	A	167	LYS
1	A	170	LYS
1	A	171	GLN
1	A	182	LEU
1	A	183	SER
1	A	184	LEU
1	A	185	THR
1	A	187	GLU
1	A	190	LYS
1	A	200	THR
1	A	210	VAL
1	B	13	SER
1	B	18	VAL
1	B	19	THR
1	B	22	CYS
1	B	23	THR
1	B	26	SER
1	B	27	SER
1	B	28	ASP
1	B	38	TYR
1	B	48	VAL
1	B	50	ILE
1	B	53	VAL
1	B	75	LEU
1	B	78	SER
1	B	80	LEU
1	B	83	GLU
1	B	96	SER
1	B	103	THR
1	B	112	GLN
1	B	118	THR

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	130	GLN
1	B	167	LYS
1	B	182	LEU
1	B	183	SER
1	B	184	LEU
1	B	185	THR
1	B	190	LYS
1	B	198	GLN
1	B	202	GLU
1	B	204	SER
1	B	207	GLU
1	B	214	GLU
1	B	215	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	40	GLN
1	A	116	ASN
1	A	201	HIS
1	B	40	GLN
1	B	71	ASN
1	B	132	ASN
1	B	198	GLN
1	B	201	HIS

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

3 non-standard protein/DNA/RNA residues are modelled in this entry.

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	DHI	P	3	2	1/1/1/2	-	-

There are no bond length outliers.

There are no bond angle outliers.

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	P	3	DHI	CA

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	P	3	DHI	1	0

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

There are no ligands in this entry.

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.