



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 09:53 AM GMT

PDB ID : 1MCB
Title : PRINCIPLES AND PITFALLS IN DESIGNING SITE DIRECTED PEP-
TIDE 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 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>

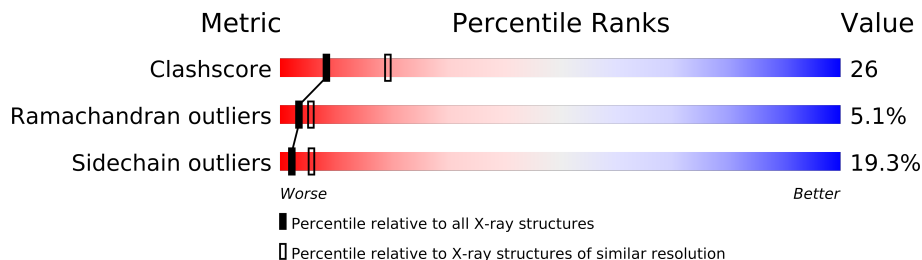
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 21963
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 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	79885	1939 (2.70-2.70)
Ramachandran outliers	78287	1905 (2.70-2.70)
Sidechain outliers	78261	1905 (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 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	216	
1	B	216	
2	P	5	

2 Entry composition

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

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

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.

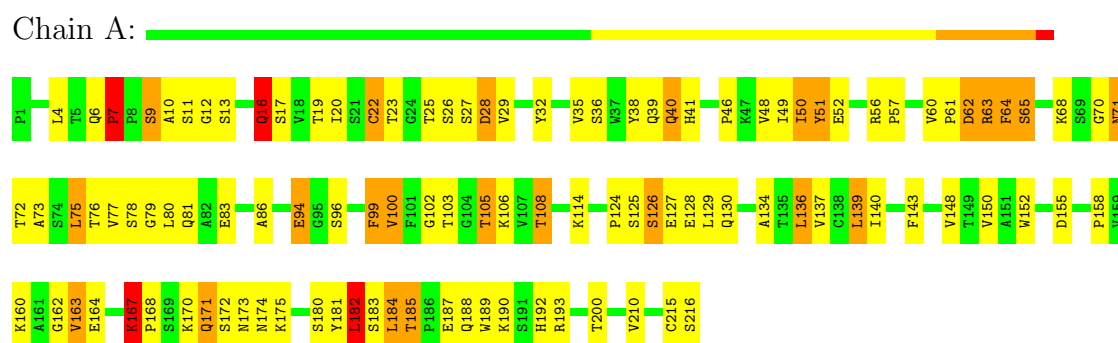
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	P	5	Total	C	N	O	0	0	0
			41	27	7	7			

3 Residue-property plots

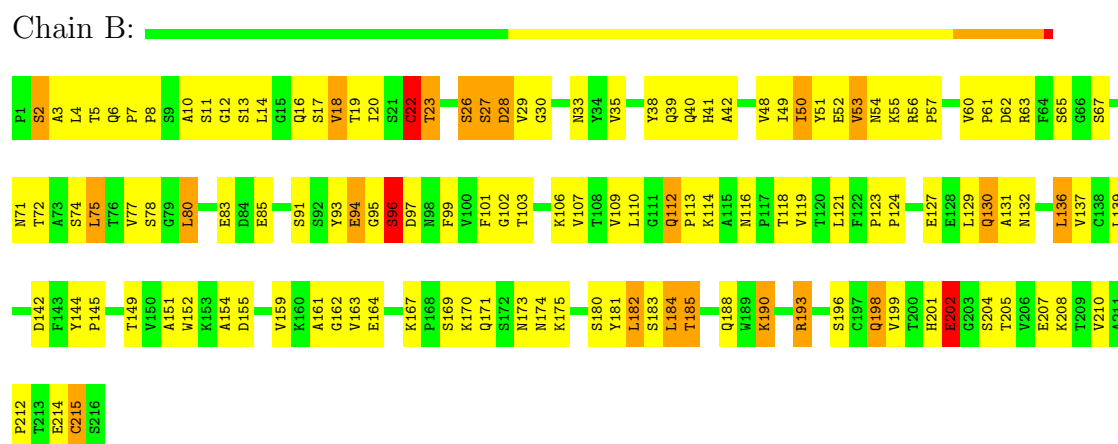
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according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

Note EDS was not executed.

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



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



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



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: DHI, DPN, DPR, ACE

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(°)	Ideal(°)
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

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
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	Mainchain,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	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

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 26.

All (168) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(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

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Atom-1	Atom-2	Distance(Å)	Clash(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:CD	1:B:106:LYS:H	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:B:182:LEU:HD22	1:B:184:LEU:HD23	1.87	0.56
1:A:9:SER:HA	1:A:106:LYS:O	2.05	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
1:A:79:GLY:O	1:A:81:GLN:HG3	2.07	0.54

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:164:GLU:O	1:B:180:SER:HA	2.08	0.54
1:A:188:GLN:O	1:A:192:HIS:ND1	2.41	0.54
1:B:112:GLN:HB3	1:B:113:PRO:HD2	1.90	0.53
1:B:119:VAL:O	1:B:208:LYS:HE2	2.08	0.53
1:A:35:VAL:HG11	1:A:73:ALA:CB	2.37	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:136:LEU:CD1	1:B:210:VAL:HG21	2.39	0.53
1:B:50:ILE:HG13	1:B:56:ARG:HG2	1.91	0.53
1:B:163:VAL:HA	1:B:181:TYR:O	2.08	0.52
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: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:B:162:GLY:O	1:B:182:LEU:HA	2.10	0.52
1:A:25:THR:HG22	1:A:26:SER:N	2.24	0.52
1:A:125:SER:O	1:A:129:LEU:HG	2.10	0.51
1:B:85:GLU:HG3	1:B:109:VAL:HG23	1.91	0.51
1:A:215:CYS:O	1:A:216:SER:CB	2.58	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:B:51:TYR:O	1:B:55:LYS:HB2	2.11	0.51
1:A:155:ASP:OD1	1:A:193:ARG:HB2	2.09	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:41:HIS:HE1	1:A:83:GLU:O	1.95	0.49
1:A:27:SER:O	1:A:32:TYR:HE1	1.96	0.49
1:B:152:TRP:CD1	1:B:163:VAL:HG11	2.48	0.49
1:B:12:GLY:O	1:B:109:VAL:HA	2.13	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
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

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
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:B:99:PHE:HB3	2:P:1:GLN:HE21	1.80	0.47
1:A:41:HIS:CE1	1:A:83:GLU:O	2.68	0.47
1:A:164:GLU:HG2	1:B:171:GLN:HG2	1.96	0.46
1:A:40:GLN:HB2	1:A:46:PRO:HA	1.97	0.46
1:B:93:TYR:O	1:B:94:GLU:O	2.32	0.46
1:B:14:LEU:CD1	1:B:109:VAL:HG11	2.45	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:B:57:PRO:O	1:B:60:VAL:HG23	2.15	0.46
1:A:124:PRO:HG3	1:A:134:ALA:HB1	1.97	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:80:LEU:HA	1:A:80:LEU:HD23	1.83	0.45
1:A:40:GLN:HE22	1:B:40:GLN:HE21	1.65	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:B:99:PHE:HE2	1:B:101:PHE:CE1	2.36	0.43
1:A:181:TYR:CD2	1:B:139:LEU:HD11	2.53	0.43
1:B:131:ALA:O	1:B:132:ASN:HB3	2.19	0.43
2:P:0:ACE:O	2:P:2:DPN:N	2.52	0.43
1:B:112:GLN:CB	1:B:113:PRO:HD2	2.49	0.43
1:A:140:ILE:HG22	1:A:143:PHE:CE2	2.54	0.43
1:A:16:GLN:CG	1:A:17:SER:H	2.30	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

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
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:B:154:ALA:CB	1:B:159:VAL:HG21	2.48	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:57:PRO:HG2	1:B:60:VAL:HG21	2.02	0.42
1:A:127:GLU:O	1:A:130:GLN:N	2.52	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:B:49:ILE:O	1:B:57:PRO:HD2	2.20	0.42
1:A:140:ILE:HD11	1:A:150:VAL:CG2	2.49	0.42
1:B:142:ASP:HA	1:B:175:LYS:HB3	2.00	0.42
1:A:11:SER:HB3	1:A:108:THR:OG1	2.19	0.42
1:B:116:ASN:ND2	1:B:201:HIS:HD2	2.11	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:A:160:LYS:HA	1:A:160:LYS:HD2	1.78	0.42
1:A:137:VAL:HG11	1:B:137:VAL:HG11	2.01	0.41
1:A:70:GLY:C	1:A:72:THR:H	2.22	0.41
1:A:28:ASP:O	1:A:32:TYR:HD1	2.03	0.41
1:A:11:SER:HA	1:A:108:THR:O	2.20	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:144:TYR:HA	1:B:145:PRO:C	2.41	0.41
1:B:99:PHE:CE2	1:B:101:PHE:CE1	3.08	0.41
1:A:215:CYS:SG	1:A:215:CYS:O	2.79	0.41
1:B:196:SER:HA	1:B:208:LYS:O	2.21	0.41
1:B:17:SER:HA	1:B:77:VAL:O	2.21	0.41
1:B:123:PRO:HA	1:B:136:LEU:HD13	2.03	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:199:VAL:O	1:B:205:THR:HA	2.21	0.41
1:B:67:SER:OG	1:B:74:SER:HB2	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:N	1:A:184:LEU:HD12	2.36	0.40
1:A:193:ARG:HG3	1:A:193:ARG:HH11	1.86	0.40
1:A:124:PRO:HD3	1:A:136:LEU:HD23	2.03	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

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	214/216 (99%)	177 (83%)	27 (13%)	10 (5%)	4	7
1	B	214/216 (99%)	177 (83%)	25 (12%)	12 (6%)	3	4
2	P	1/5 (20%)	0	1 (100%)	0	100	100
All	All	429/437 (98%)	354 (82%)	53 (12%)	22 (5%)	3	5

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
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 ⓘ

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%)	2	5
1	B	181/181 (100%)	148 (82%)	33 (18%)	2	6
2	P	1/1 (100%)	1 (100%)	0	100	100
All	All	363/363 (100%)	293 (81%)	70 (19%)	2	5

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
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

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Mol	Chain	Res	Type
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
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

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Mol	Chain	Res	Type
1	B	215	CYS

Some 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 ⓘ

There are no RNA chains in this entry.

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

3 non-standard protein/DNA/RNA residues 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	DPN	P	2	2	11,11,12	5.22	2 (18%)	11,13,15	2.89	1 (9%)
2	DHI	P	3	2	10,10,11	5.90	1 (10%)	10,12,14	2.44	1 (10%)
2	DPR	P	4	2	8,8,8	2.08	3 (37%)	10,10,10	1.70	3 (30%)

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	DPN	P	2	2	-	0/4/6/8	0/1/1/1
2	DHI	P	3	2	1/1/1/2	0/4/6/8	0/1/1/1
2	DPR	P	4	2	-	0/4/11/11	0/1/1/1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	P	3	DHI	O-C	18.45	1.24	1.11
2	P	2	DPN	O-C	16.92	1.23	1.11
2	P	4	DPR	OXT-C	-3.83	1.16	1.30
2	P	4	DPR	CB-CA	-2.87	1.42	1.54
2	P	2	DPN	CA-C	2.84	1.53	1.48
2	P	4	DPR	CA-C	2.64	1.59	1.52

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	P	2	DPN	C-CA-N	-9.35	104.49	113.83
2	P	3	DHI	C-CA-N	-7.45	106.38	113.83
2	P	4	DPR	OXT-C-CA	3.28	126.21	113.90
2	P	4	DPR	C-CA-N	-2.62	95.52	106.47
2	P	4	DPR	O-C-CA	-2.44	112.95	121.86

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.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

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 ⓘ

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.