



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:43 AM UTC

PDB ID : 1MCB / pdb_00001mcb
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 X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

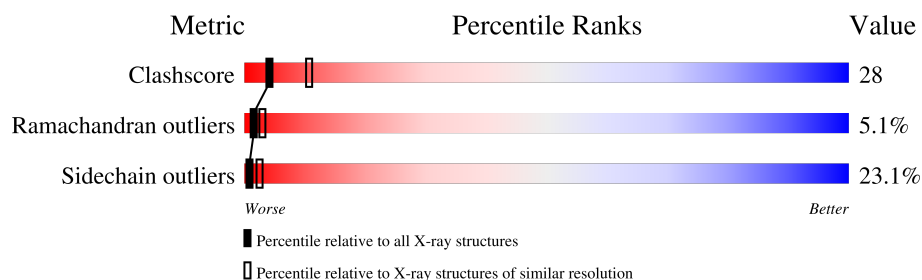
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (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 $\leq 5\%$.

Note EDS was not executed.

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

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

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
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.

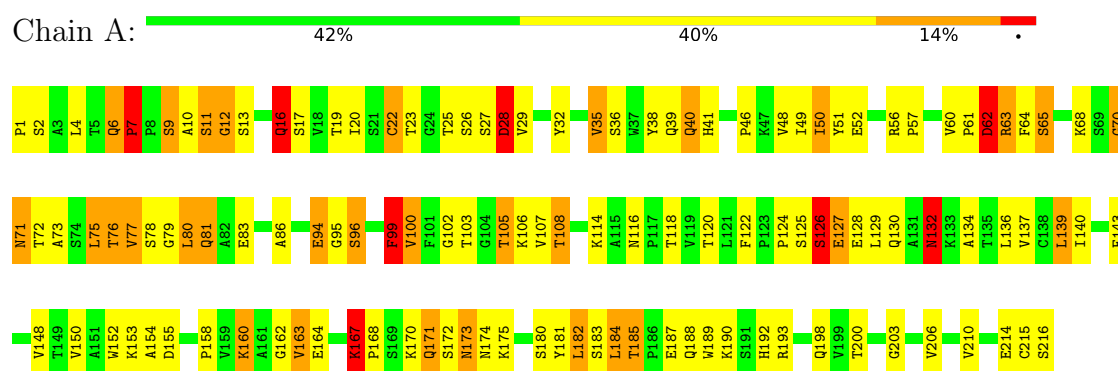
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	P	5	Total	C	N	O	0	0	0
			41	27	7	7			

3 Residue-property plots

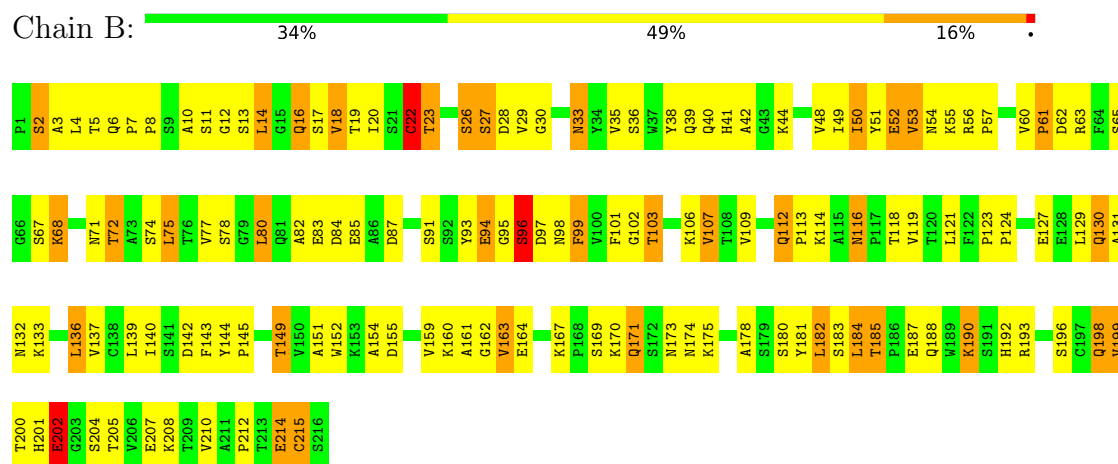
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded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

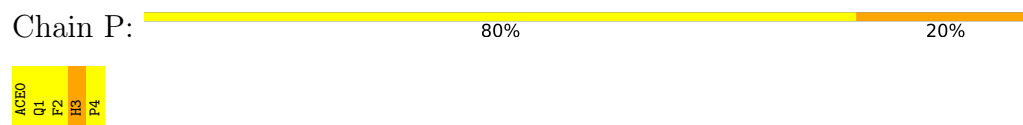
• Molecule 1: IMMUNOGLOBULIN LAMBDA DIMER MCG (LIGHT CHAIN)



• 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 is therefore incomplete.

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	72.30 Å 72.30 Å 185.90 Å 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: DPN, DPR, ACE, 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	1.13	2/1644 (0.1%)	1.82	32/2241 (1.4%)
1	B	1.09	1/1644 (0.1%)	1.80	28/2241 (1.2%)
2	P	0.65	0/9	2.09	0/11
All	All	1.11	3/3297 (0.1%)	1.81	60/4493 (1.3%)

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	A	0	1
2	P	1	2
All	All	1	3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	12	GLY	N-CA	7.69	1.53	1.44
1	A	6	GLN	CA-C	5.83	1.57	1.52
1	B	33	ASN	C-O	5.25	1.30	1.24

All (60) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	33	ASN	N-CA-C	14.44	130.35	113.02
1	A	99	PHE	CA-CB-CG	9.18	122.98	113.80
1	B	22	CYS	CB-CA-C	9.02	123.32	110.79
1	B	87	ASP	CA-CB-CG	8.27	120.87	112.60
1	B	99	PHE	CA-CB-CG	7.88	121.68	113.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	94	GLU	CA-CB-CG	7.88	129.85	114.10
1	A	122	PHE	CA-CB-CG	7.63	121.43	113.80
1	A	7	PRO	N-CA-C	7.10	119.36	110.70
1	B	143	PHE	CA-CB-CG	7.07	120.87	113.80
1	B	33	ASN	CA-C-O	-7.02	110.94	119.43
1	B	202	GLU	CB-CG-CD	7.01	124.51	112.60
1	B	192	HIS	CA-CB-CG	-7.00	106.80	113.80
1	A	100	VAL	CB-CA-C	6.85	120.97	111.38
1	A	7	PRO	N-CA-CB	-6.62	96.66	103.08
1	A	173	ASN	CA-CB-CG	6.61	119.21	112.60
1	A	28	ASP	CA-CB-CG	6.51	119.11	112.60
1	A	116	ASN	CA-CB-CG	6.43	119.03	112.60
1	A	62	ASP	CA-CB-CG	6.41	119.01	112.60
1	B	52	GLU	CB-CG-CD	6.32	123.33	112.60
1	A	126	SER	CA-C-O	6.28	127.59	121.00
1	A	16	GLN	CB-CG-CD	6.21	123.15	112.60
1	A	2	SER	N-CA-C	6.16	118.28	110.33
1	A	107	VAL	O-C-N	6.14	129.40	123.14
1	A	1	PRO	CA-C-N	5.99	131.61	120.95
1	A	1	PRO	C-N-CA	5.99	131.61	120.95
1	B	163	VAL	N-CA-C	5.83	116.87	108.48
1	A	35	VAL	O-C-N	5.79	129.20	123.18
1	A	80	LEU	CA-C-N	5.78	129.69	121.42
1	A	80	LEU	C-N-CA	5.78	129.69	121.42
1	B	116	ASN	CA-CB-CG	5.75	118.35	112.60
1	B	61	PRO	CA-C-N	5.64	132.32	121.54
1	B	61	PRO	C-N-CA	5.64	132.32	121.54
1	A	77	VAL	O-C-N	5.63	129.81	122.94
1	A	11	SER	CA-C-N	-5.61	115.59	122.15
1	A	11	SER	C-N-CA	-5.61	115.59	122.15
1	A	132	ASN	CA-CB-CG	-5.55	107.05	112.60
1	A	203	GLY	N-CA-C	-5.49	108.16	114.69
1	A	128	GLU	CA-CB-CG	5.45	125.00	114.10
1	A	206	VAL	N-CA-C	-5.42	100.68	108.48
1	B	178	ALA	CA-C-O	-5.37	115.48	121.23
1	A	189	TRP	CA-CB-CG	5.37	123.81	113.60
1	B	214	GLU	CB-CG-CD	5.37	121.72	112.60
1	B	68	LYS	N-CA-C	5.36	116.18	107.23
1	B	103	THR	N-CA-C	5.33	118.94	112.23
1	B	187	GLU	N-CA-CB	5.32	118.03	110.16
1	A	167	LYS	CA-CB-CG	5.30	124.69	114.10
1	B	140	ILE	O-C-N	5.28	128.80	123.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	70	GLY	N-CA-C	-5.26	100.70	113.18
1	B	16	GLN	CB-CG-CD	5.21	121.46	112.60
1	B	136	LEU	CA-CB-CG	5.21	134.54	116.30
1	B	129	LEU	CB-CA-C	5.19	120.24	110.63
1	A	152	TRP	CA-CB-CG	5.19	123.45	113.60
1	B	215	CYS	N-CA-C	5.17	117.66	111.71
1	B	133	LYS	CA-C-O	-5.17	115.06	121.06
1	A	22	CYS	CA-CB-SG	5.12	126.19	114.40
1	B	199	VAL	N-CA-C	5.11	115.32	108.17
1	B	22	CYS	CA-CB-SG	5.08	126.09	114.40
1	A	132	ASN	N-CA-C	5.08	118.77	112.58
1	B	171	GLN	CB-CG-CD	5.06	121.20	112.60
1	B	98	ASN	CA-CB-CG	-5.02	107.58	112.60

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

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	79	0
1	B	1605	0	1540	104	0
2	P	41	0	33	9	0
All	All	3251	0	3113	175	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 28.

All (175) 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
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:16:GLN:CD	1:B:17:SER:H	2.14	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:154:ALA:HB2	1:B:159:VAL:HG21	1.87	0.55
1:B:26:SER:O	1:B:27:SER:C	2.51	0.54
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
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:33:ASN:OD1	1:B:68:LYS:NZ	2.43	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: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: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:A:81:GLN:HB3	1:A:83:GLU:CD	2.37	0.48
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:173:ASN:C	1:B:175:LYS:H	2.22	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
1:B:50:ILE:CG1	1:B:56:ARG:HG2	2.44	0.47
1:B:95:GLY:O	1:B:96:SER:HB3	2.13	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:A:70:GLY:C	1:A:72:THR:H	2.22	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:B:144:TYR:HA	1:B:145:PRO:C	2.41	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:A:171:GLN:C	1:A:173:ASN:H	2.25	0.44
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: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: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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:A:95:GLY:O	1:A:96:SER:C	2.61	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
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:A:167:LYS:HD3	1:B:169:SER:HG	1.84	0.42
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:A:214:GLU:O	1:A:215:CYS:C	2.62	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: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: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:154:ALA:O	1:A:155:ASP:C	2.63	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: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:A:130:GLN:C	1:A:132:ASN:H	2.28	0.41
1:B:18:VAL:HG13	1:B:80:LEU:HD13	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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: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:B:82:ALA:C	1:B:84:ASP:H	2.29	0.40
1:A:124:PRO:HD3	1:A:136:LEU:HD23	2.03	0.40
1:A:173:ASN:C	1:A:175:LYS:H	2.28	0.40
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	3
1	B	214/216 (99%)	177 (83%)	25 (12%)	12 (6%)	1	2
2	P	1/5 (20%)	0	1 (100%)	0	100	100
All	All	429/437 (98%)	354 (82%)	53 (12%)	22 (5%)	1	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

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Mol	Chain	Res	Type
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%)	137 (76%)	44 (24%)	1	2
1	B	181/181 (100%)	141 (78%)	40 (22%)	1	3
2	P	1/1 (100%)	1 (100%)	0	100	100
All	All	363/363 (100%)	279 (77%)	84 (23%)	1	2

All (84) 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

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Mol	Chain	Res	Type
1	A	36	SER
1	A	50	ILE
1	A	52	GLU
1	A	62	ASP
1	A	63	ARG
1	A	65	SER
1	A	75	LEU
1	A	76	THR
1	A	81	GLN
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	118	THR
1	A	120	THR
1	A	126	SER
1	A	127	GLU
1	A	132	ASN
1	A	139	LEU
1	A	148	VAL
1	A	153	LYS
1	A	158	PRO
1	A	160	LYS
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	198	GLN
1	A	200	THR
1	A	210	VAL
1	B	13	SER
1	B	14	LEU
1	B	18	VAL
1	B	19	THR

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Mol	Chain	Res	Type
1	B	22	CYS
1	B	23	THR
1	B	26	SER
1	B	27	SER
1	B	36	SER
1	B	38	TYR
1	B	44	LYS
1	B	48	VAL
1	B	50	ILE
1	B	53	VAL
1	B	72	THR
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	107	VAL
1	B	112	GLN
1	B	118	THR
1	B	130	GLN
1	B	149	THR
1	B	160	LYS
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	200	THR
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 (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	201	HIS
1	B	40	GLN

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

5.3.3 RNA ⓘ

There are no RNA molecules 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 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 ⓘ

There are no oligosaccharides 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 is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.