



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 15, 2020 – 10:17 pm BST

PDB ID : 1MCQ
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 ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtrriage (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) : 2.11

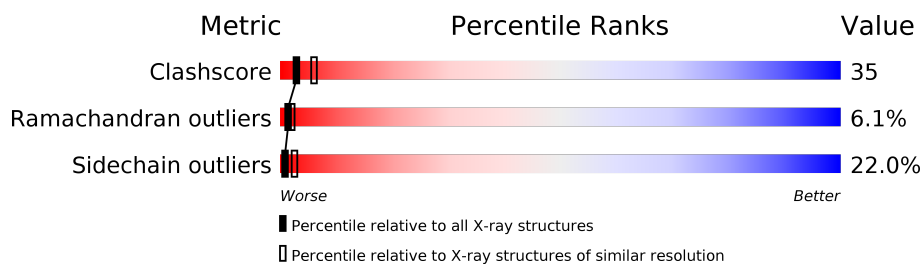
1 Overall quality at a glance

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	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 on 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	4	

2 Entry composition i

There are 2 unique types of molecules in this entry. The entry contains 3231 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 44 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	20	ILE	PHE	CONFLICT	PIR S14675
A	23	THR	SER	CONFLICT	PIR S14675
A	29	VAL	ILE	CONFLICT	PIR S14675
A	31	GLY	ASN	CONFLICT	PIR S14675
A	39	GLN	ARG	CONFLICT	PIR S14675
A	42	ALA	PRO	CONFLICT	PIR S14675
A	48	VAL	LEU	CONFLICT	PIR S14675
A	49	ILE	MET	CONFLICT	PIR S14675
A	54	ASN	THR	CONFLICT	PIR S14675
A	62	ASP	ASN	CONFLICT	PIR S14675
A	94	GLU	ALA	CONFLICT	PIR S14675
A	97	ASP	ASN	CONFLICT	PIR S14675
A	98	ASN	SER	CONFLICT	PIR S14675
A	99	PHE	LEU	CONFLICT	PIR S14675
A	100	VAL	ILE	CONFLICT	PIR S14675
A	103	THR	GLY	CONFLICT	PIR S14675
A	106	LYS	ARG	CONFLICT	PIR S14675
A	107	VAL	LEU	CONFLICT	PIR S14675
A	116	ASN	ALA	CONFLICT	PIR S14675
A	118	THR	SER	CONFLICT	PIR S14675
A	156	GLY	SER	CONFLICT	PIR S14675
A	167	LYS	THR	CONFLICT	PIR S14675
B	20	ILE	PHE	CONFLICT	PIR S14675
B	23	THR	SER	CONFLICT	PIR S14675

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Chain	Residue	Modelled	Actual	Comment	Reference
B	29	VAL	ILE	CONFLICT	PIR S14675
B	31	GLY	ASN	CONFLICT	PIR S14675
B	39	GLN	ARG	CONFLICT	PIR S14675
B	42	ALA	PRO	CONFLICT	PIR S14675
B	48	VAL	LEU	CONFLICT	PIR S14675
B	49	ILE	MET	CONFLICT	PIR S14675
B	54	ASN	THR	CONFLICT	PIR S14675
B	62	ASP	ASN	CONFLICT	PIR S14675
B	94	GLU	ALA	CONFLICT	PIR S14675
B	97	ASP	ASN	CONFLICT	PIR S14675
B	98	ASN	SER	CONFLICT	PIR S14675
B	99	PHE	LEU	CONFLICT	PIR S14675
B	100	VAL	ILE	CONFLICT	PIR S14675
B	103	THR	GLY	CONFLICT	PIR S14675
B	106	LYS	ARG	CONFLICT	PIR S14675
B	107	VAL	LEU	CONFLICT	PIR S14675
B	116	ASN	ALA	CONFLICT	PIR S14675
B	118	THR	SER	CONFLICT	PIR S14675
B	156	GLY	SER	CONFLICT	PIR S14675
B	167	LYS	THR	CONFLICT	PIR S14675

- Molecule 2 is a protein called PEPTIDE N-ACETYL-L-HIS-D-PRO-NH₂.

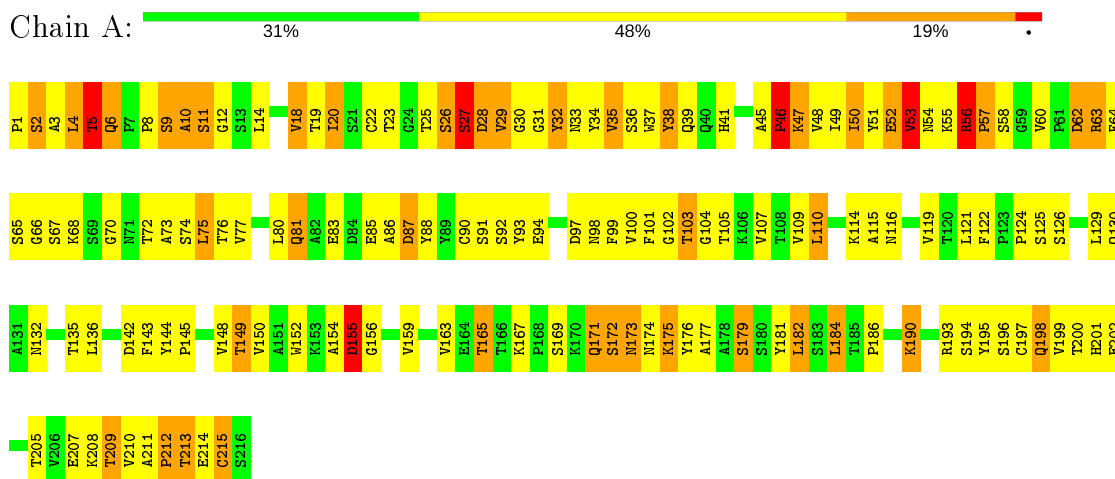
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	P	4	21	13	5	3	0	0	1

3 Residue-property plots

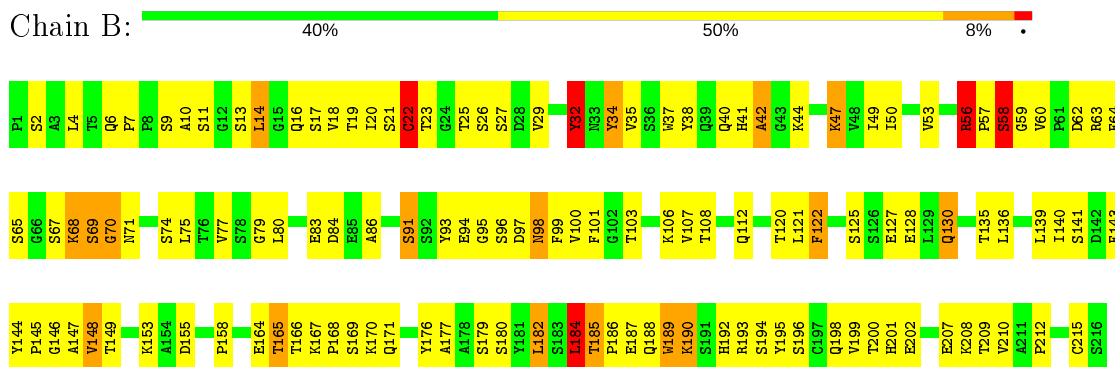
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded according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

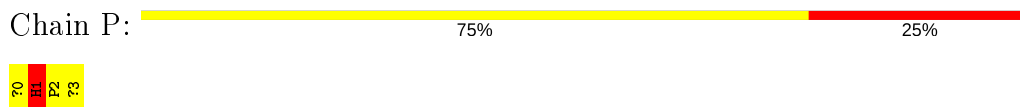
- Molecule 1: IMMUNOGLOBULIN LAMBDA DIMER MCG (LIGHT CHAIN)



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



- Molecule 2: PEPTIDE N-ACETYL-L-HIS-D-PRO-NH2



4 Data and refinement statistics

Xtrriage (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.	Xtrriage
Total number of atoms	3231	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, NH2

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	0.85	0/1644	1.41	7/2241 (0.3%)
1	B	0.87	0/1644	1.38	10/2241 (0.4%)
2	P	0.76	0/11	2.21	1/14 (7.1%)
All	All	0.86	0/3299	1.40	18/4496 (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	B	0	1

There are no bond length outliers.

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	52	GLU	CA-CB-CG	9.27	133.80	113.40
1	B	202	GLU	CA-CB-CG	9.00	133.20	113.40
1	B	32	TYR	CA-CB-CG	-7.54	99.07	113.40
1	B	22	CYS	CA-CB-SG	7.36	127.25	114.00
1	A	198	GLN	CA-CB-CG	6.41	127.50	113.40
1	A	46	PRO	N-CA-C	6.36	128.64	112.10
1	B	215	CYS	CA-CB-SG	6.36	125.44	114.00
1	B	22	CYS	CB-CA-C	6.32	123.03	110.40
1	A	174	ASN	CB-CA-C	6.22	122.84	110.40
1	B	189	TRP	CA-CB-CG	6.11	125.31	113.70
2	P	1	HIS	CA-CB-CG	5.91	123.65	113.60
1	B	184	LEU	CA-CB-CG	5.59	128.16	115.30
1	B	136	LEU	CA-CB-CG	5.56	128.08	115.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	215	CYS	N-CA-C	-5.50	96.16	111.00
1	A	56	ARG	NE-CZ-NH1	5.27	122.94	120.30
1	B	56	ARG	NE-CZ-NH1	5.24	122.92	120.30
1	B	215	CYS	N-CA-CB	5.05	119.69	110.60
1	A	214	GLU	CA-C-N	-5.01	106.17	117.20

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	32	TYR	Sidechain

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	132	0
1	B	1605	0	1540	93	0
2	P	21	0	16	7	0
All	All	3231	0	3096	221	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 35.

All (221) 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:B:10:ALA:HB1	1:B:18:VAL:HG11	1.51	0.92
1:B:9:SER:HB2	1:B:147:ALA:HB3	1.51	0.91
1:A:135:THR:HG22	1:B:122:PHE:HZ	1.37	0.89
1:B:187:GLU:HA	1:B:190:LYS:HE2	1.57	0.87
1:A:18:VAL:HG12	1:A:77:VAL:HB	1.58	0.85
1:B:58:SER:C	1:B:60:VAL:H	1.86	0.79
1:A:190:LYS:HZ1	1:A:212:PRO:HB3	1.48	0.78
1:A:150:VAL:HG22	1:A:199:VAL:HG22	1.65	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:56:ARG:NH1	1:A:56:ARG:HB3	1.99	0.78
1:A:49:ILE:O	1:A:57:PRO:HD2	1.82	0.77
1:B:13:SER:O	1:B:14:LEU:HB2	1.85	0.75
1:A:51:TYR:CD1	1:A:57:PRO:HD3	2.22	0.74
1:A:14:LEU:HD23	1:A:109:VAL:HG11	1.70	0.73
1:A:66:GLY:HA2	1:A:74:SER:O	1.90	0.72
1:B:194:SER:HB2	1:B:209:THR:CG2	2.20	0.72
1:A:190:LYS:NZ	1:A:212:PRO:HB3	2.04	0.71
1:A:115:ALA:HB3	1:A:175:LYS:HZ1	1.55	0.71
1:A:6:GLN:HE22	1:A:90:CYS:HB3	1.56	0.70
1:A:10:ALA:HB3	1:A:107:VAL:HG22	1.73	0.69
1:A:65:SER:O	1:A:75:LEU:HA	1.94	0.68
1:B:169:SER:O	1:B:176:TYR:HA	1.94	0.67
1:B:6:GLN:HG2	1:B:22:CYS:HB2	1.75	0.67
1:A:5:THR:HB	1:A:23:THR:HB	1.75	0.66
1:B:49:ILE:HD11	1:B:75:LEU:HD11	1.77	0.66
1:A:12:GLY:O	1:A:110:LEU:HG	1.95	0.66
1:A:212:PRO:O	1:A:213:THR:HG23	1.96	0.65
1:B:38:TYR:HE1	1:B:91:SER:HB2	1.61	0.65
1:A:102:GLY:O	1:A:104:GLY:N	2.31	0.64
1:B:32:TYR:CE2	1:B:34:TYR:HB2	2.33	0.64
1:A:115:ALA:HB3	1:A:175:LYS:NZ	2.11	0.63
1:B:4:LEU:HD21	1:B:29:VAL:HG21	1.80	0.63
1:B:10:ALA:HB3	1:B:107:VAL:HG22	1.80	0.63
1:A:154:ALA:O	1:A:155:ASP:C	2.37	0.63
1:A:39:GLN:HG3	1:A:88:TYR:CE2	2.34	0.63
1:A:171:GLN:NE2	1:A:177:ALA:HB2	2.15	0.62
1:A:63:ARG:NH1	1:A:63:ARG:HA	2.14	0.62
1:A:64:PHE:HA	1:A:76:THR:O	1.99	0.62
1:A:184:LEU:HD21	1:A:195:TYR:CE2	2.34	0.62
1:A:3:ALA:HA	1:A:101:PHE:O	2.01	0.61
1:A:26:SER:O	1:A:27:SER:HB2	1.99	0.60
1:A:64:PHE:CD1	1:A:77:VAL:HG22	2.36	0.60
1:A:171:GLN:HE21	1:A:177:ALA:HB2	1.67	0.60
1:B:146:GLY:O	1:B:168:PRO:HG2	2.02	0.60
1:B:23:THR:HG23	1:B:71:ASN:O	2.02	0.59
1:B:127:GLU:O	1:B:130:GLN:N	2.36	0.59
1:B:40:GLN:O	1:B:86:ALA:HB1	2.01	0.59
1:A:150:VAL:HG22	1:A:199:VAL:CG2	2.32	0.59
1:A:65:SER:OG	1:A:76:THR:HB	2.03	0.59
1:A:135:THR:HG22	1:B:122:PHE:CZ	2.29	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:149:THR:HG23	1:A:200:THR:HB	1.83	0.58
1:A:6:GLN:NE2	1:A:90:CYS:HB3	2.18	0.58
1:A:35:VAL:HA	1:A:91:SER:O	2.03	0.58
1:A:116:ASN:OD1	1:A:201:HIS:HD2	1.87	0.58
1:B:165:THR:HG22	1:B:166:THR:H	1.69	0.58
1:B:35:VAL:HA	1:B:91:SER:O	2.04	0.57
1:B:148:VAL:HG12	1:B:201:HIS:HB2	1.87	0.57
1:A:163:VAL:HA	1:A:181:TYR:O	2.05	0.57
2:P:1:HIS:O	2:P:3:NH2:N	2.36	0.57
1:B:185:THR:O	1:B:188:GLN:HG2	2.05	0.57
1:A:148:VAL:HG23	1:A:201:HIS:HB2	1.87	0.57
1:A:49:ILE:HG13	1:A:50:ILE:N	2.20	0.57
1:A:34:TYR:O	1:A:92:SER:HA	2.05	0.56
1:B:32:TYR:CE2	1:B:93:TYR:O	2.59	0.56
1:B:149:THR:HB	1:B:200:THR:HB	1.88	0.56
1:A:50:ILE:HD12	1:A:56:ARG:HG3	1.88	0.56
1:A:6:GLN:NE2	1:A:22:CYS:SG	2.79	0.55
1:B:170:LYS:HE3	1:B:176:TYR:CZ	2.41	0.55
1:B:4:LEU:HD12	1:B:100:VAL:HG12	1.88	0.55
1:B:64:PHE:CE1	1:B:77:VAL:HG22	2.42	0.54
1:A:56:ARG:CZ	1:A:56:ARG:HB3	2.37	0.54
1:A:154:ALA:O	1:A:156:GLY:N	2.41	0.54
1:B:32:TYR:HE2	1:B:34:TYR:HB2	1.71	0.54
1:A:28:ASP:HA	1:A:92:SER:OG	2.08	0.54
1:A:93:TYR:HE1	1:A:97:ASP:HA	1.73	0.53
1:B:195:TYR:HB2	1:B:210:VAL:HG23	1.90	0.53
1:A:119:VAL:HG12	1:A:208:LYS:HG3	1.91	0.53
1:A:28:ASP:O	1:A:30:GLY:N	2.42	0.53
1:B:193:ARG:O	1:B:212:PRO:HD2	2.08	0.52
1:A:20:ILE:O	1:A:74:SER:HA	2.09	0.52
1:B:171:GLN:OE1	1:B:177:ALA:HB2	2.09	0.52
1:A:27:SER:HA	1:A:32:TYR:CE1	2.44	0.52
1:A:81:GLN:OE1	1:A:83:GLU:HB2	2.09	0.52
1:A:121:LEU:HD11	1:A:152:TRP:CZ3	2.44	0.52
1:A:9:SER:O	1:A:10:ALA:HB2	2.10	0.51
1:A:14:LEU:CD2	1:A:109:VAL:HG11	2.40	0.51
1:B:4:LEU:CD2	1:B:29:VAL:HG21	2.40	0.51
1:A:159:VAL:HG11	1:A:182:LEU:HD21	1.93	0.51
1:B:194:SER:HB2	1:B:209:THR:HG22	1.91	0.51
1:A:56:ARG:HH11	1:A:56:ARG:CB	2.23	0.51
1:B:170:LYS:HE3	1:B:176:TYR:OH	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:121:LEU:HG	1:B:210:VAL:HG13	1.93	0.51
1:B:96:SER:O	1:B:97:ASP:HB3	2.11	0.51
1:A:56:ARG:HG2	1:A:60:VAL:CG1	2.41	0.50
1:A:184:LEU:HD21	1:A:195:TYR:CZ	2.45	0.50
1:A:37:TRP:H	1:A:50:ILE:HG22	1.76	0.50
1:A:52:GLU:OE1	1:A:55:LYS:HD2	2.11	0.50
1:B:50:ILE:HD12	1:B:75:LEU:HD12	1.94	0.50
1:A:11:SER:HB2	1:A:110:LEU:HD21	1.92	0.50
1:A:36:SER:CB	1:A:51:TYR:HA	2.41	0.50
1:A:41:HIS:CE1	1:A:86:ALA:HB2	2.46	0.50
1:B:10:ALA:HB1	1:B:18:VAL:CG1	2.35	0.50
1:B:34:TYR:CE1	2:P:1:HIS:HB3	2.46	0.50
1:B:166:THR:HG22	1:B:167:LYS:O	2.12	0.50
1:A:194:SER:HB3	1:A:209:THR:CG2	2.42	0.49
1:B:112:GLN:HG2	1:B:144:TYR:CE2	2.47	0.49
1:A:122:PHE:O	1:A:136:LEU:HD23	2.12	0.49
1:B:47:LYS:HB3	1:B:47:LYS:NZ	2.26	0.49
1:B:184:LEU:HD12	1:B:188:GLN:OE1	2.12	0.49
1:B:164:GLU:O	1:B:180:SER:HA	2.13	0.49
1:A:121:LEU:HD13	1:A:197:CYS:HB2	1.95	0.49
1:B:58:SER:C	1:B:60:VAL:N	2.58	0.49
1:B:18:VAL:CG2	1:B:80:LEU:HD11	2.43	0.48
1:A:132:ASN:ND2	1:A:186:PRO:HG2	2.27	0.48
1:A:144:TYR:HB2	1:A:175:LYS:NZ	2.28	0.48
1:B:11:SER:OG	1:B:145:PRO:HG2	2.14	0.48
1:A:36:SER:OG	1:A:51:TYR:HA	2.13	0.48
1:B:10:ALA:HB3	1:B:107:VAL:HA	1.95	0.48
1:A:145:PRO:O	1:A:201:HIS:HE1	1.97	0.48
1:A:56:ARG:HB3	1:A:56:ARG:HH11	1.72	0.48
1:A:33:ASN:HA	1:A:68:LYS:HZ3	1.79	0.48
1:A:48:VAL:HG21	1:B:99:PHE:HD2	1.79	0.48
1:A:51:TYR:HD1	1:A:57:PRO:HD3	1.73	0.48
1:A:19:THR:HA	1:A:75:LEU:O	2.13	0.47
1:A:211:ALA:O	1:A:213:THR:N	2.47	0.47
1:B:20:ILE:O	1:B:74:SER:HA	2.14	0.47
1:A:197:CYS:O	1:A:207:GLU:HA	2.15	0.47
1:A:25:THR:HB	1:A:28:ASP:OD2	2.14	0.47
1:B:34:TYR:HD1	1:B:34:TYR:HA	1.58	0.47
1:B:53:VAL:HG13	1:B:68:LYS:HB2	1.97	0.47
1:B:10:ALA:CB	1:B:107:VAL:HG22	2.44	0.47
1:A:38:TYR:CE2	1:A:48:VAL:HG22	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:46:PRO:O	1:A:47:LYS:HB2	2.15	0.47
1:A:46:PRO:HG2	1:B:101:PHE:CE2	2.50	0.47
1:A:62:ASP:O	1:A:63:ARG:HD2	2.15	0.46
1:B:18:VAL:HG23	1:B:80:LEU:HD11	1.97	0.46
1:A:126:SER:HA	1:A:129:LEU:HD12	1.96	0.46
1:B:196:SER:HA	1:B:208:LYS:O	2.16	0.46
1:B:155:ASP:OD1	1:B:192:HIS:HB3	2.16	0.46
1:A:87:ASP:HA	1:A:105:THR:O	2.15	0.46
1:B:145:PRO:O	1:B:201:HIS:HE1	1.99	0.46
1:A:121:LEU:HD23	1:A:210:VAL:HG13	1.96	0.46
1:A:38:TYR:CD1	1:A:38:TYR:N	2.83	0.46
1:A:124:PRO:HD3	1:A:136:LEU:HG	1.98	0.46
1:B:153:LYS:HA	1:B:158:PRO:HA	1.97	0.46
1:A:38:TYR:CD2	1:A:48:VAL:HG22	2.50	0.45
1:B:140:ILE:HG12	1:B:199:VAL:HG21	1.98	0.45
2:P:0:ACE:O	2:P:1:HIS:CG	2.69	0.45
1:A:29:VAL:HG11	1:A:72:THR:CA	2.46	0.45
1:B:187:GLU:HA	1:B:190:LYS:CE	2.38	0.45
1:A:99:PHE:CE2	2:P:2:DPR:HB2	2.52	0.45
1:A:37:TRP:N	1:A:50:ILE:HG22	2.31	0.45
1:A:173:ASN:CG	1:A:175:LYS:HG3	2.36	0.44
1:A:37:TRP:HD1	1:A:50:ILE:CG2	2.30	0.44
1:A:27:SER:O	1:A:32:TYR:HD1	2.01	0.44
1:B:186:PRO:O	1:B:189:TRP:HB3	2.17	0.44
1:A:53:VAL:HG13	1:A:54:ASN:N	2.33	0.44
1:A:56:ARG:HG2	1:A:60:VAL:HG13	1.98	0.44
1:A:63:ARG:HH11	1:A:63:ARG:HA	1.82	0.44
1:A:25:THR:H	1:A:28:ASP:HB2	1.83	0.44
1:A:37:TRP:CE2	1:A:75:LEU:HB2	2.52	0.44
1:A:6:GLN:HG2	1:A:105:THR:OG1	2.17	0.44
1:A:169:SER:O	1:A:176:TYR:HA	2.16	0.44
1:B:94:GLU:HB3	1:B:98:ASN:HB2	1.99	0.44
1:B:153:LYS:HD2	1:B:207:GLU:OE2	2.17	0.44
1:B:6:GLN:HA	1:B:7:PRO:HD2	1.76	0.44
1:A:20:ILE:H	1:A:20:ILE:HG12	1.70	0.44
1:B:40:GLN:HA	1:B:44:LYS:HZ2	1.83	0.44
1:A:85:GLU:HA	1:A:107:VAL:O	2.19	0.43
1:B:47:LYS:HD2	1:B:47:LYS:O	2.18	0.43
1:A:49:ILE:CG1	1:A:50:ILE:N	2.82	0.43
1:A:65:SER:O	1:A:75:LEU:HD23	2.19	0.43
1:A:85:GLU:HG2	1:A:109:VAL:HG23	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:20:ILE:HG22	1:B:37:TRP:CH2	2.53	0.43
1:B:41:HIS:HB2	1:B:44:LYS:HD2	2.00	0.43
1:B:41:HIS:O	1:B:42:ALA:C	2.56	0.43
1:B:108:THR:HG21	1:B:145:PRO:HB3	2.01	0.43
1:A:48:VAL:HG21	1:B:99:PHE:CD2	2.54	0.43
1:A:56:ARG:NH1	1:A:56:ARG:CB	2.75	0.42
1:B:47:LYS:O	1:B:49:ILE:HG23	2.18	0.42
1:B:185:THR:O	1:B:188:GLN:N	2.52	0.42
1:A:194:SER:HB3	1:A:209:THR:HG22	2.02	0.42
1:B:32:TYR:HH	1:B:93:TYR:HD2	1.64	0.42
1:B:57:PRO:O	1:B:59:GLY:N	2.47	0.42
1:B:49:ILE:CD1	1:B:75:LEU:HD11	2.46	0.42
1:B:19:THR:C	1:B:20:ILE:HD12	2.39	0.42
1:A:99:PHE:HZ	2:P:1:HIS:HD2	1.66	0.42
1:A:2:SER:O	1:A:100:VAL:HG22	2.19	0.42
1:B:40:GLN:HA	1:B:44:LYS:NZ	2.34	0.42
1:B:69:SER:O	1:B:70:GLY:C	2.57	0.42
1:B:94:GLU:CB	1:B:98:ASN:HB2	2.50	0.42
1:A:99:PHE:HZ	2:P:1:HIS:CD2	2.37	0.42
1:A:105:THR:HG22	1:A:107:VAL:HG23	2.02	0.42
1:A:27:SER:H	1:A:31:GLY:H	1.68	0.42
1:B:20:ILE:HD13	1:B:75:LEU:HD23	2.02	0.41
1:B:56:ARG:HA	1:B:57:PRO:HD3	1.85	0.41
1:A:171:GLN:O	1:A:172:SER:HB2	2.20	0.41
1:A:45:ALA:HA	1:A:46:PRO:HD3	1.87	0.41
1:B:63:ARG:NH2	1:B:84:ASP:OD1	2.52	0.41
1:A:165:THR:HA	1:A:179:SER:O	2.21	0.41
1:A:37:TRP:HD1	1:A:50:ILE:HG21	1.84	0.41
1:B:135:THR:HG23	1:B:182:LEU:O	2.20	0.41
1:B:153:LYS:HE3	1:B:198:GLN:OE1	2.21	0.41
1:A:135:THR:HA	1:A:182:LEU:O	2.20	0.41
1:A:53:VAL:CG1	1:A:54:ASN:N	2.84	0.41
1:A:99:PHE:CZ	2:P:1:HIS:HD2	2.38	0.41
1:B:56:ARG:O	1:B:56:ARG:HD2	2.20	0.41
1:A:27:SER:HA	1:A:32:TYR:HE1	1.84	0.41
1:A:68:LYS:HA	1:A:73:ALA:HA	2.02	0.41
1:A:182:LEU:HD23	1:A:182:LEU:HA	1.89	0.41
1:A:1:PRO:O	1:A:2:SER:CB	2.69	0.41
1:A:29:VAL:HG12	1:A:70:GLY:O	2.21	0.41
1:A:14:LEU:HA	1:A:80:LEU:O	2.21	0.41
1:A:51:TYR:CE1	1:A:57:PRO:HD3	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:4:LEU:HB3	1:A:22:CYS:SG	2.60	0.40
1:B:120:THR:OG1	1:B:139:LEU:HD22	2.21	0.40
1:A:167:LYS:HG2	1:B:169:SER:OG	2.21	0.40
1:B:187:GLU:CA	1:B:190:LYS:HE2	2.40	0.40
1:B:141:SER:HB2	1:B:171:GLN:OE1	2.22	0.40
1:B:38:TYR:HH	1:B:99:PHE:HE2	1.65	0.40
1:A:136:LEU:O	1:A:181:TYR:HA	2.22	0.40
1:A:50:ILE:HG13	1:A:55:LYS:O	2.22	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%)	165 (77%)	32 (15%)	17 (8%)	1	1
1	B	214/216 (99%)	181 (85%)	25 (12%)	8 (4%)	3	7
2	P	1/4 (25%)	0	0	1 (100%)	0	0
All	All	429/436 (98%)	346 (81%)	57 (13%)	26 (6%)	1	2

All (26) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	2	SER
1	A	10	ALA
1	A	27	SER
1	A	29	VAL
1	A	103	THR
1	A	142	ASP
1	A	155	ASP
1	B	14	LEU
1	B	42	ALA

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Mol	Chain	Res	Type
1	B	95	GLY
1	A	4	LEU
1	A	215	CYS
1	B	58	SER
1	B	70	GLY
2	P	1	HIS
1	A	57	PRO
1	A	202	GLU
1	A	212	PRO
1	A	213	THR
1	B	2	SER
1	B	26	SER
1	A	172	SER
1	A	26	SER
1	A	53	VAL
1	A	5	THR
1	B	79	GLY

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%)	134 (74%)	47 (26%)	0 1
1	B	181/181 (100%)	149 (82%)	32 (18%)	2 4
2	P	1/1 (100%)	0	1 (100%)	0 0
All	All	363/363 (100%)	283 (78%)	80 (22%)	1 2

All (80) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	THR
1	A	6	GLN
1	A	8	PRO
1	A	9	SER
1	A	11	SER

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Mol	Chain	Res	Type
1	A	18	VAL
1	A	20	ILE
1	A	27	SER
1	A	28	ASP
1	A	32	TYR
1	A	35	VAL
1	A	38	TYR
1	A	46	PRO
1	A	47	LYS
1	A	50	ILE
1	A	53	VAL
1	A	56	ARG
1	A	58	SER
1	A	62	ASP
1	A	63	ARG
1	A	67	SER
1	A	75	LEU
1	A	81	GLN
1	A	87	ASP
1	A	94	GLU
1	A	98	ASN
1	A	103	THR
1	A	110	LEU
1	A	114	LYS
1	A	125	SER
1	A	130	GLN
1	A	143	PHE
1	A	149	THR
1	A	155	ASP
1	A	165	THR
1	A	171	GLN
1	A	173	ASN
1	A	175	LYS
1	A	179	SER
1	A	182	LEU
1	A	184	LEU
1	A	190	LYS
1	A	193	ARG
1	A	196	SER
1	A	198	GLN
1	A	205	THR
1	A	209	THR

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Mol	Chain	Res	Type
1	B	16	GLN
1	B	17	SER
1	B	21	SER
1	B	22	CYS
1	B	25	THR
1	B	27	SER
1	B	34	TYR
1	B	47	LYS
1	B	56	ARG
1	B	58	SER
1	B	62	ASP
1	B	65	SER
1	B	67	SER
1	B	68	LYS
1	B	69	SER
1	B	83	GLU
1	B	91	SER
1	B	98	ASN
1	B	103	THR
1	B	106	LYS
1	B	122	PHE
1	B	125	SER
1	B	128	GLU
1	B	130	GLN
1	B	143	PHE
1	B	148	VAL
1	B	165	THR
1	B	179	SER
1	B	182	LEU
1	B	184	LEU
1	B	185	THR
1	B	190	LYS
2	P	1	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	39	GLN
1	A	41	HIS
1	A	98	ASN
1	A	171	GLN
1	A	173	ASN

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Mol	Chain	Res	Type
1	A	198	GLN
1	A	201	HIS
1	B	40	GLN
1	B	98	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](#)

1 non-standard protein/DNA/RNA residue is modelled in this entry.

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.

5.5 Carbohydrates [i](#)

There are no carbohydrates 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.