



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 08:26 PM GMT

PDB ID : 1KCS  
Title : CRYSTAL STRUCTURE OF ANTIBODY PC282 IN COMPLEX WITH PS1 PEPTIDE  
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Deposited on : 2001-11-11  
Resolution : 2.50 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](mailto:validation@mail.wwpdb.org)  
A user guide is available at  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

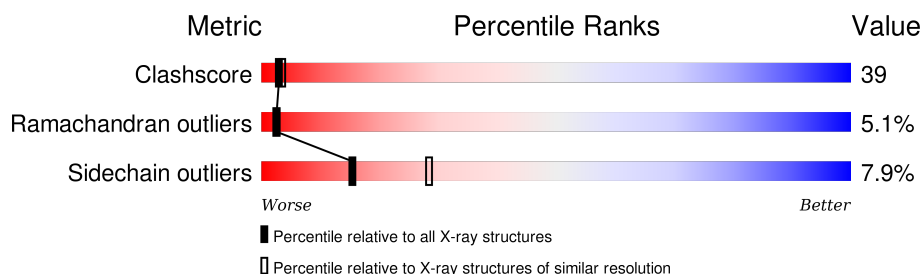
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.50 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	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	L	214	
2	H	217	
3	P	15	

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3348 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 PC282 IMMUNOGLOBULIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	214	Total	C	N	O	S	0	0	0
			1611	996	269	337	9			

- Molecule 2 is a protein called PC282 IMMUNOGLOBULIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	217	Total	C	N	O	S	0	0	0
			1600	1010	260	325	5			

- Molecule 3 is a protein called PS1 peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	P	7	Total	C	N	O	0	0	0
			52	34	8	10			

- Molecule 4 is water.

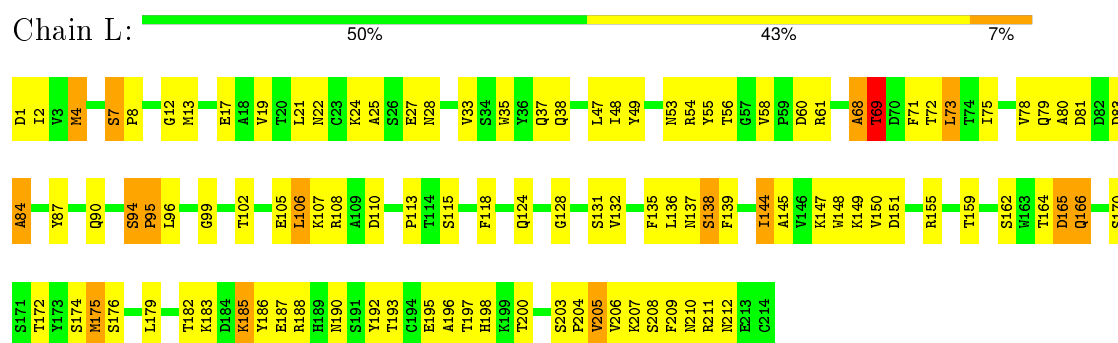
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	H	43	Total	O	0	0
			43	43		
4	L	41	Total	O	0	0
			41	41		
4	P	1	Total	O	0	0
			1	1		

### 3 Residue-property plots

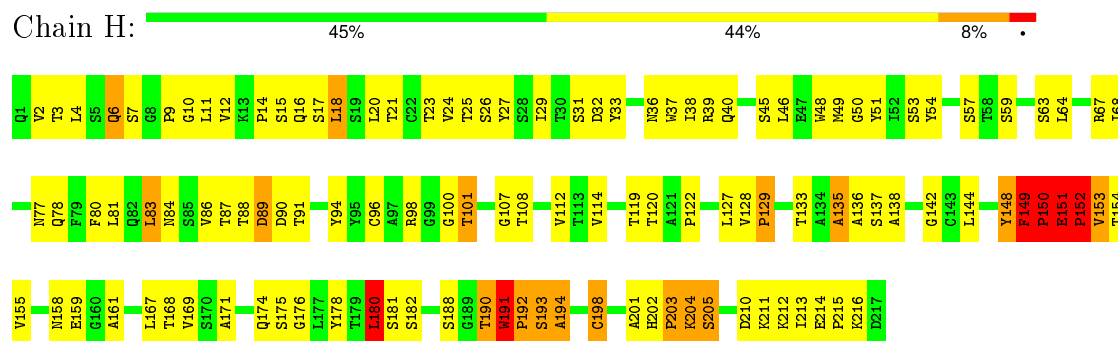
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according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

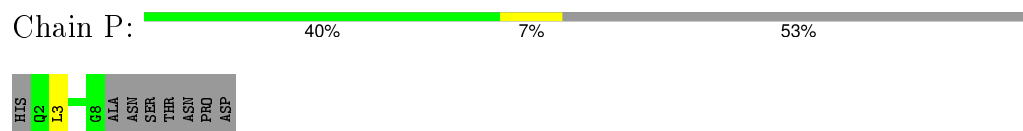
#### • Molecule 1: PC282 IMMUNOGLOBULIN



#### • Molecule 2: PC282 IMMUNOGLOBULIN



#### • Molecule 3: PS1 peptide



## 4 Data and refinement statistics

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

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	52.60 Å 69.52 Å 59.36 Å 90.00° 95.62° 90.00°	Depositor
Resolution (Å)	50.00 – 2.50	Depositor
% Data completeness (in resolution range)	84.3 (50.00-2.50)	Depositor
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	CNS 0.5	Depositor
R, $R_{free}$	0.209 , 0.277	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3348	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	31.0	wwPDB-VP

## 5 Model quality

### 5.1 Standard geometry

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	L	0.40	0/1646	0.68	0/2233
2	H	0.57	1/1640 (0.1%)	1.16	14/2249 (0.6%)
3	P	0.48	0/53	0.55	0/71
All	All	0.49	1/3339 (0.0%)	0.94	14/4553 (0.3%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	191	TRP	CB-CG	-7.66	1.36	1.50

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	151	GLU	C-N-CD	-15.80	85.85	120.60
2	H	152	PRO	CA-N-CD	-14.72	90.89	111.50
2	H	151	GLU	C-N-CA	13.03	176.71	122.00
2	H	152	PRO	N-CD-CG	10.27	118.61	103.20
2	H	152	PRO	N-CA-CB	9.40	114.58	103.30
2	H	190	THR	C-N-CA	-8.08	101.51	121.70
2	H	191	TRP	CB-CA-C	7.32	125.04	110.40
2	H	150	PRO	CA-N-CD	-7.11	101.55	111.50
2	H	192	PRO	CA-N-CD	-6.84	101.93	111.50
2	H	149	PHE	C-N-CA	-6.08	96.48	122.00
2	H	148	TYR	C-N-CA	-5.94	106.84	121.70
2	H	149	PHE	CA-C-N	5.73	133.16	117.10
2	H	180	LEU	CA-CB-CG	5.62	128.22	115.30
2	H	191	TRP	CE2-CD2-CG	-5.00	103.30	107.30

There are no chirality outliers.

There are no planarity outliers.

## 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	L	1611	0	1535	136	0
2	H	1600	0	1555	125	0
3	P	52	0	46	1	0
4	H	43	0	0	13	0
4	L	41	0	0	11	0
4	P	1	0	0	0	0
All	All	3348	0	3136	250	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 39.

All (250) 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:L:94:SER:OG	1:L:95:PRO:HD3	1.29	1.27
1:L:94:SER:OG	1:L:95:PRO:CD	1.86	1.24
1:L:87:TYR:HD1	4:L:222:HOH:O	1.34	1.06
1:L:38:GLN:HE22	2:H:40:GLN:HE22	1.07	1.02
1:L:7:SER:HB3	1:L:8:PRO:HD3	1.42	1.00
1:L:37:GLN:HB2	1:L:47:LEU:HD11	1.45	0.96
2:H:3:THR:HG23	2:H:25:THR:HB	1.45	0.95
2:H:122:PRO:HB3	2:H:148:TYR:HB3	1.48	0.95
2:H:91:THR:HG22	2:H:114:VAL:H	1.29	0.95
2:H:152:PRO:O	2:H:153:VAL:HG13	1.75	0.86
1:L:137:ASN:HD21	2:H:167:LEU:HD13	1.40	0.84
1:L:144:ILE:HG12	4:L:234:HOH:O	1.78	0.83
2:H:24:VAL:HG13	2:H:77:ASN:ND2	1.94	0.83
2:H:149:PHE:O	2:H:150:PRO:O	1.96	0.82
1:L:94:SER:OG	1:L:95:PRO:HD2	1.77	0.82
1:L:68:ALA:O	1:L:69:THR:HB	1.79	0.81
2:H:88:THR:O	2:H:91:THR:HG23	1.81	0.80
1:L:13:MET:HG3	1:L:19:VAL:HG12	1.63	0.80
2:H:136:ALA:O	2:H:188:SER:HB3	1.84	0.78
1:L:38:GLN:NE2	2:H:40:GLN:HE22	1.81	0.78
1:L:113:PRO:HB3	1:L:139:PHE:HB3	1.63	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:1:ASP:CG	1:L:2:ILE:H	1.85	0.78
1:L:94:SER:HG	1:L:95:PRO:HD3	1.46	0.77
1:L:28:ASN:HD22	1:L:68:ALA:HB1	1.49	0.77
1:L:38:GLN:HE22	2:H:40:GLN:NE2	1.83	0.77
2:H:180:LEU:HD23	2:H:181:SER:N	1.99	0.77
1:L:137:ASN:ND2	2:H:167:LEU:HD13	2.02	0.75
2:H:98:ARG:HB2	4:H:226:HOH:O	1.87	0.75
2:H:151:GLU:HG3	2:H:178:TYR:CE2	2.22	0.74
1:L:136:LEU:HD23	1:L:175:MET:CE	2.18	0.74
1:L:7:SER:CB	1:L:8:PRO:HD3	2.18	0.74
1:L:190:ASN:HD21	1:L:212:ASN:HB3	1.53	0.73
2:H:101:THR:O	2:H:101:THR:CG2	2.35	0.73
2:H:10:GLY:HA3	2:H:204:LYS:HE2	1.71	0.73
2:H:204:LYS:N	2:H:204:LYS:HD2	2.04	0.72
2:H:149:PHE:O	2:H:150:PRO:C	2.12	0.71
2:H:4:LEU:HD11	4:H:226:HOH:O	1.91	0.71
1:L:1:ASP:CG	1:L:95:PRO:HG2	2.11	0.70
2:H:16:GLN:O	2:H:86:VAL:HG22	1.92	0.69
2:H:149:PHE:C	2:H:150:PRO:O	2.31	0.69
1:L:113:PRO:HB3	1:L:139:PHE:CD2	2.28	0.69
2:H:91:THR:HG22	2:H:114:VAL:N	2.07	0.68
1:L:145:ALA:HB3	1:L:197:THR:CG2	2.23	0.68
1:L:196:ALA:HB1	4:L:234:HOH:O	1.94	0.67
1:L:94:SER:CB	1:L:95:PRO:HD3	2.23	0.67
1:L:12:GLY:HA2	1:L:105:GLU:O	1.93	0.67
2:H:149:PHE:O	2:H:149:PHE:CD1	2.48	0.67
2:H:67:ARG:HH22	2:H:90:ASP:CG	1.97	0.67
1:L:164:THR:HG23	1:L:174:SER:O	1.94	0.66
1:L:179:LEU:HD11	4:L:231:HOH:O	1.94	0.66
2:H:108:THR:HG22	2:H:108:THR:O	1.94	0.66
2:H:191:TRP:HZ3	2:H:213:ILE:O	1.78	0.66
2:H:119:THR:HG23	2:H:150:PRO:HD2	1.78	0.66
2:H:202:HIS:ND1	2:H:205:SER:HB3	2.10	0.65
1:L:7:SER:HB3	1:L:8:PRO:CD	2.24	0.65
2:H:100:GLY:HA2	3:P:3:LEU:HD22	1.78	0.65
1:L:136:LEU:HD23	1:L:175:MET:HE1	1.78	0.65
1:L:145:ALA:HB3	1:L:197:THR:HG22	1.79	0.65
1:L:4:MET:SD	1:L:90:GLN:HB3	2.37	0.64
2:H:36:ASN:OD1	2:H:51:TYR:HB3	1.97	0.64
1:L:8:PRO:HD2	4:L:227:HOH:O	1.97	0.64
1:L:28:ASN:HD22	1:L:68:ALA:CB	2.10	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:17:SER:CB	2:H:84:ASN:HA	2.27	0.64
1:L:1:ASP:HB2	1:L:95:PRO:HG2	1.77	0.64
1:L:48:ILE:HD13	1:L:54:ARG:HA	1.80	0.63
2:H:17:SER:HB3	2:H:84:ASN:HA	1.80	0.63
2:H:4:LEU:HD21	4:H:226:HOH:O	1.97	0.63
2:H:6:GLN:OE1	2:H:107:GLY:HA3	1.99	0.63
1:L:150:VAL:O	1:L:151:ASP:HB2	1.98	0.63
2:H:18:LEU:HD22	2:H:20:LEU:HD21	1.79	0.63
1:L:17:GLU:O	1:L:78:VAL:HG23	1.99	0.63
1:L:165:ASP:OD2	1:L:165:ASP:N	2.32	0.62
2:H:168:THR:HG22	2:H:169:VAL:N	2.14	0.62
1:L:128:GLY:HA2	1:L:183:LYS:HZ2	1.63	0.62
2:H:204:LYS:CD	2:H:204:LYS:N	2.62	0.62
1:L:203:SER:HB2	1:L:204:PRO:HD2	1.81	0.62
2:H:39:ARG:HB3	2:H:49:MET:HE1	1.82	0.61
1:L:1:ASP:CB	1:L:95:PRO:HG2	2.28	0.61
1:L:25:ALA:HB3	1:L:69:THR:HA	1.81	0.61
2:H:191:TRP:CZ3	2:H:213:ILE:O	2.54	0.61
2:H:2:VAL:HG21	2:H:98:ARG:NH2	2.16	0.61
2:H:18:LEU:HD23	2:H:112:VAL:HG11	1.82	0.61
2:H:128:VAL:HG22	2:H:213:ILE:CG2	2.30	0.60
1:L:165:ASP:O	1:L:166:GLN:O	2.18	0.60
2:H:14:PRO:O	2:H:15:SER:HB2	2.01	0.60
1:L:137:ASN:HD21	2:H:167:LEU:CD1	2.10	0.60
2:H:6:GLN:NE2	2:H:6:GLN:H	2.00	0.60
1:L:164:THR:HG22	2:H:168:THR:O	2.01	0.60
2:H:39:ARG:HB3	2:H:49:MET:CE	2.32	0.60
2:H:149:PHE:CG	2:H:149:PHE:O	2.53	0.59
1:L:8:PRO:O	1:L:102:THR:HG23	2.02	0.59
1:L:19:VAL:HG21	1:L:75:ILE:HD12	1.84	0.59
2:H:101:THR:HG22	2:H:101:THR:O	2.01	0.59
1:L:4:MET:O	1:L:99:GLY:HA2	2.01	0.59
2:H:127:LEU:HB2	2:H:142:GLY:CA	2.32	0.59
1:L:1:ASP:CG	1:L:2:ILE:N	2.55	0.59
1:L:136:LEU:HD23	1:L:175:MET:HE2	1.84	0.59
2:H:158:ASN:HB3	2:H:161:ALA:HB3	1.84	0.59
1:L:200:THR:HA	4:L:250:HOH:O	2.02	0.58
2:H:24:VAL:HG11	2:H:29:ILE:HG23	1.84	0.58
2:H:10:GLY:CA	2:H:204:LYS:HE2	2.34	0.58
1:L:128:GLY:HA2	1:L:183:LYS:NZ	2.19	0.58
1:L:110:ASP:HB3	1:L:200:THR:HG22	1.85	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:32:ASP:O	2:H:33:TYR:HB2	2.03	0.57
2:H:3:THR:CG2	2:H:25:THR:HB	2.27	0.57
1:L:79:GLN:HE21	1:L:79:GLN:HA	1.69	0.57
2:H:212:LYS:HE2	2:H:214:GLU:HG2	1.86	0.57
1:L:108:ARG:NH1	1:L:172:THR:HG23	2.19	0.57
1:L:78:VAL:HG12	1:L:106:LEU:HD11	1.85	0.57
1:L:137:ASN:ND2	1:L:138:SER:H	2.04	0.56
1:L:182:THR:OG1	1:L:185:LYS:HG3	2.05	0.56
2:H:20:LEU:N	2:H:20:LEU:HD22	2.21	0.56
2:H:9:PRO:HD2	4:H:228:HOH:O	2.06	0.55
1:L:149:LYS:HB2	1:L:193:THR:HB	1.88	0.55
1:L:19:VAL:HG22	1:L:75:ILE:HB	1.88	0.55
2:H:127:LEU:HB2	2:H:142:GLY:HA3	1.88	0.55
1:L:4:MET:HE3	1:L:90:GLN:HG2	1.88	0.55
1:L:118:PHE:HE2	2:H:129:PRO:HD3	1.71	0.55
1:L:147:LYS:HD2	1:L:149:LYS:NZ	2.21	0.55
1:L:198:HIS:HD2	1:L:200:THR:OG1	1.90	0.55
2:H:53:SER:OG	2:H:57:SER:HB2	2.07	0.55
2:H:88:THR:HG23	2:H:89:ASP:OD1	2.07	0.55
1:L:49:TYR:O	1:L:53:ASN:HB2	2.07	0.54
1:L:196:ALA:O	1:L:205:VAL:HG23	2.08	0.54
1:L:68:ALA:O	1:L:69:THR:CB	2.55	0.54
1:L:80:ALA:HA	1:L:106:LEU:HG	1.90	0.54
1:L:83:ASP:O	1:L:84:ALA:HB2	2.09	0.53
2:H:48:TRP:CZ2	2:H:50:GLY:HA2	2.43	0.53
2:H:154:THR:HB	2:H:201:ALA:HB3	1.91	0.53
1:L:78:VAL:CG1	1:L:106:LEU:HD11	2.39	0.53
1:L:162:SER:O	1:L:175:MET:HB2	2.08	0.52
1:L:72:THR:HG23	4:L:242:HOH:O	2.08	0.52
2:H:135:ALA:HB3	2:H:138:ALA:O	2.09	0.52
1:L:113:PRO:HB3	1:L:139:PHE:HD2	1.73	0.52
1:L:54:ARG:HD3	1:L:58:VAL:O	2.09	0.52
1:L:113:PRO:HB3	1:L:139:PHE:CB	2.39	0.52
2:H:129:PRO:O	2:H:216:LYS:HD3	2.10	0.51
1:L:21:LEU:HD12	4:L:225:HOH:O	2.10	0.51
1:L:135:PHE:CE2	1:L:137:ASN:HB2	2.45	0.51
2:H:7:SER:HA	4:H:219:HOH:O	2.10	0.51
1:L:7:SER:CB	1:L:8:PRO:CD	2.88	0.51
1:L:113:PRO:CB	1:L:139:PHE:HB3	2.36	0.50
2:H:10:GLY:HA3	2:H:204:LYS:CE	2.40	0.50
2:H:37:TRP:CD1	2:H:81:LEU:HB2	2.45	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:152:PRO:O	2:H:153:VAL:CG1	2.52	0.50
2:H:24:VAL:HG11	2:H:29:ILE:CG2	2.40	0.50
2:H:128:VAL:CG2	2:H:213:ILE:HG23	2.42	0.50
1:L:108:ARG:HD2	1:L:170:SER:O	2.11	0.50
1:L:49:TYR:CD2	2:H:101:THR:HG23	2.46	0.50
2:H:168:THR:HG22	2:H:169:VAL:H	1.75	0.50
2:H:101:THR:HG23	2:H:101:THR:O	2.10	0.50
2:H:6:GLN:HE21	2:H:6:GLN:H	1.59	0.50
2:H:23:THR:HA	2:H:78:GLN:HG2	1.94	0.50
1:L:124:GLN:HE22	1:L:131:SER:H	1.60	0.50
2:H:40:GLN:HB2	2:H:46:LEU:HD23	1.93	0.49
1:L:186:TYR:HA	1:L:192:TYR:OH	2.12	0.49
1:L:193:THR:HG23	1:L:208:SER:HB2	1.93	0.49
2:H:169:VAL:O	2:H:180:LEU:HG	2.11	0.49
1:L:1:ASP:HB2	1:L:95:PRO:CG	2.41	0.49
1:L:4:MET:HA	1:L:24:LYS:O	2.12	0.49
1:L:206:VAL:HG23	1:L:206:VAL:O	2.11	0.49
2:H:128:VAL:HG22	2:H:213:ILE:HG23	1.95	0.49
1:L:61:ARG:HG3	1:L:75:ILE:HG23	1.95	0.48
2:H:38:ILE:C	2:H:49:MET:HE2	2.34	0.48
2:H:94:TYR:CD1	2:H:94:TYR:N	2.82	0.48
1:L:195:GLU:HG3	1:L:206:VAL:HG12	1.94	0.48
1:L:193:THR:HG23	1:L:207:LYS:O	2.14	0.48
1:L:28:ASN:ND2	1:L:68:ALA:CB	2.76	0.48
2:H:67:ARG:HB3	2:H:83:LEU:CD2	2.43	0.48
2:H:193:SER:OG	2:H:194:ALA:N	2.47	0.48
1:L:61:ARG:O	1:L:75:ILE:HA	2.13	0.48
2:H:216:LYS:HB3	4:H:242:HOH:O	2.14	0.48
1:L:80:ALA:HA	1:L:106:LEU:CD2	2.44	0.48
2:H:168:THR:CG2	2:H:169:VAL:N	2.77	0.48
1:L:211:ARG:HH11	1:L:211:ARG:HG2	1.79	0.48
1:L:54:ARG:NE	1:L:60:ASP:HA	2.28	0.47
1:L:183:LYS:HB3	1:L:183:LYS:NZ	2.30	0.47
1:L:118:PHE:CE2	2:H:129:PRO:HD3	2.49	0.47
2:H:63:SER:OG	2:H:64:LEU:HD12	2.13	0.47
1:L:137:ASN:ND2	1:L:138:SER:N	2.62	0.47
1:L:185:LYS:O	1:L:188:ARG:HG2	2.14	0.47
2:H:26:SER:O	2:H:27:TYR:HB2	2.13	0.47
1:L:49:TYR:CG	2:H:101:THR:HG23	2.49	0.47
2:H:10:GLY:N	2:H:204:LYS:HE2	2.29	0.47
1:L:175:MET:HE3	1:L:176:SER:CA	2.43	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:36:ASN:O	2:H:96:CYS:HA	2.14	0.47
1:L:137:ASN:HD22	1:L:138:SER:H	1.61	0.47
2:H:108:THR:O	2:H:108:THR:CG2	2.62	0.47
2:H:67:ARG:NH2	4:H:229:HOH:O	2.48	0.47
1:L:33:VAL:HG21	1:L:71:PHE:CD1	2.49	0.47
2:H:11:LEU:HD13	2:H:12:VAL:N	2.31	0.46
2:H:190:THR:O	2:H:190:THR:CG2	2.62	0.46
4:L:226:HOH:O	2:H:181:SER:HB3	2.15	0.46
2:H:21:THR:HG23	2:H:80:PHE:CE2	2.51	0.46
2:H:49:MET:HE1	2:H:94:TYR:CD2	2.51	0.46
1:L:209:PHE:C	1:L:210:ASN:HD22	2.19	0.46
2:H:67:ARG:CZ	4:H:229:HOH:O	2.63	0.46
1:L:115:SER:HA	1:L:135:PHE:O	2.15	0.45
1:L:79:GLN:NE2	1:L:79:GLN:HA	2.31	0.45
1:L:183:LYS:HG2	1:L:187:GLU:OE2	2.16	0.45
2:H:17:SER:HB2	2:H:84:ASN:HA	1.97	0.45
1:L:83:ASP:CG	4:L:241:HOH:O	2.55	0.45
1:L:4:MET:CE	1:L:25:ALA:HB2	2.47	0.45
1:L:2:ILE:HG12	1:L:27:GLU:HG2	1.99	0.44
1:L:13:MET:HG3	1:L:19:VAL:CG1	2.40	0.44
1:L:35:TRP:CE2	1:L:73:LEU:HB2	2.52	0.44
2:H:14:PRO:O	2:H:15:SER:CB	2.66	0.44
2:H:45:SER:HA	4:H:250:HOH:O	2.16	0.44
2:H:86:VAL:HA	4:H:229:HOH:O	2.17	0.44
2:H:51:TYR:CD1	2:H:51:TYR:C	2.91	0.44
2:H:198:CYS:SG	2:H:211:LYS:HB3	2.58	0.44
2:H:6:GLN:HG3	2:H:96:CYS:SG	2.57	0.43
2:H:193:SER:N	4:H:227:HOH:O	2.27	0.43
1:L:94:SER:CB	1:L:95:PRO:CD	2.77	0.43
2:H:171:ALA:HB1	2:H:178:TYR:HD2	1.84	0.43
1:L:55:TYR:CG	1:L:56:THR:N	2.86	0.43
1:L:144:ILE:CG2	1:L:144:ILE:O	2.66	0.43
1:L:207:LYS:O	1:L:208:SER:HB2	2.18	0.42
1:L:190:ASN:ND2	1:L:212:ASN:HB3	2.27	0.42
1:L:132:VAL:HG12	1:L:148:TRP:CH2	2.54	0.42
1:L:19:VAL:HG13	1:L:78:VAL:CG2	2.50	0.42
2:H:24:VAL:HG12	2:H:77:ASN:O	2.19	0.42
2:H:18:LEU:HA	4:H:258:HOH:O	2.19	0.42
1:L:115:SER:HB3	1:L:136:LEU:CD1	2.49	0.42
2:H:67:ARG:NH2	2:H:90:ASP:CG	2.70	0.42
2:H:168:THR:CG2	2:H:169:VAL:H	2.31	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:127:LEU:HD21	2:H:144:LEU:HB2	2.02	0.42
1:L:79:GLN:C	1:L:81:ASP:N	2.73	0.42
1:L:48:ILE:HG22	1:L:49:TYR:N	2.35	0.42
2:H:155:VAL:HG11	2:H:182:SER:HB2	2.01	0.42
2:H:190:THR:HG22	2:H:190:THR:O	2.19	0.41
2:H:67:ARG:NH2	2:H:90:ASP:OD2	2.53	0.41
1:L:4:MET:HE1	1:L:25:ALA:HB2	2.02	0.41
1:L:19:VAL:CG2	1:L:75:ILE:HD12	2.49	0.41
1:L:22:ASN:HD21	1:L:24:LYS:HE2	1.86	0.41
2:H:31:SER:O	2:H:32:ASP:HB2	2.20	0.41
1:L:87:TYR:CD1	4:L:222:HOH:O	2.25	0.41
2:H:153:VAL:HA	2:H:201:ALA:O	2.21	0.41
1:L:96:LEU:HD12	1:L:96:LEU:H	1.85	0.41
1:L:115:SER:OG	1:L:207:LYS:HG2	2.20	0.41
1:L:13:MET:O	1:L:107:LYS:N	2.52	0.41
1:L:4:MET:HE2	1:L:25:ALA:CB	2.51	0.41
1:L:183:LYS:HB3	1:L:183:LYS:HZ3	1.84	0.41
1:L:207:LYS:HD3	1:L:207:LYS:HA	1.88	0.41
1:L:79:GLN:O	1:L:106:LEU:HD21	2.20	0.41
2:H:67:ARG:NE	4:H:229:HOH:O	2.52	0.41
2:H:32:ASP:HA	2:H:54:TYR:CD2	2.56	0.41
1:L:185:LYS:HE3	1:L:185:LYS:HB3	1.18	0.40
2:H:64:LEU:O	2:H:68:ILE:HG22	2.20	0.40
2:H:174:GLN:O	2:H:175:SER:OG	2.28	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	L	212/214 (99%)	187 (88%)	16 (8%)	9 (4%)	<b>3</b> <b>4</b>

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	H	215/217 (99%)	186 (86%)	16 (7%)	13 (6%)	2	2
3	P	5/15 (33%)	3 (60%)	2 (40%)	0	100	100
All	All	432/446 (97%)	376 (87%)	34 (8%)	22 (5%)	2	2

All (22) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	L	7	SER
1	L	166	GLN
2	H	152	PRO
2	H	193	SER
2	H	150	PRO
1	L	68	ALA
1	L	69	THR
2	H	137	SER
1	L	94	SER
1	L	138	SER
2	H	133	THR
2	H	194	ALA
2	H	215	PRO
1	L	84	ALA
2	H	135	ALA
2	H	176	GLY
2	H	203	PRO
1	L	95	PRO
2	H	149	PHE
2	H	129	PRO
2	H	153	VAL
1	L	144	ILE

### 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	L	183/183 (100%)	173 (94%)	10 (6%)	27	48

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	H	179/180 (99%)	160 (89%)	19 (11%)	8	16
3	P	5/12 (42%)	5 (100%)	0	100	100
All	All	367/375 (98%)	338 (92%)	29 (8%)	15	28

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	4	MET
1	L	69	THR
1	L	73	LEU
1	L	106	LEU
1	L	155	ARG
1	L	159	THR
1	L	165	ASP
1	L	175	MET
1	L	185	LYS
1	L	205	VAL
2	H	6	GLN
2	H	18	LEU
2	H	59	SER
2	H	83	LEU
2	H	87	THR
2	H	89	ASP
2	H	101	THR
2	H	120	THR
2	H	151	GLU
2	H	152	PRO
2	H	159	GLU
2	H	180	LEU
2	H	191	TRP
2	H	192	PRO
2	H	198	CYS
2	H	203	PRO
2	H	204	LYS
2	H	205	SER
2	H	210	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	22	ASN
1	L	28	ASN
1	L	38	GLN
1	L	79	GLN
1	L	124	GLN
1	L	137	ASN
1	L	166	GLN
1	L	190	ASN
1	L	198	HIS
1	L	210	ASN
1	L	212	ASN
2	H	44	GLN
2	H	82	GLN
2	H	111	ASN
2	H	174	GLN
3	P	2	GLN

### 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 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 ⓘ

### 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.