



Full wwPDB NMR Structure Validation Report ⓘ

Mar 9, 2026 – 12:48 AM UTC

PDB ID : 2DVJ / pdb_00002dvj
Title : phosphorylated Crk-II
Authors : Kobashigawa, Y.; Inagaki, F.
Deposited on : 2006-07-31

This is a Full wwPDB NMR 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/NMRValidationReportHelp>

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)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
wwPDB-RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
wwPDB-ShiftChecker	:	v1.2
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

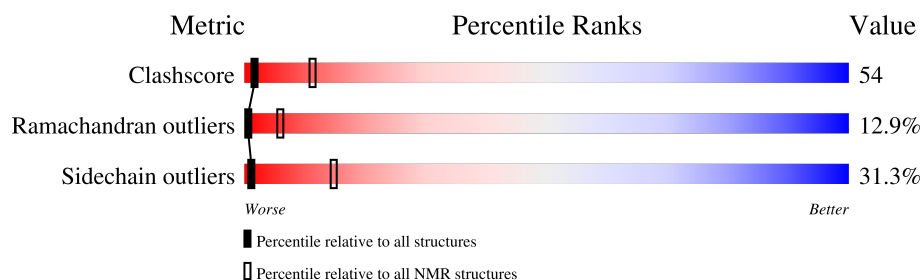
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

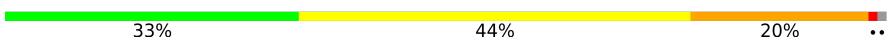
The overall completeness of chemical shifts assignment was not calculated.

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)	NMR archive (#Entries)
Clashscore	229148	14424
Ramachandran outliers	224038	12848
Sidechain outliers	223484	12823

The table below summarises the geometric issues observed across the polymeric chains and their fit to the experimental data. The red, orange, yellow and green segments indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A cyan segment indicates the fraction of residues that are not part of the well-defined cores, and 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 $\leq 5\%$

Mol	Chain	Length	Quality of chain
1	A	230	

2 Ensemble composition and analysis ⓘ

This entry contains 1 models. Identification of well-defined residues and clustering analysis are not possible.

3 Entry composition [i](#)

There is only 1 type of molecule in this entry. The entry contains 3508 atoms, of which 1715 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called V-crk sarcoma virus CT10 oncogene homolog, isoform a.

Mol	Chain	Residues	Atoms								Trace
1	A	228	Total	C	H	N	O	P	S		0
			3508	1121	1715	316	353	1	2		

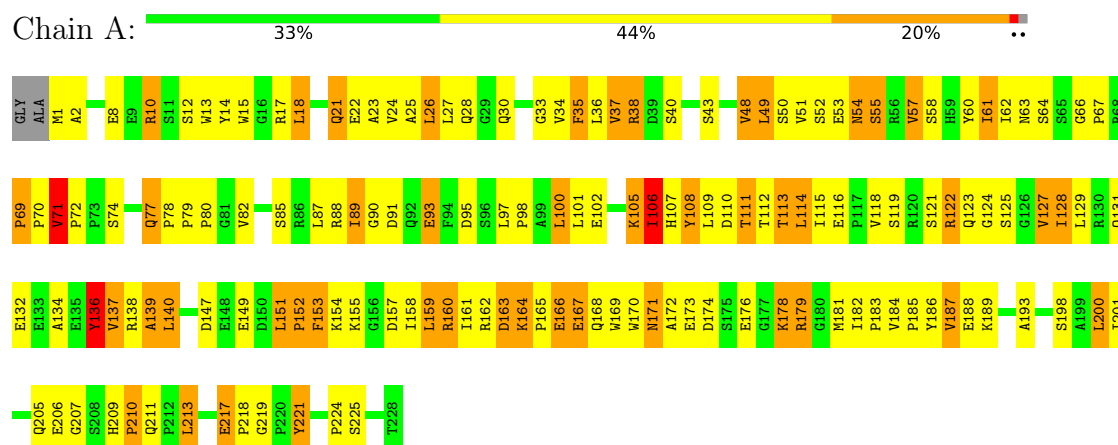
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP Q96HJ0
A	0	ALA	-	expression tag	UNP Q96HJ0

4 Residue-property plots

These plots are provided for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic is the same as shown in the summary in section 1 of this report. The second graphic shows the sequence where residues are colour-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 outliers are shown as green connectors. Residues which are classified as ill-defined in the NMR ensemble, are shown in cyan with an underline colour-coded according to the previous scheme. Residues which were present in the experimental sample, but not modelled in the final structure are shown in grey.

- Molecule 1: V-crk sarcoma virus CT10 oncogene homolog, isoform a



5 Refinement protocol and experimental data overview ⓘ

Of the 20 calculated structures, 1 were deposited, based on the following criterion: *structures with the lowest energy*.

The following table shows the software used for structure solution, optimisation and refinement.

Software name	Classification	Version
CYANA	structure solution	2.0
CNS	refinement	1.1

No chemical shift data was provided.

6 Model quality

6.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PTR

There are no covalent bond-length or bond-angle outliers.

There are no bond-length outliers.

There are no bond-angle outliers.

There are no chirality outliers.

There are no planarity outliers.

6.2 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 each 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 averaged over the ensemble.

Mol	Chain	Non-H	H(model)	H(added)	Clashes
1	A	1793	1715	1712	188
All	All	1793	1715	1712	188

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 54.

All clashes are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Clash(Å)	Distance(Å)
1:A:161:ILE:HD13	1:A:172:ALA:HB2	0.93	1.40
1:A:136:TYR:O	1:A:137:VAL:HG13	0.86	1.70
1:A:35:PHE:CE2	1:A:114:LEU:HD11	0.77	2.15
1:A:171:ASN:N	1:A:181:MET:HE1	0.77	1.94
1:A:26:LEU:HD22	1:A:27:LEU:N	0.76	1.95
1:A:23:ALA:O	1:A:36:LEU:HD21	0.75	1.80
1:A:22:GLU:O	1:A:26:LEU:HD12	0.74	1.82
1:A:13:TRP:CZ3	1:A:37:VAL:HG21	0.73	2.19
1:A:70:PRO:O	1:A:71:VAL:HG13	0.72	1.84
1:A:37:VAL:HG12	1:A:48:VAL:O	0.71	1.85
1:A:89:ILE:HG23	1:A:90:GLY:H	0.70	1.47
1:A:152:PRO:O	1:A:153:PHE:CG	0.70	2.44
1:A:139:ALA:HB1	1:A:153:PHE:CE1	0.69	2.21

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Atom-1	Atom-2	Clash(Å)	Distance(Å)
1:A:18:LEU:HD13	1:A:23:ALA:HA	0.69	1.65
1:A:87:LEU:CD1	1:A:100:LEU:HD21	0.69	2.17
1:A:169:TRP:CA	1:A:184:VAL:HG23	0.69	2.17
1:A:30:GLN:HG3	1:A:118:VAL:HG23	0.68	1.66
1:A:13:TRP:CZ3	1:A:101:LEU:HD11	0.68	2.24
1:A:24:VAL:HG23	1:A:57:VAL:HG13	0.67	1.67
1:A:15:TRP:CG	1:A:18:LEU:HD12	0.66	2.25
1:A:109:LEU:HD12	1:A:110:ASP:N	0.66	2.05
1:A:127:VAL:HG12	1:A:165:PRO:HB3	0.66	1.66
1:A:115:ILE:HG22	1:A:116:GLU:H	0.66	1.49
1:A:109:LEU:HD12	1:A:110:ASP:H	0.65	1.52
1:A:27:LEU:HD23	1:A:57:VAL:HG11	0.65	1.67
1:A:33:GLY:HA3	1:A:115:ILE:HG21	0.64	1.68
1:A:127:VAL:HG21	1:A:171:ASN:OD1	0.64	1.92
1:A:70:PRO:O	1:A:71:VAL:HG22	0.64	1.92
1:A:209:HIS:N	1:A:210:PRO:CD	0.64	2.60
1:A:26:LEU:HD22	1:A:26:LEU:C	0.63	2.19
1:A:82:VAL:O	1:A:82:VAL:HG23	0.63	1.94
1:A:106:ILE:HD13	1:A:106:ILE:N	0.63	2.08
1:A:137:VAL:HG12	1:A:188:GLU:O	0.62	1.94
1:A:169:TRP:HA	1:A:184:VAL:HG23	0.62	1.71
1:A:171:ASN:N	1:A:181:MET:CE	0.61	2.63
1:A:70:PRO:C	1:A:71:VAL:HG13	0.61	2.20
1:A:161:ILE:HD13	1:A:172:ALA:CB	0.61	2.24
1:A:26:LEU:CD2	1:A:118:VAL:HG21	0.61	2.26
1:A:137:VAL:O	1:A:158:ILE:HG23	0.61	1.96
1:A:71:VAL:HG22	1:A:72:PRO:HD3	0.60	1.73
1:A:34:VAL:CG1	1:A:118:VAL:HB	0.58	2.29
1:A:51:VAL:HG23	1:A:51:VAL:O	0.57	1.98
1:A:151:LEU:HB3	1:A:182:ILE:HG22	0.57	1.76
1:A:137:VAL:O	1:A:158:ILE:HA	0.57	1.99
1:A:162:ARG:N	1:A:171:ASN:O	0.57	2.38
1:A:110:ASP:O	1:A:111:THR:CB	0.56	2.53
1:A:181:MET:HE3	1:A:181:MET:HA	0.56	1.75
1:A:36:LEU:HD12	1:A:37:VAL:N	0.56	2.16
1:A:15:TRP:CD1	1:A:36:LEU:HD22	0.56	2.35
1:A:36:LEU:HD12	1:A:36:LEU:C	0.56	2.25
1:A:13:TRP:CZ3	1:A:37:VAL:CG2	0.56	2.88
1:A:26:LEU:HD13	1:A:27:LEU:H	0.55	1.62
1:A:110:ASP:O	1:A:111:THR:HB	0.55	2.01
1:A:109:LEU:HD13	1:A:224:PRO:HB3	0.54	1.79
1:A:78:PRO:N	1:A:79:PRO:HD2	0.54	2.18

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Atom-1	Atom-2	Clash(Å)	Distance(Å)
1:A:35:PHE:CD2	1:A:114:LEU:HD11	0.54	2.38
1:A:36:LEU:O	1:A:49:LEU:HD22	0.54	2.03
1:A:139:ALA:CB	1:A:153:PHE:CE1	0.54	2.90
1:A:159:LEU:HD13	1:A:159:LEU:N	0.54	2.18
1:A:170:TRP:O	1:A:182:ILE:HG12	0.53	2.02
1:A:169:TRP:C	1:A:184:VAL:HG23	0.53	2.29
1:A:49:LEU:HD13	1:A:49:LEU:C	0.53	2.28
1:A:49:LEU:CD1	1:A:51:VAL:HG13	0.53	2.33
1:A:49:LEU:HD11	1:A:51:VAL:CG1	0.53	2.33
1:A:137:VAL:HA	1:A:188:GLU:O	0.53	2.03
1:A:138:ARG:CD	1:A:158:ILE:HD11	0.52	2.34
1:A:77:GLN:HB2	1:A:78:PRO:HD3	0.52	1.82
1:A:105:LYS:HA	1:A:114:LEU:HB3	0.52	1.81
1:A:136:TYR:N	1:A:136:TYR:CD1	0.52	2.76
1:A:70:PRO:C	1:A:72:PRO:CD	0.52	2.83
1:A:140:LEU:HG	1:A:187:VAL:HA	0.52	1.82
1:A:60:TYR:CD1	1:A:224:PRO:HD3	0.51	2.40
1:A:128:ILE:HD11	1:A:163:ASP:OD2	0.51	2.05
1:A:138:ARG:O	1:A:140:LEU:CD2	0.51	2.59
1:A:153:PHE:C	1:A:153:PHE:CD1	0.51	2.89
1:A:54:ASN:O	1:A:55:SER:C	0.50	2.53
1:A:27:LEU:HB2	1:A:36:LEU:HD23	0.50	1.83
1:A:140:LEU:HD12	1:A:186:TYR:O	0.50	2.05
1:A:100:LEU:HD12	1:A:101:LEU:HD23	0.50	1.81
1:A:170:TRP:O	1:A:171:ASN:C	0.50	2.54
1:A:38:ARG:HB2	1:A:48:VAL:HG13	0.50	1.84
1:A:138:ARG:HD3	1:A:158:ILE:HD11	0.50	1.82
1:A:159:LEU:HA	1:A:173:GLU:O	0.50	2.07
1:A:70:PRO:O	1:A:71:VAL:CG1	0.49	2.58
1:A:106:ILE:N	1:A:106:ILE:CD1	0.49	2.74
1:A:30:GLN:HG3	1:A:118:VAL:CG2	0.49	2.36
1:A:221:PTR:N	1:A:221:PTR:HD1	0.49	2.23
1:A:27:LEU:CD1	1:A:34:VAL:HG12	0.49	2.38
1:A:118:VAL:HG13	1:A:119:SER:N	0.49	2.23
1:A:170:TRP:CG	1:A:184:VAL:HG22	0.49	2.43
1:A:27:LEU:O	1:A:30:GLN:HG2	0.49	2.07
1:A:111:THR:HG23	1:A:111:THR:O	0.49	2.06
1:A:136:TYR:O	1:A:137:VAL:CG1	0.49	2.54
1:A:127:VAL:HG12	1:A:165:PRO:CB	0.49	2.37
1:A:138:ARG:HA	1:A:158:ILE:HD13	0.49	1.84
1:A:183:PRO:HB2	1:A:186:TYR:CD1	0.48	2.43
1:A:50:SER:OG	1:A:57:VAL:HG21	0.48	2.08

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Atom-1	Atom-2	Clash(Å)	Distance(Å)
1:A:33:GLY:CA	1:A:115:ILE:HG21	0.48	2.36
1:A:10:ARG:N	1:A:10:ARG:CD	0.48	2.76
1:A:158:ILE:C	1:A:159:LEU:HD13	0.48	2.34
1:A:87:LEU:HD11	1:A:97:LEU:HD22	0.47	1.84
1:A:49:LEU:CB	1:A:62:ILE:HD11	0.47	2.39
1:A:36:LEU:CD1	1:A:37:VAL:N	0.47	2.77
1:A:49:LEU:HD11	1:A:51:VAL:HG13	0.47	1.85
1:A:169:TRP:CG	1:A:181:MET:HE2	0.47	2.44
1:A:14:TYR:HA	1:A:37:VAL:HG23	0.47	1.86
1:A:28:GLN:O	1:A:28:GLN:CG	0.47	2.62
1:A:87:LEU:C	1:A:87:LEU:HD13	0.47	2.33
1:A:87:LEU:HD12	1:A:100:LEU:HD21	0.47	1.86
1:A:187:VAL:HG23	1:A:188:GLU:N	0.47	2.24
1:A:139:ALA:O	1:A:140:LEU:HD23	0.47	2.10
1:A:161:ILE:HA	1:A:171:ASN:O	0.47	2.10
1:A:121:SER:OG	1:A:122:ARG:N	0.46	2.48
1:A:151:LEU:CD2	1:A:178:LYS:CD	0.46	2.94
1:A:110:ASP:O	1:A:111:THR:HG22	0.46	2.11
1:A:165:PRO:HG2	1:A:169:TRP:CD1	0.46	2.46
1:A:152:PRO:O	1:A:153:PHE:CD2	0.46	2.68
1:A:164:LYS:HG3	1:A:170:TRP:CE2	0.46	2.46
1:A:181:MET:HE3	1:A:181:MET:CA	0.46	2.41
1:A:173:GLU:HA	1:A:179:ARG:HA	0.46	1.88
1:A:60:TYR:OH	1:A:110:ASP:HB3	0.46	2.11
1:A:70:PRO:O	1:A:71:VAL:CG2	0.46	2.64
1:A:160:ARG:CD	1:A:160:ARG:C	0.46	2.86
1:A:166:GLU:CG	1:A:169:TRP:CD1	0.46	2.99
1:A:170:TRP:CD1	1:A:184:VAL:HG21	0.45	2.46
1:A:166:GLU:HG2	1:A:169:TRP:CD1	0.45	2.47
1:A:217:GLU:N	1:A:218:PRO:HD3	0.45	2.26
1:A:140:LEU:HD21	1:A:188:GLU:CD	0.45	2.36
1:A:165:PRO:HD2	1:A:170:TRP:HA	0.45	1.88
1:A:170:TRP:CD1	1:A:184:VAL:CG2	0.45	2.99
1:A:151:LEU:HD23	1:A:179:ARG:C	0.44	2.38
1:A:1:MET:O	1:A:2:ALA:HB3	0.44	2.13
1:A:21:GLN:O	1:A:25:ALA:HB2	0.44	2.12
1:A:35:PHE:CD1	1:A:35:PHE:C	0.44	2.94
1:A:209:HIS:CG	1:A:210:PRO:HD3	0.44	2.48
1:A:62:ILE:CG2	1:A:63:ASN:N	0.44	2.81
1:A:89:ILE:HG23	1:A:90:GLY:N	0.44	2.23
1:A:217:GLU:O	1:A:219:GLY:N	0.44	2.51
1:A:49:LEU:HB2	1:A:62:ILE:HD11	0.43	1.88

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Atom-1	Atom-2	Clash(Å)	Distance(Å)
1:A:70:PRO:O	1:A:72:PRO:HD3	0.43	2.13
1:A:26:LEU:C	1:A:26:LEU:CD2	0.43	2.91
1:A:60:TYR:CG	1:A:224:PRO:HD3	0.43	2.48
1:A:110:ASP:O	1:A:111:THR:CG2	0.43	2.67
1:A:161:ILE:HG21	1:A:164:LYS:CE	0.43	2.44
1:A:27:LEU:HD23	1:A:50:SER:OG	0.43	2.13
1:A:79:PRO:CB	1:A:80:PRO:CD	0.43	2.96
1:A:91:ASP:CB	1:A:225:SER:CB	0.43	2.97
1:A:13:TRP:CE3	1:A:37:VAL:CG2	0.43	3.02
1:A:15:TRP:CD1	1:A:18:LEU:HD12	0.43	2.49
1:A:151:LEU:HD11	1:A:174:ASP:HB2	0.42	1.91
1:A:166:GLU:CG	1:A:169:TRP:NE1	0.42	2.82
1:A:184:VAL:N	1:A:185:PRO:HD2	0.42	2.29
1:A:209:HIS:N	1:A:210:PRO:HD2	0.42	2.29
1:A:82:VAL:O	1:A:82:VAL:CG2	0.42	2.65
1:A:88:ARG:CB	1:A:93:GLU:HB3	0.42	2.45
1:A:106:ILE:HD13	1:A:106:ILE:H	0.42	1.73
1:A:152:PRO:O	1:A:153:PHE:CD1	0.42	2.72
1:A:172:ALA:HB3	1:A:182:ILE:CG2	0.42	2.44
1:A:27:LEU:CD2	1:A:50:SER:OG	0.42	2.67
1:A:134:ALA:C	1:A:160:ARG:CD	0.42	2.92
1:A:78:PRO:CD	1:A:79:PRO:HD2	0.42	2.44
1:A:127:VAL:HG11	1:A:171:ASN:HB2	0.42	1.92
1:A:33:GLY:C	1:A:34:VAL:HG23	0.42	2.39
1:A:33:GLY:C	1:A:34:VAL:CG2	0.42	2.93
1:A:62:ILE:HG22	1:A:63:ASN:N	0.42	2.30
1:A:132:GLU:O	1:A:200:LEU:HD11	0.42	2.15
1:A:35:PHE:CE2	1:A:101:LEU:HD13	0.41	2.49
1:A:149:GLU:O	1:A:179:ARG:O	0.41	2.37
1:A:61:ILE:HG12	1:A:221:PTR:CG	0.41	2.45
1:A:87:LEU:CD1	1:A:100:LEU:CD2	0.41	2.95
1:A:97:LEU:CB	1:A:98:PRO:CD	0.41	2.98
1:A:26:LEU:O	1:A:30:GLN:NE2	0.41	2.53
1:A:164:LYS:HE3	1:A:170:TRP:CE3	0.41	2.50
1:A:164:LYS:HD2	1:A:170:TRP:CG	0.41	2.51
1:A:10:ARG:N	1:A:10:ARG:HD2	0.41	2.30
1:A:87:LEU:HD12	1:A:100:LEU:CD2	0.41	2.45
1:A:168:GLN:O	1:A:185:PRO:HD2	0.41	2.15
1:A:61:ILE:HG12	1:A:221:PTR:CD1	0.41	2.45
1:A:36:LEU:O	1:A:49:LEU:CD2	0.41	2.69
1:A:61:ILE:CB	1:A:221:PTR:HB3	0.41	2.46
1:A:88:ARG:HB2	1:A:93:GLU:HB3	0.41	1.92

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Atom-1	Atom-2	Clash(Å)	Distance(Å)
1:A:71:VAL:CG2	1:A:72:PRO:HD3	0.41	2.44
1:A:167:GLU:O	1:A:184:VAL:HG11	0.41	2.16
1:A:69:PRO:CB	1:A:70:PRO:CD	0.41	2.99
1:A:138:ARG:O	1:A:139:ALA:O	0.40	2.39
1:A:34:VAL:HG12	1:A:118:VAL:HB	0.40	1.92
1:A:173:GLU:HG2	1:A:174:ASP:N	0.40	2.32
1:A:151:LEU:HD22	1:A:151:LEU:HA	0.40	1.81

6.3 Torsion angles [i](#)

6.3.1 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 PDB entries followed by that with respect to all NMR entries. 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	225/230 (98%)	139 (62%)	57 (25%)	29 (13%)	0	6
All	All	225/230 (98%)	139 (62%)	57 (25%)	29 (13%)	0	6

All 29 Ramachandran outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type
1	A	40	SER
1	A	52	SER
1	A	57	VAL
1	A	66	GLY
1	A	67	PRO
1	A	69	PRO
1	A	71	VAL
1	A	74	SER
1	A	77	GLN
1	A	89	ILE
1	A	106	ILE
1	A	108	TYR
1	A	111	THR
1	A	112	THR
1	A	113	THR
1	A	122	ARG

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Mol	Chain	Res	Type
1	A	124	GLY
1	A	125	SER
1	A	127	VAL
1	A	136	TYR
1	A	137	VAL
1	A	139	ALA
1	A	152	PRO
1	A	153	PHE
1	A	193	ALA
1	A	207	GLY
1	A	210	PRO
1	A	211	GLN
1	A	213	LEU

6.3.2 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 PDB entries followed by that with respect to all NMR entries. 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	195/195 (100%)	134 (69%)	61 (31%)	1	15
All	All	195/195 (100%)	134 (69%)	61 (31%)	1	15

All 61 residues with a non-rotameric sidechain are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type
1	A	8	GLU
1	A	10	ARG
1	A	12	SER
1	A	17	ARG
1	A	18	LEU
1	A	21	GLN
1	A	26	LEU
1	A	35	PHE
1	A	37	VAL
1	A	38	ARG
1	A	43	SER
1	A	48	VAL

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Mol	Chain	Res	Type
1	A	49	LEU
1	A	53	GLU
1	A	54	ASN
1	A	55	SER
1	A	58	SER
1	A	61	ILE
1	A	64	SER
1	A	71	VAL
1	A	85	SER
1	A	93	GLU
1	A	95	ASP
1	A	100	LEU
1	A	102	GLU
1	A	105	LYS
1	A	106	ILE
1	A	107	HIS
1	A	108	TYR
1	A	113	THR
1	A	114	LEU
1	A	123	GLN
1	A	128	ILE
1	A	129	LEU
1	A	131	GLN
1	A	136	TYR
1	A	140	LEU
1	A	147	ASP
1	A	151	LEU
1	A	154	LYS
1	A	155	LYS
1	A	157	ASP
1	A	159	LEU
1	A	160	ARG
1	A	163	ASP
1	A	164	LYS
1	A	166	GLU
1	A	167	GLU
1	A	171	ASN
1	A	176	GLU
1	A	178	LYS
1	A	179	ARG
1	A	187	VAL
1	A	189	LYS

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Mol	Chain	Res	Type
1	A	198	SER
1	A	200	LEU
1	A	201	ILE
1	A	205	GLN
1	A	206	GLU
1	A	213	LEU
1	A	217	GLU

6.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

In the following table, the Counts columns list the number of bonds for which Mogul statistics could be retrieved, the number of bonds that are observed in the model and the number of bonds 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 is the number of standard deviations the observed value is removed from the expected value. A bond length with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the average root-mean-square of all Z scores of the bond lengths.

Mol	Type	Chain	Res	Link	Bond lengths		
					Counts	RMSZ	#Z>2
1	PTR	A	221	1	15,16,17	0.98	1 (6%)

In the following table, the Counts columns list the number of angles for which Mogul statistics could be retrieved, the number of angles that are observed in the model and the number of 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 angle is the number of standard deviations the observed value is removed from the expected value. A bond angle with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the average root-mean-square of all Z scores of the bond angles.

Mol	Type	Chain	Res	Link	Bond angles		
					Counts	RMSZ	#Z>2
1	PTR	A	221	1	17,22,24	0.53	0 (0%)

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
1	PTR	A	221	1	-	0,10,11,13	0,1,1,1

All bond outliers are listed below.

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	221	PTR	P-OH	2.26	1.63	1.59

There are no bond-angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.6 Ligand geometry [i](#)

There are no ligands in this entry.

6.7 Other polymers [i](#)

There are no such molecules in this entry.

6.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

7 Chemical shift validation

No chemical shift data were provided