



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:24 AM UTC

PDB ID : 3LLL / pdb_00003lll
Title : Crystal structure of mouse pacsin2 F-BAR domain
Authors : Rudolph, M.G.
Deposited on : 2010-01-29
Resolution : 3.30 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)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

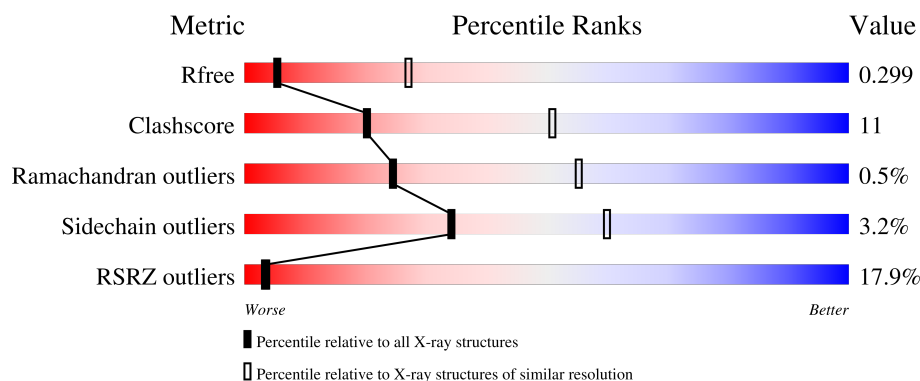
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.30 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))
R_{free}	180053	1169 (3.32-3.28)
Clashscore	190562	1209 (3.32-3.28)
Ramachandran outliers	187476	1188 (3.32-3.28)
Sidechain outliers	187428	1187 (3.32-3.28)
RSRZ outliers	180081	1169 (3.32-3.28)

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 $\leq 5\%$. 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.

Mol	Chain	Length	Quality of chain
1	A	287	
1	B	287	

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 4806 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.

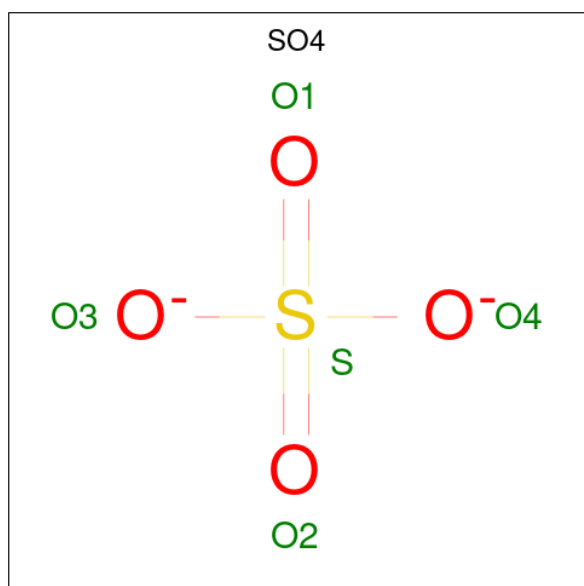
- Molecule 1 is a protein called Protein kinase C and casein kinase substrate in neurons protein 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	287	Total	C	N	O	S	0	0	0
			2386	1502	427	443	14			
1	B	287	Total	C	N	O	S	0	0	0
			2386	1502	427	443	14			

- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Ca	0	0
			1	1		
2	B	1	Total	Ca	0	0
			1	1		

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total O S 5 4 1	0	0
3	A	1	Total O S 5 4 1	0	0
3	A	1	Total O S 5 4 1	0	0
3	A	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0

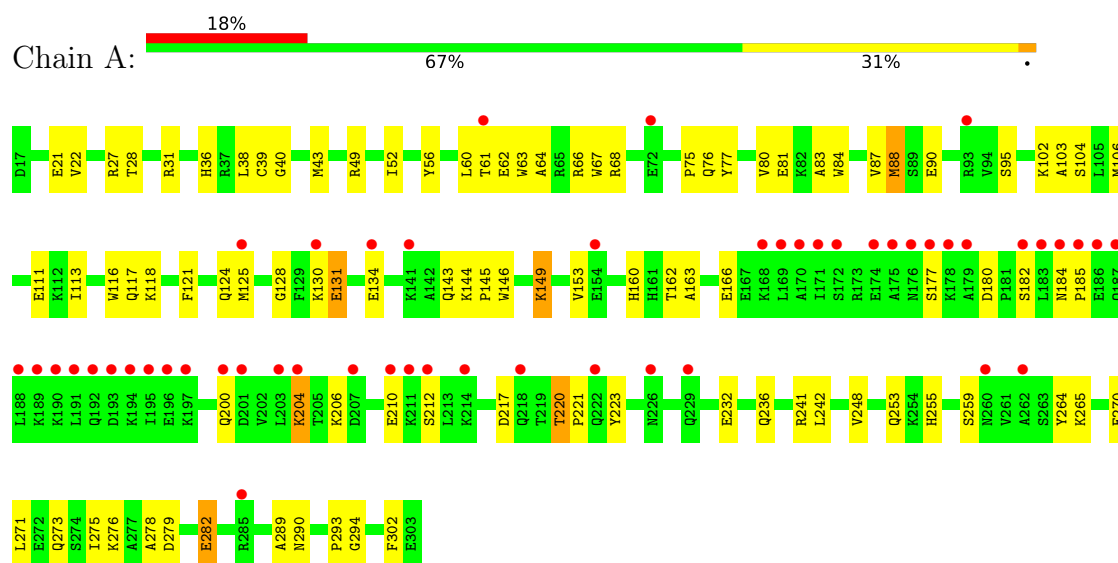
- Molecule 4 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	2	Total Cl 2 2	0	0

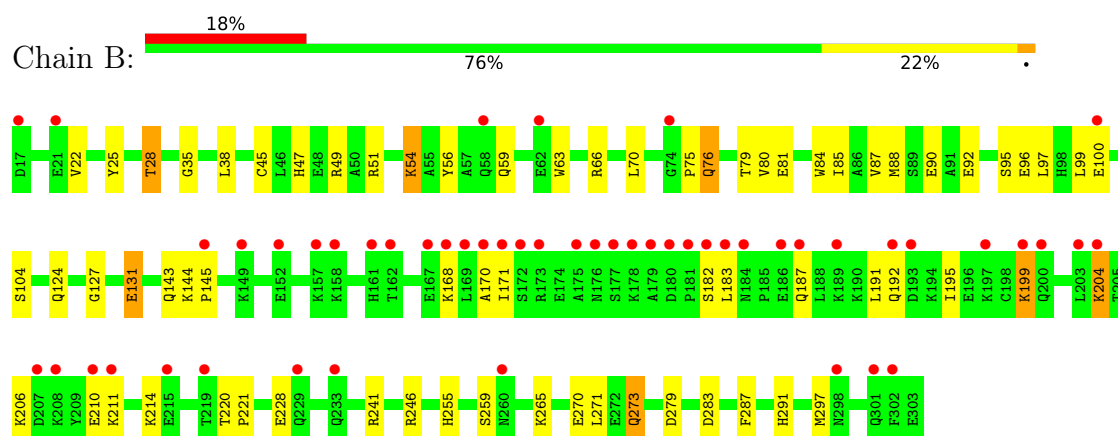
3 Residue-property plots [i](#)

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 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.

- Molecule 1: Protein kinase C and casein kinase substrate in neurons protein 2



- Molecule 1: Protein kinase C and casein kinase substrate in neurons protein 2



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	104.39Å 106.40Å 125.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	52.20 – 3.30 52.20 – 3.30	Depositor EDS
% Data completeness (in resolution range)	83.3 (52.20-3.30) 83.4 (52.20-3.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.03	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.48 (at 3.33Å)	Xtriage
Refinement program	PHENIX (phenix.refine)	Depositor
R, R_{free}	0.256 , 0.297 0.257 , 0.299	Depositor DCC
R_{free} test set	993 reflections (4.61%)	wwPDB-VP
Wilson B-factor (Å ²)	41.1	Xtriage
Anisotropy	0.330	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 95.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	0.067 for k,h,-l	Xtriage
F_o, F_c correlation	0.83	EDS
Total number of atoms	4806	wwPDB-VP
Average B, all atoms (Å ²)	98.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.24% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CA, SO4, CL

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	0.25	0/2434	0.66	0/3259
1	B	0.25	0/2434	0.65	0/3259
All	All	0.25	0/4868	0.66	0/6518

There are no bond length outliers.

There are no bond angle outliers.

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	A	2386	0	2365	72	0
1	B	2386	0	2365	55	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	20	0	0	0	0
3	B	10	0	0	0	0
4	B	2	0	0	0	0
All	All	4806	0	4730	108	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 11.

All (108) 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:153:VAL:HG22	1:A:212:SER:HB3	1.59	0.84
1:A:289:ALA:HA	1:A:294:GLY:HA3	1.71	0.71
1:A:40:GLY:HA2	1:A:43:MET:HE2	1.72	0.71
1:A:265:LYS:HG2	1:B:259:SER:HB2	1.73	0.69
1:B:35:GLY:HA2	1:B:241:ARG:HH21	1.56	0.69
1:A:38:LEU:HD21	1:B:75:PRO:HD2	1.76	0.68
1:A:242:LEU:HD22	1:B:80:VAL:HG21	1.77	0.66
1:B:144:LYS:HB3	1:B:145:PRO:HD3	1.78	0.65
1:A:259:SER:HB2	1:B:265:LYS:HG2	1.78	0.64
1:B:87:VAL:O	1:B:90:GLU:HG3	1.98	0.64
1:A:184:ASN:HB3	1:A:185:PRO:HD2	1.79	0.63
1:A:289:ALA:C	1:A:290:ASN:HD22	2.08	0.62
1:A:62:GLU:O	1:A:66:ARG:HB2	1.99	0.62
1:B:104:SER:HB3	1:B:255:HIS:CD2	2.35	0.61
1:A:52:ILE:HD13	1:B:59:GLN:HB3	1.83	0.60
1:A:206:LYS:O	1:A:210:GLU:HG2	2.03	0.59
1:A:88:MET:N	1:A:88:MET:HE3	2.18	0.58
1:B:35:GLY:HA2	1:B:241:ARG:NH2	2.20	0.57
1:B:170:ALA:HB1	1:B:195:ILE:HG22	1.87	0.57
1:B:211:LYS:HA	1:B:214:LYS:HE2	1.87	0.56
1:B:88:MET:N	1:B:88:MET:HE3	2.20	0.56
1:A:149:LYS:HA	1:A:149:LYS:HE2	1.87	0.55
1:A:160:HIS:NE2	1:A:206:LYS:HB2	2.22	0.55
1:A:220:THR:H	1:A:221:PRO:CD	2.20	0.55
1:B:95:SER:O	1:B:99:LEU:HB2	2.07	0.54
1:A:217:ASP:O	1:A:220:THR:HG22	2.08	0.54
1:A:104:SER:HB3	1:A:255:HIS:CD2	2.43	0.53
1:B:79:THR:HG21	1:B:279:ASP:N	2.24	0.53
1:B:79:THR:HG22	1:B:283:ASP:OD1	2.08	0.53
1:B:92:GLU:O	1:B:96:GLU:HG3	2.09	0.53
1:A:113:ILE:HD11	1:A:248:VAL:HG11	1.91	0.53
1:A:270:GLU:O	1:A:273:GLN:HG2	2.10	0.52
1:A:220:THR:HA	1:A:223:TYR:HB3	1.90	0.52
1:A:102:LYS:HD3	1:A:103:ALA:N	2.24	0.52
1:B:38:LEU:HD12	1:B:241:ARG:NH2	2.24	0.52
1:A:22:VAL:HA	1:A:143:GLN:NE2	2.25	0.51
1:B:204:LYS:HB2	1:B:204:LYS:NZ	2.24	0.51
1:B:187:GLN:HE21	1:B:191:LEU:HD11	1.76	0.50
1:B:270:GLU:HA	1:B:273:GLN:HG3	1.94	0.50
1:A:38:LEU:HD13	1:B:76:GLN:HG2	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:25:TYR:O	1:B:28:THR:HG22	2.13	0.49
1:B:47:HIS:CE1	1:B:51:ARG:HH11	2.30	0.49
1:A:279:ASP:CG	1:A:282:GLU:HB2	2.38	0.49
1:A:220:THR:N	1:A:221:PRO:CD	2.76	0.48
1:A:83:ALA:HB2	1:A:278:ALA:HB2	1.95	0.48
1:B:38:LEU:HD12	1:B:241:ARG:HH22	1.79	0.48
1:A:144:LYS:HB3	1:A:145:PRO:HD3	1.96	0.48
1:A:184:ASN:HB3	1:A:185:PRO:CD	2.43	0.48
1:A:264:TYR:CD2	1:B:259:SER:HB3	2.48	0.48
1:A:146:TRP:HZ3	1:A:220:THR:HB	1.78	0.48
1:B:22:VAL:HA	1:B:143:GLN:NE2	2.28	0.48
1:A:177:SER:HA	1:A:180:ASP:OD1	2.13	0.48
1:A:279:ASP:HB3	1:A:282:GLU:HB2	1.96	0.48
1:A:220:THR:CG2	1:A:221:PRO:HD3	2.44	0.47
1:B:220:THR:HB	1:B:221:PRO:HD3	1.94	0.47
1:A:276:LYS:HA	1:B:246:ARG:NH1	2.29	0.47
1:A:56:TYR:CD1	1:B:56:TYR:HB2	2.49	0.47
1:A:27:ARG:HG2	1:B:291:HIS:ND1	2.30	0.47
1:A:87:VAL:O	1:A:90:GLU:HG3	2.14	0.47
1:B:192:GLN:HA	1:B:195:ILE:HG12	1.96	0.47
1:A:220:THR:H	1:A:221:PRO:HD3	1.79	0.47
1:A:200:GLN:O	1:A:204:LYS:HG2	2.16	0.46
1:A:204:LYS:HB2	1:A:204:LYS:NZ	2.31	0.46
1:A:39:CYS:O	1:A:43:MET:HG3	2.16	0.46
1:A:220:THR:HG22	1:A:221:PRO:HD3	1.98	0.46
1:A:293:PRO:HG2	1:B:228:GLU:HG2	1.97	0.46
1:B:84:TRP:C	1:B:84:TRP:CD1	2.94	0.46
1:B:124:GLN:HB3	1:B:127:GLY:O	2.16	0.46
1:A:232:GLU:O	1:A:236:GLN:HG2	2.16	0.45
1:A:131:GLU:HA	1:A:134:GLU:HB2	1.99	0.44
1:A:63:TRP:CE2	1:B:49:ARG:HB2	2.53	0.44
1:B:168:LYS:HD3	1:B:171:ILE:HD12	1.99	0.43
1:A:271:LEU:O	1:A:275:ILE:HG13	2.18	0.43
1:A:56:TYR:O	1:A:60:LEU:HG	2.19	0.43
1:A:61:THR:HG23	1:A:95:SER:OG	2.18	0.43
1:A:39:CYS:HB2	1:A:117:GLN:HG3	2.01	0.43
1:B:97:LEU:O	1:B:100:GLU:HB3	2.18	0.43
1:A:302:PHE:HZ	1:B:206:LYS:HE2	1.84	0.43
1:B:183:LEU:O	1:B:187:GLN:HB2	2.19	0.43
1:A:88:MET:HE3	1:A:88:MET:H	1.84	0.43
1:A:290:ASN:HD22	1:A:290:ASN:N	2.17	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:270:GLU:O	1:B:273:GLN:HB2	2.18	0.42
1:A:77:TYR:HA	1:A:81:GLU:OE2	2.18	0.42
1:A:67:TRP:CE2	1:B:45:CYS:HA	2.55	0.42
1:B:47:HIS:HE1	1:B:51:ARG:HH11	1.66	0.42
1:A:49:ARG:HD3	1:B:63:TRP:CH2	2.55	0.42
1:B:79:THR:HG22	1:B:283:ASP:CG	2.45	0.42
1:B:66:ARG:O	1:B:70:LEU:HG	2.20	0.42
1:A:125:MET:SD	1:A:125:MET:N	2.93	0.42
1:B:54:LYS:NZ	1:B:54:LYS:HB3	2.35	0.42
1:A:31:ARG:NH2	1:B:287:PHE:CD1	2.88	0.42
1:A:124:GLN:HG3	1:A:128:GLY:O	2.20	0.42
1:A:64:ALA:O	1:A:68:ARG:HB2	2.20	0.41
1:A:163:ALA:HA	1:A:166:GLU:HB3	2.01	0.41
1:A:84:TRP:C	1:A:84:TRP:CD1	2.97	0.41
1:A:130:LYS:O	1:A:134:GLU:HB2	2.20	0.41
1:A:278:ALA:HB3	1:B:246:ARG:HG3	2.01	0.41
1:B:81:GLU:O	1:B:85:ILE:HG13	2.21	0.41
1:A:36:HIS:HB2	1:A:121:PHE:CE1	2.56	0.41
1:A:76:GLN:NE2	1:A:80:VAL:HB	2.34	0.41
1:A:160:HIS:C	1:A:162:THR:H	2.28	0.41
1:B:131:GLU:CD	1:B:131:GLU:H	2.27	0.41
1:B:206:LYS:O	1:B:210:GLU:HG2	2.20	0.41
1:A:75:PRO:HD2	1:B:38:LEU:HD21	2.01	0.41
1:A:106:MET:O	1:A:111:GLU:HG2	2.20	0.41
1:B:199:LYS:HA	1:B:199:LYS:HE3	2.02	0.41
1:A:116:TRP:CH2	1:A:241:ARG:HG3	2.56	0.41
1:A:21:GLU:HG2	1:B:297:MET:HE1	2.03	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	285/287 (99%)	268 (94%)	15 (5%)	2 (1%)	18	49
1	B	285/287 (99%)	267 (94%)	17 (6%)	1 (0%)	30	60
All	All	570/574 (99%)	535 (94%)	32 (6%)	3 (0%)	24	55

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	182	SER
1	A	182	SER
1	A	220	THR

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	252/252 (100%)	244 (97%)	8 (3%)	34	60
1	B	252/252 (100%)	244 (97%)	8 (3%)	34	60
All	All	504/504 (100%)	488 (97%)	16 (3%)	34	60

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	THR
1	A	88	MET
1	A	118	LYS
1	A	131	GLU
1	A	149	LYS
1	A	204	LYS
1	A	253	GLN
1	A	282	GLU
1	B	28	THR
1	B	54	LYS
1	B	76	GLN
1	B	131	GLU
1	B	199	LYS

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Mol	Chain	Res	Type
1	B	204	LYS
1	B	271	LEU
1	B	273	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	36	HIS
1	A	47	HIS
1	A	58	GLN
1	A	124	GLN
1	A	187	GLN
1	A	192	GLN
1	A	222	GLN
1	A	229	GLN
1	A	253	GLN
1	A	255	HIS
1	A	273	GLN
1	A	290	ASN
1	A	298	ASN
1	B	24	ASN
1	B	36	HIS
1	B	44	ASN
1	B	47	HIS
1	B	69	GLN
1	B	117	GLN
1	B	143	GLN
1	B	176	ASN
1	B	187	GLN
1	B	218	GLN
1	B	222	GLN
1	B	235	GLN
1	B	253	GLN
1	B	273	GLN
1	B	290	ASN

5.3.3 RNA ⓘ

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 10 ligands modelled in this entry, 4 are monoatomic - leaving 6 for Mogul analysis.

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	$\# Z > 2$
3	SO4	A	2	-	4,4,4	0.24	0	6,6,6	0.05	0
3	SO4	B	3	-	4,4,4	0.24	0	6,6,6	0.10	0
3	SO4	A	7	-	4,4,4	0.24	0	6,6,6	0.07	0
3	SO4	B	5	-	4,4,4	0.24	0	6,6,6	0.07	0
3	SO4	A	1	-	4,4,4	0.23	0	6,6,6	0.07	0
3	SO4	A	4	-	4,4,4	0.24	0	6,6,6	0.07	0

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.

5.7 Other polymers [i](#)

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

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	287/287 (100%)	1.04	51 (17%)  	42, 89, 182, 306	0
1	B	287/287 (100%)	1.01	52 (18%)  	40, 85, 165, 278	0
All	All	574/574 (100%)	1.02	103 (17%)  	40, 87, 178, 306	0

All (103) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	195	ILE	6.6
1	A	179	ALA	6.6
1	A	169	LEU	5.4
1	B	193	ASP	5.2
1	A	171	ILE	5.1
1	B	177	SER	4.5
1	A	170	ALA	4.3
1	B	260	ASN	4.3
1	B	180	ASP	4.1
1	A	285	ARG	4.1
1	B	158	LYS	4.0
1	B	203	LEU	3.9
1	A	184	ASN	3.9
1	B	298	ASN	3.9
1	A	201	ASP	3.9
1	B	167	GLU	3.8
1	A	182	SER	3.8
1	B	179	ALA	3.8
1	A	187	GLN	3.7
1	B	204	LYS	3.6
1	B	187	GLN	3.6
1	A	61	THR	3.5
1	A	193	ASP	3.5
1	A	197	LYS	3.5

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Mol	Chain	Res	Type	RSRZ
1	A	183	LEU	3.5
1	B	17	ASP	3.5
1	A	125	MET	3.4
1	B	170	ALA	3.4
1	B	171	ILE	3.3
1	B	182	SER	3.3
1	B	186	GLU	3.3
1	B	183	LEU	3.3
1	A	218	GLN	3.2
1	B	173	ARG	3.2
1	A	172	SER	3.2
1	B	162	THR	3.2
1	A	190	LYS	3.1
1	A	189	LYS	3.1
1	B	207	ASP	3.1
1	A	186	GLU	3.1
1	A	192	GLN	3.1
1	B	176	ASN	3.1
1	B	178	LYS	3.1
1	A	191	LEU	3.0
1	B	74	GLY	3.0
1	A	260	ASN	3.0
1	A	203	LEU	3.0
1	A	210	GLU	3.0
1	A	229	GLN	2.9
1	B	208	LYS	2.9
1	A	211	LYS	2.8
1	A	188	LEU	2.7
1	A	174	GLU	2.7
1	B	229	GLN	2.7
1	A	141	LYS	2.6
1	A	194	LYS	2.6
1	B	233	GLN	2.6
1	A	185	PRO	2.6
1	A	130	LYS	2.6
1	B	175	ALA	2.5
1	B	302	PHE	2.5
1	B	197	LYS	2.5
1	B	210	GLU	2.5
1	A	222	GLN	2.5
1	B	21	GLU	2.4
1	B	219	THR	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	181	PRO	2.4
1	A	200	GLN	2.4
1	B	192	GLN	2.4
1	B	62	GLU	2.4
1	A	93	ARG	2.4
1	B	215	GLU	2.4
1	A	175	ALA	2.3
1	B	152	GLU	2.3
1	B	168	LYS	2.3
1	B	301	GLN	2.3
1	A	72	GLU	2.3
1	B	169	LEU	2.3
1	A	214	LYS	2.3
1	B	58	GLN	2.3
1	A	168	LYS	2.2
1	B	211	LYS	2.2
1	A	176	ASN	2.2
1	A	207	ASP	2.2
1	A	226	ASN	2.2
1	A	262	ALA	2.2
1	B	100	GLU	2.2
1	A	204	LYS	2.2
1	A	178	LYS	2.1
1	B	157	LYS	2.1
1	A	134	GLU	2.1
1	A	196	GLU	2.1
1	B	145	PRO	2.1
1	B	172	SER	2.1
1	B	149	LYS	2.1
1	A	212	SER	2.1
1	B	161	HIS	2.1
1	B	200	GLN	2.1
1	B	184	ASN	2.0
1	A	154	GLU	2.0
1	B	189	LYS	2.0
1	B	199	LYS	2.0
1	A	177	SER	2.0

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	CA	A	500	1/1	0.77	0.24	104,104,104,104	0
3	SO4	A	7	5/5	0.80	0.18	151,151,151,151	0
3	SO4	A	4	5/5	0.82	0.15	131,131,131,131	0
3	SO4	B	5	5/5	0.84	0.16	145,145,145,145	0
3	SO4	A	2	5/5	0.86	0.16	103,103,103,103	0
3	SO4	B	3	5/5	0.90	0.20	85,85,85,85	0
2	CA	B	500	1/1	0.95	0.06	71,71,71,71	0
4	CL	B	1	1/1	0.95	0.06	61,61,61,61	0
4	CL	B	2	1/1	0.95	0.07	81,81,81,81	0
3	SO4	A	1	5/5	0.97	0.09	78,78,78,78	0

6.5 Other polymers [i](#)

There are no such residues in this entry.