



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 08:23 PM UTC

PDB ID : 8OQW / pdb_00008oqw
Title : Crystal structure of Tannerella forsythia MurNAc kinase MurK
Authors : Gogler, K.; Fink, P.; Stasiak, A.C.; Stehle, T.; Zocher, G.
Deposited on : 2023-04-12
Resolution : 2.05 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

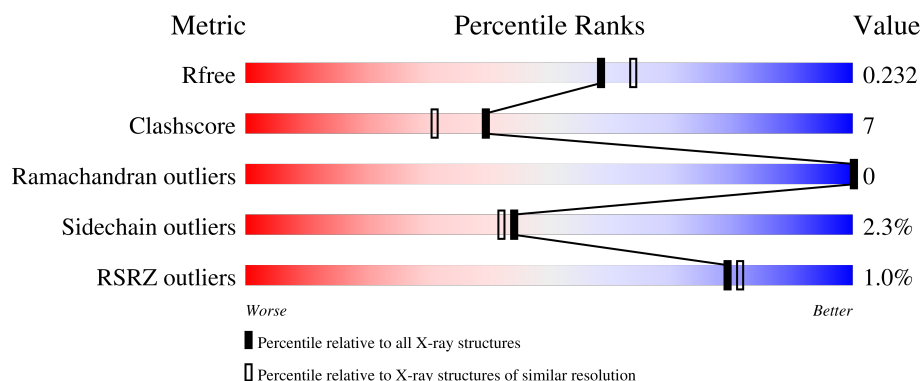
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.05 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	2260 (2.04-2.04)
Clashscore	190562	2333 (2.04-2.04)
Ramachandran outliers	187476	2318 (2.04-2.04)
Sidechain outliers	187428	2318 (2.04-2.04)
RSRZ outliers	180081	2260 (2.04-2.04)

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	283	 83% 15% ..
1	B	283	 84% 16%
1	C	283	 82% 17% .
1	D	283	 83% 14% ..

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 9079 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 ATPase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	278	Total	C	N	O	S	0	1	0
			2163	1381	369	399	14			
1	B	283	Total	C	N	O	S	0	0	0
			2186	1399	368	406	13			
1	C	283	Total	C	N	O	S	0	0	0
			2181	1396	367	405	13			
1	D	278	Total	C	N	O	S	0	0	0
			2155	1377	365	400	13			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	20	GLN	ARG	conflict	UNP G8UQH9
A	22	ARG	GLN	conflict	UNP G8UQH9
A	24	ILE	VAL	conflict	UNP G8UQH9
A	76	VAL	ILE	conflict	UNP G8UQH9
A	130	ALA	THR	conflict	UNP G8UQH9
A	282	LYS	GLU	conflict	UNP G8UQH9
B	20	GLN	ARG	conflict	UNP G8UQH9
B	22	ARG	GLN	conflict	UNP G8UQH9
B	24	ILE	VAL	conflict	UNP G8UQH9
B	76	VAL	ILE	conflict	UNP G8UQH9
B	130	ALA	THR	conflict	UNP G8UQH9
B	282	LYS	GLU	conflict	UNP G8UQH9
C	20	GLN	ARG	conflict	UNP G8UQH9
C	22	ARG	GLN	conflict	UNP G8UQH9
C	24	ILE	VAL	conflict	UNP G8UQH9
C	76	VAL	ILE	conflict	UNP G8UQH9
C	130	ALA	THR	conflict	UNP G8UQH9
C	282	LYS	GLU	conflict	UNP G8UQH9
D	20	GLN	ARG	conflict	UNP G8UQH9
D	22	ARG	GLN	conflict	UNP G8UQH9
D	24	ILE	VAL	conflict	UNP G8UQH9

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
D	76	VAL	ILE	conflict	UNP G8UQH9
D	130	ALA	THR	conflict	UNP G8UQH9
D	282	LYS	GLU	conflict	UNP G8UQH9

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	D	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	D	1	Total	C	O	0	0
			6	3	3		

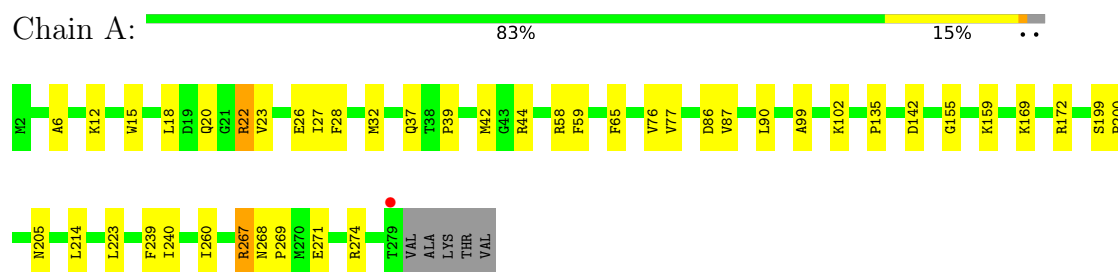
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	106	Total	O	0	0
			106	106		
4	B	102	Total	O	0	0
			102	102		
4	C	85	Total	O	0	0
			85	85		
4	D	65	Total	O	0	0
			65	65		

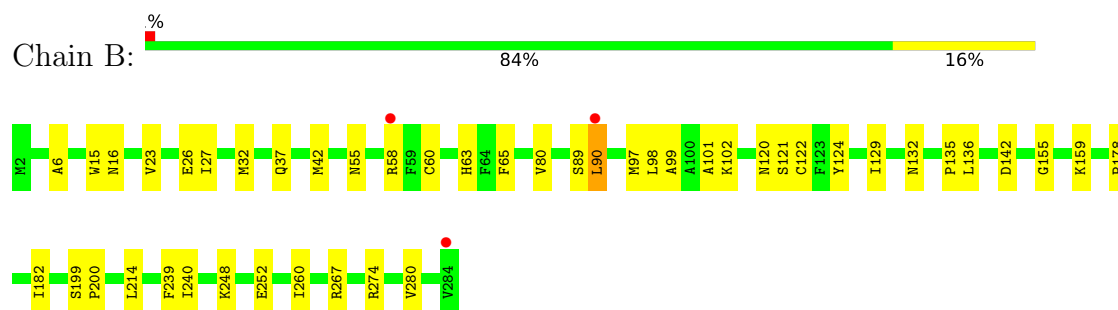
3 Residue-property plots [i](#)

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.

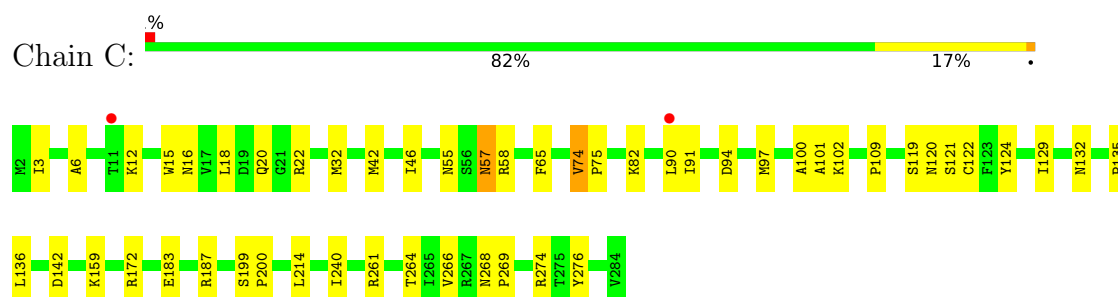
• Molecule 1: ATPase



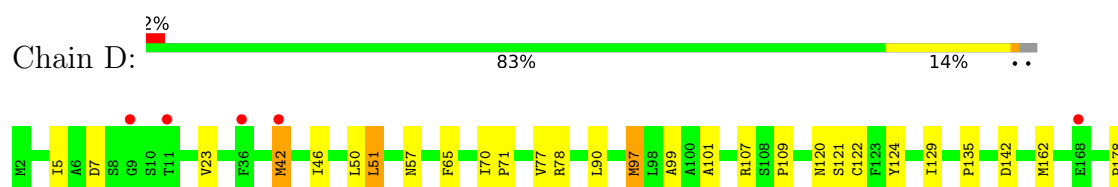
• Molecule 1: ATPase



• Molecule 1: ATPase



• Molecule 1: ATPase



I182	I198	S199	P200	A210	L214	L223	M228	Y232	I260	N268	P269	M270	R274	T279	VAL	ALA	LYS	THR	VAL
------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	-----	-----	-----	-----	-----

4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	58.76Å 59.51Å 96.71Å 74.30° 75.27° 65.95°	Depositor
Resolution (Å)	49.50 – 2.05 49.50 – 2.05	Depositor EDS
% Data completeness (in resolution range)	96.2 (49.50-2.05) 96.2 (49.50-2.05)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.77 (at 2.05Å)	Xtriage
Refinement program	REFMAC 5.8.0352	Depositor
R, R_{free}	0.201 , 0.249 (Not available) , 0.232	Depositor DCC
R_{free} test set	1716 reflections (2.50%)	wwPDB-VP
Wilson B-factor (Å ²)	30.6	Xtriage
Anisotropy	0.322	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 37.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.019 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	9079	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.25% 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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.52	0/2208	0.96	1/2986 (0.0%)
1	B	0.52	0/2231	1.02	1/3019 (0.0%)
1	C	0.53	0/2226	1.02	0/3013
1	D	0.50	0/2200	0.97	0/2976
All	All	0.52	0/8865	0.99	2/11994 (0.0%)

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	3
1	B	0	2
1	C	0	2
1	D	0	1
All	All	0	8

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	55	ASN	CB-CA-C	5.40	119.05	109.65
1	A	271	GLU	CB-CG-CD	5.01	121.12	112.60

There are no chirality outliers.

All (8) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	172	ARG	Sidechain
1	A	267	ARG	Sidechain
1	A	44	ARG	Sidechain
1	B	267	ARG	Sidechain
1	B	274	ARG	Sidechain
1	C	172	ARG	Sidechain
1	C	261	ARG	Sidechain
1	D	107	ARG	Sidechain

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	2163	0	2158	30	0
1	B	2186	0	2183	25	0
1	C	2181	0	2175	31	0
1	D	2155	0	2147	33	0
2	A	10	0	0	0	0
2	B	5	0	0	0	0
2	C	10	0	0	0	0
2	D	5	0	0	0	0
3	D	6	0	8	0	0
4	A	106	0	0	3	0
4	B	102	0	0	0	0
4	C	85	0	0	0	0
4	D	65	0	0	1	0
All	All	9079	0	8671	117	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 7.

All (117) 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:D:23:VAL:HG21	1:D:274:ARG:HD3	1.61	0.82
1:B:42:MET:HG2	1:B:80:VAL:HG11	1.66	0.77
1:D:5:ILE:HG22	1:D:270:MET:HE1	1.74	0.69

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:82:LYS:HG3	1:C:91:ILE:HD12	1.76	0.66
1:A:23:VAL:CG2	1:A:274:ARG:HD3	2.25	0.66
1:B:32:MET:HG2	1:B:37:GLN:HE22	1.61	0.65
1:D:214:LEU:C	1:D:214:LEU:HD23	2.22	0.65
1:A:199:SER:N	1:A:200:PRO:CD	2.61	0.64
1:D:7:ASP:HB2	1:D:270:MET:HE3	1.81	0.63
1:D:120:ASN:HD22	1:D:121:SER:H	1.45	0.63
1:C:122:CYS:SG	1:C:129:ILE:HD12	2.39	0.63
1:D:135:PRO:HB2	1:D:142:ASP:OD1	1.99	0.62
1:C:57:ASN:HD22	1:C:57:ASN:H	1.48	0.62
1:B:248:LYS:HE3	1:B:252:GLU:OE2	2.00	0.61
1:D:97:MET:HE1	1:D:121:SER:C	2.25	0.61
1:A:155:GLY:O	1:A:159:LYS:HB2	2.02	0.60
1:D:46:ILE:HG23	1:D:50:LEU:HD23	1.82	0.60
1:B:37:GLN:HE21	1:B:42:MET:HE2	1.67	0.60
1:D:199:SER:N	1:D:200:PRO:CD	2.66	0.59
1:A:90:LEU:HD23	4:A:442:HOH:O	2.04	0.58
1:A:214:LEU:HD23	1:A:214:LEU:C	2.29	0.57
1:A:23:VAL:HG22	1:A:274:ARG:HD3	1.86	0.57
1:C:97:MET:HE2	1:C:120:ASN:CG	2.29	0.57
1:C:240:ILE:HG12	1:C:266:VAL:HG12	1.87	0.57
1:B:155:GLY:O	1:B:159:LYS:HB2	2.04	0.57
1:C:97:MET:HE1	1:C:121:SER:CA	2.35	0.56
1:A:23:VAL:HG21	1:A:274:ARG:HD3	1.88	0.56
1:B:16:ASN:ND2	1:B:26:GLU:HG3	2.21	0.55
1:B:97:MET:HE3	1:B:122:CYS:HB3	1.87	0.55
1:C:97:MET:HE3	1:C:122:CYS:HB3	1.88	0.55
1:C:214:LEU:C	1:C:214:LEU:HD23	2.32	0.55
1:A:135:PRO:HB2	1:A:142:ASP:OD1	2.07	0.54
1:A:169:LYS:NZ	1:A:205:ASN:OD1	2.36	0.54
1:B:214:LEU:C	1:B:214:LEU:HD23	2.33	0.54
1:D:51:LEU:HD21	1:D:57:ASN:HB3	1.89	0.54
1:D:23:VAL:CG2	1:D:274:ARG:HD3	2.35	0.53
1:C:32:MET:HG2	1:C:42:MET:HG2	1.91	0.53
1:B:135:PRO:HB2	1:B:142:ASP:OD1	2.09	0.53
1:A:12:LYS:HD2	1:A:28:PHE:CD1	2.44	0.52
1:A:199:SER:N	1:A:200:PRO:HD3	2.25	0.52
1:A:58:ARG:HG2	1:A:86:ASP:HB3	1.90	0.52
1:C:199:SER:N	1:C:200:PRO:CD	2.74	0.51
1:C:32:MET:HE1	1:C:46:ILE:HD11	1.93	0.51
1:B:199:SER:N	1:B:200:PRO:CD	2.74	0.50

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:178:PRO:O	1:D:182:ILE:HG12	2.11	0.50
1:D:268:ASN:HB2	1:D:269:PRO:HD2	1.92	0.50
1:A:20:GLN:HB2	1:A:22:ARG:NH2	2.26	0.50
1:A:39:PRO:HG3	1:A:76:VAL:HG13	1.95	0.49
1:D:214:LEU:HD23	1:D:214:LEU:O	2.12	0.49
1:B:239:PHE:C	1:B:240:ILE:HD12	2.38	0.48
1:D:162:MET:HG2	1:D:210:ALA:HB1	1.95	0.48
1:D:97:MET:HE1	1:D:121:SER:CA	2.43	0.48
1:D:120:ASN:HD22	1:D:121:SER:N	2.10	0.48
1:B:63:HIS:CE1	1:B:90:LEU:HD11	2.48	0.48
1:A:26:GLU:C	1:A:27:ILE:HG13	2.38	0.47
1:C:20:GLN:O	1:C:22:ARG:NH1	2.47	0.47
1:D:223:LEU:HB3	1:D:260:ILE:HD12	1.95	0.47
1:D:65:PHE:CB	1:D:99:ALA:HB2	2.44	0.47
1:D:42:MET:HE1	1:D:77:VAL:HG13	1.96	0.47
1:A:223:LEU:HB3	1:A:260:ILE:HD12	1.97	0.47
1:B:98:LEU:HG	1:B:102:LYS:HD2	1.97	0.47
1:B:120:ASN:HD22	1:B:121:SER:H	1.61	0.47
1:C:32:MET:CE	1:C:46:ILE:HD11	2.45	0.47
1:D:223:LEU:HB3	1:D:260:ILE:CD1	2.45	0.47
1:A:65:PHE:CB	1:A:99:ALA:HB2	2.44	0.47
1:C:268:ASN:HB2	1:C:269:PRO:HD2	1.97	0.47
1:C:6:ALA:HB2	1:C:15:TRP:CE3	2.49	0.46
1:A:12:LYS:HB3	1:A:12:LYS:NZ	2.30	0.46
1:A:18:LEU:HA	1:A:22:ARG:O	2.16	0.46
1:C:3:ILE:HG12	1:C:18:LEU:HB2	1.98	0.45
1:A:159:LYS:NZ	1:B:135:PRO:O	2.50	0.45
1:D:78:ARG:HD2	4:D:601:HOH:O	2.16	0.45
1:A:268:ASN:HB2	1:A:269:PRO:HD2	1.99	0.45
1:A:18:LEU:HD21	1:A:274:ARG:HG2	1.98	0.45
1:A:135:PRO:O	1:B:159:LYS:NZ	2.50	0.45
1:D:228:MET:HG2	1:D:232:TYR:CE1	2.52	0.45
1:C:32:MET:HB3	1:C:42:MET:HE3	1.98	0.45
1:B:6:ALA:HB2	1:B:15:TRP:CE3	2.53	0.44
1:C:97:MET:HE1	1:C:121:SER:HA	1.99	0.44
1:C:100:ALA:HA	1:C:240:ILE:HD12	1.98	0.44
1:B:65:PHE:CB	1:B:99:ALA:HB2	2.48	0.44
1:D:101:ALA:HB1	1:D:124:TYR:CG	2.52	0.44
1:A:239:PHE:C	1:A:240:ILE:HD12	2.43	0.44
1:C:101:ALA:HB1	1:C:124:TYR:CG	2.53	0.44
1:C:97:MET:HE1	1:C:121:SER:C	2.43	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:7:ASP:CB	1:D:270:MET:HE3	2.45	0.43
1:A:32:MET:HG2	1:A:37:GLN:HE22	1.82	0.43
1:C:135:PRO:O	1:C:136:LEU:HB2	2.18	0.43
1:C:16:ASN:OD1	1:C:274:ARG:HD3	2.19	0.43
1:C:65:PHE:HA	1:C:94:ASP:O	2.18	0.43
1:A:6:ALA:HB2	1:A:15:TRP:CE3	2.54	0.42
1:A:59:PHE:O	1:A:87:VAL:HA	2.19	0.42
1:A:42:MET:HE1	1:A:77:VAL:HG13	2.01	0.42
1:D:97:MET:CE	1:D:121:SER:C	2.90	0.42
1:C:135:PRO:HB2	1:C:142:ASP:OD1	2.20	0.42
1:D:214:LEU:C	1:D:214:LEU:CD2	2.90	0.42
1:B:178:PRO:O	1:B:182:ILE:HG12	2.19	0.42
1:D:70:ILE:O	1:D:71:PRO:C	2.63	0.42
1:D:122:CYS:HB2	1:D:129:ILE:HG23	2.02	0.42
1:B:60:CYS:O	1:B:89:SER:HB2	2.20	0.42
1:B:42:MET:CG	1:B:80:VAL:HG11	2.42	0.41
1:C:74:VAL:HG22	1:C:75:PRO:HD3	2.02	0.41
1:C:119:SER:O	1:C:135:PRO:HB3	2.20	0.41
1:D:109:PRO:HA	1:D:124:TYR:O	2.20	0.41
1:A:102:LYS:NZ	4:A:405:HOH:O	2.47	0.41
1:C:183:GLU:OE1	1:C:187:ARG:NH2	2.43	0.41
1:D:198:ILE:C	1:D:200:PRO:HD2	2.46	0.41
1:C:57:ASN:HD22	1:C:57:ASN:N	2.11	0.41
1:C:109:PRO:HA	1:C:124:TYR:O	2.20	0.41
1:B:135:PRO:O	1:B:136:LEU:HB2	2.21	0.41
1:B:122:CYS:SG	1:B:129:ILE:HD12	2.61	0.41
1:A:90:LEU:CD2	4:A:442:HOH:O	2.66	0.40
1:D:97:MET:HE3	1:D:122:CYS:HB3	2.01	0.40
1:B:101:ALA:HB1	1:B:124:TYR:CG	2.56	0.40
1:C:102:LYS:HD3	1:C:276:TYR:CE2	2.57	0.40
1:D:42:MET:HE3	1:D:42:MET:HB3	1.90	0.40
1:B:15:TRP:HB2	1:B:27:ILE:HB	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	277/283 (98%)	269 (97%)	8 (3%)	0	100	100
1	B	281/283 (99%)	275 (98%)	6 (2%)	0	100	100
1	C	281/283 (99%)	275 (98%)	6 (2%)	0	100	100
1	D	276/283 (98%)	266 (96%)	10 (4%)	0	100	100
All	All	1115/1132 (98%)	1085 (97%)	30 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	232/237 (98%)	230 (99%)	2 (1%)	70	75
1	B	234/237 (99%)	228 (97%)	6 (3%)	40	37
1	C	233/237 (98%)	224 (96%)	9 (4%)	28	23
1	D	231/237 (98%)	227 (98%)	4 (2%)	53	53
All	All	930/948 (98%)	909 (98%)	21 (2%)	44	42

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	ARG
1	A	267	ARG
1	B	23	VAL
1	B	58	ARG
1	B	90	LEU
1	B	132	ASN
1	B	260	ILE
1	B	280	VAL

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	C	12	LYS
1	C	55	ASN
1	C	57	ASN
1	C	58	ARG
1	C	74	VAL
1	C	90	LEU
1	C	132	ASN
1	C	159	LYS
1	C	264	THR
1	D	42	MET
1	D	51	LEU
1	D	90	LEU
1	D	97	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	37	GLN
1	A	53	GLN
1	B	37	GLN
1	B	55	ASN
1	B	120	ASN
1	B	132	ASN
1	C	57	ASN
1	C	79	ASN
1	C	120	ASN
1	D	120	ASN
1	D	188	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

7 ligands are modelled in this entry.

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	$\# Z > 2$
2	SO4	B	301	-	4,4,4	0.35	0	6,6,6	0.08	0
2	SO4	D	502	-	4,4,4	0.34	0	6,6,6	0.10	0
2	SO4	C	302	-	4,4,4	0.31	0	6,6,6	0.07	0
2	SO4	C	301	-	4,4,4	0.33	0	6,6,6	0.07	0
2	SO4	A	302	-	4,4,4	0.32	0	6,6,6	0.11	0
2	SO4	A	301	-	4,4,4	0.31	0	6,6,6	0.12	0
3	GOL	D	501	-	5,5,5	0.09	0	5,5,5	0.34	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
3	GOL	D	501	-	-	4/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	D	501	GOL	O1-C1-C2-C3
3	D	501	GOL	C1-C2-C3-O3
3	D	501	GOL	O2-C2-C3-O3
3	D	501	GOL	O1-C1-C2-O2

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 [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	278/283 (98%)	-0.11	1 (0%) 88 90	16, 32, 48, 59	1 (0%)
1	B	283/283 (100%)	-0.02	3 (1%) 78 80	22, 33, 50, 64	0
1	C	283/283 (100%)	0.03	2 (0%) 84 85	25, 36, 57, 68	0
1	D	278/283 (98%)	0.14	5 (1%) 67 69	25, 38, 67, 89	0
All	All	1122/1132 (99%)	0.01	11 (0%) 79 81	16, 35, 56, 89	1 (0%)

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	90	LEU	2.9
1	C	90	LEU	2.8
1	D	9	GLY	2.5
1	D	11	THR	2.5
1	B	284	VAL	2.5
1	A	279	THR	2.2
1	B	58	ARG	2.2
1	D	42	MET	2.1
1	C	11	THR	2.1
1	D	168	GLU	2.0
1	D	36	PHE	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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(Å ²)	Q<0.9
2	SO4	A	302	5/5	0.81	0.11	68,71,78,81	0
3	GOL	D	501	6/6	0.88	0.12	41,49,55,58	0
2	SO4	D	502	5/5	0.90	0.09	61,73,77,83	0
2	SO4	B	301	5/5	0.93	0.08	55,56,62,72	0
2	SO4	A	301	5/5	0.94	0.09	57,57,61,62	0
2	SO4	C	301	5/5	0.95	0.09	54,54,63,73	0
2	SO4	C	302	5/5	0.96	0.08	46,56,66,78	0

6.5 Other polymers [i](#)

There are no such residues in this entry.