



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 11:58 PM UTC

PDB ID : 6O60 / pdb_00006o60
Title : Crystal structure of GGTase3-FBXL2-SKP1 complex
Authors : Wang, H.; Zheng, N.
Deposited on : 2019-03-04
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

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
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

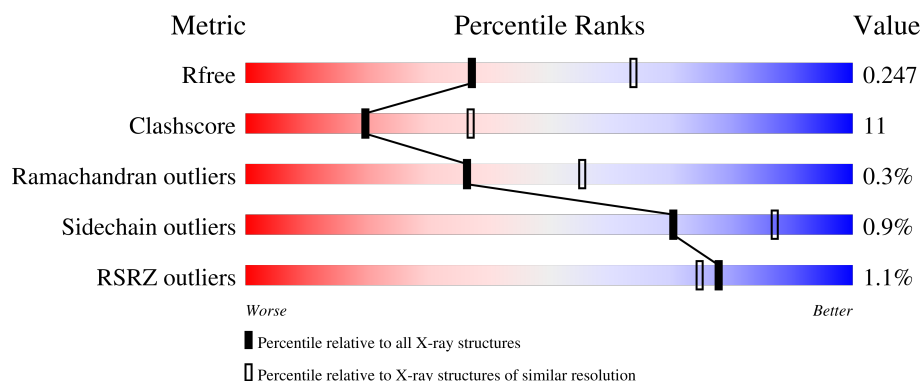
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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))
R_{free}	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (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 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	407	55% 19% 26%
2	B	332	82% 14% •
3	C	424	67% 25% 8%
4	D	164	64% 24% • 10%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 9299 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 prenyltransferase alpha subunit repeat-containing protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	300	Total	C	N	O	S	0	0	0
			2504	1609	450	436	9			

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-4	SER	-	expression tag	UNP Q7Z6K3
A	-3	GLY	-	expression tag	UNP Q7Z6K3
A	-2	SER	-	expression tag	UNP Q7Z6K3
A	-1	GLY	-	expression tag	UNP Q7Z6K3
A	0	GLY	-	expression tag	UNP Q7Z6K3

- Molecule 2 is a protein called Geranylgeranyl transferase type-2 subunit beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	320	Total	C	N	O	S	0	0	0
			2510	1603	416	471	20			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	0	SER	-	expression tag	UNP P53611

- Molecule 3 is a protein called F-box/LRR-repeat protein 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	389	Total	C	N	O	S	0	0	0
			3029	1883	552	563	31			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	0	SER	-	expression tag	UNP Q9UKC9

- Molecule 4 is a protein called S-phase kinase-associated protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	148	Total	C	N	O	S	0	0	0
			1185	747	193	239	6			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	0	SER	-	expression tag	UNP P63208

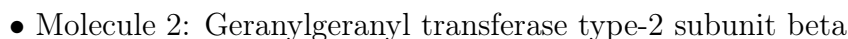
- Molecule 5 is ZINC ION (CCD ID: ZN) (formula: Zn).

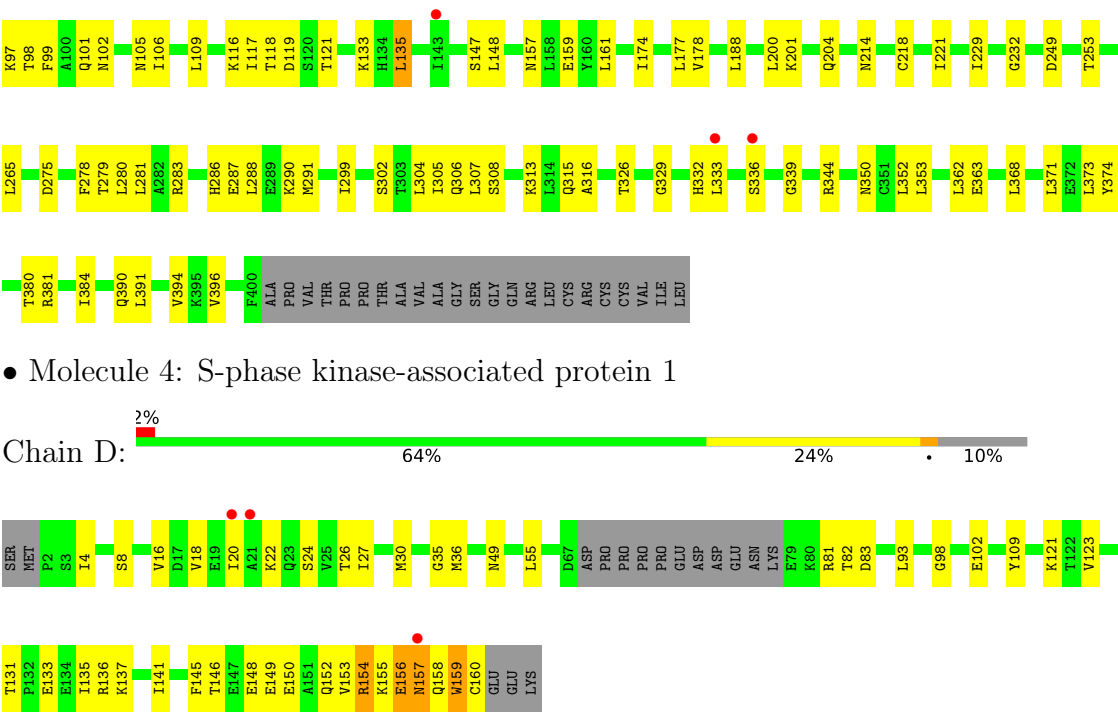
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	1	Total	Zn	0	0
			1	1		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	15	Total	O	0	0
			15	15		
6	B	34	Total	O	0	0
			34	34		
6	C	20	Total	O	0	0
			20	20		
6	D	1	Total	O	0	0
			1	1		

- Molecule 1: Protein prenyltransferase alpha subunit repeat-containing protein 1





● Molecule 4: S-phase kinase-associated protein 1

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	86.84Å 99.32Å 151.20Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.45 – 2.50 49.45 – 2.50	Depositor EDS
% Data completeness (in resolution range)	90.2 (49.45-2.50) 90.2 (49.45-2.50)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.60 (at 2.51Å)	Xtriage
Refinement program	PHENIX (1.13_2998: ???)	Depositor
R, R_{free}	0.194 , 0.247 0.195 , 0.247	Depositor DCC
R_{free} test set	2069 reflections (4.52%)	wwPDB-VP
Wilson B-factor (Å ²)	50.6	Xtriage
Anisotropy	0.289	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 30.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	9299	wwPDB-VP
Average B, all atoms (Å ²)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.59% 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: ZN

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	0.38	0/2568	0.58	0/3482
2	B	0.40	0/2567	0.59	0/3478
3	C	0.36	0/3065	0.64	0/4131
4	D	0.45	0/1203	0.76	3/1625 (0.2%)
All	All	0.39	0/9403	0.63	3/12716 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	154	ARG	N-CA-C	-9.82	99.85	113.20
4	D	159	TRP	N-CA-C	-6.40	97.17	110.80
4	D	159	TRP	CB-CA-C	5.34	121.04	110.42

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	2504	0	2505	61	0
2	B	2510	0	2450	38	0
3	C	3029	0	3095	78	0
4	D	1185	0	1173	43	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	B	1	0	0	0	0
6	A	15	0	0	2	0
6	B	34	0	0	1	0
6	C	20	0	0	1	0
6	D	1	0	0	0	0
All	All	9299	0	9223	212	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 (212) 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)
2:B:71:ILE:O	2:B:75:ILE:HD12	1.53	1.06
4:D:156:GLU:O	4:D:159:TRP:HD1	1.42	1.03
4:D:158:GLN:O	4:D:159:TRP:C	2.05	0.94
2:B:313:PHE:HB3	2:B:315:MET:HE3	1.50	0.94
4:D:153:VAL:C	4:D:155:LYS:H	1.76	0.94
3:C:279:THR:HG22	3:C:306:GLN:HG3	1.50	0.92
2:B:71:ILE:O	2:B:75:ILE:CD1	2.18	0.91
3:C:315:GLN:HE21	3:C:344:ARG:HD3	1.35	0.91
3:C:12:ASN:HD21	3:C:42:TRP:HE1	1.20	0.89
3:C:363:GLU:HG2	3:C:390:GLN:HG2	1.58	0.86
1:A:171:GLU:HB3	1:A:174:GLN:HG3	1.57	0.84
1:A:12:VAL:HG13	1:A:64:LEU:HD12	1.60	0.81
4:D:156:GLU:O	4:D:159:TRP:CD1	2.31	0.80
3:C:302:SER:HA	3:C:305:ILE:HD12	1.66	0.78
1:A:77:THR:HG23	1:A:79:LYS:HG3	1.64	0.78
4:D:153:VAL:C	4:D:155:LYS:N	2.43	0.75
3:C:200:LEU:HD22	3:C:229:ILE:HD11	1.70	0.74
4:D:158:GLN:O	4:D:160:CYS:N	2.22	0.71
1:A:367:LEU:HD11	1:A:397:LEU:HD22	1.73	0.71
3:C:67:ARG:HH22	3:C:74:LYS:HZ1	1.39	0.70
3:C:283:ARG:O	3:C:286:HIS:NE2	2.25	0.70
1:A:32:ILE:CG2	1:A:128:LEU:HD21	2.21	0.70
3:C:336:SER:HB2	3:C:339:GLY:H	1.55	0.69
3:C:304:LEU:HB3	3:C:333:LEU:HD21	1.75	0.69
4:D:146:THR:HG23	4:D:149:GLU:H	1.59	0.68
1:A:68:VAL:HG12	1:A:94:LEU:HG	1.75	0.68
1:A:202:ILE:HD13	1:A:242:PHE:CD2	2.29	0.67
2:B:52:TRP:HA	2:B:292:LEU:HD22	1.77	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:133:LYS:HE2	3:C:159:GLU:CD	2.20	0.67
3:C:99:PHE:HZ	3:C:106:ILE:HD13	1.60	0.66
1:A:231:VAL:HG13	1:A:312:LEU:HD11	1.78	0.66
4:D:156:GLU:HG3	4:D:157:ASN:N	2.11	0.66
1:A:42:TYR:CD2	1:A:44:ARG:HB2	2.30	0.66
4:D:81:ARG:HD2	4:D:83:ASP:OD1	1.96	0.66
3:C:67:ARG:NE	3:C:71:ASN:ND2	2.43	0.65
1:A:148:GLN:HE21	1:A:151:ILE:HD11	1.62	0.65
3:C:249:ASP:O	3:C:253:THR:HG23	1.97	0.64
1:A:32:ILE:HG21	1:A:128:LEU:HD21	1.80	0.63
3:C:68:VAL:O	3:C:72:ILE:HG12	1.97	0.63
3:C:200:LEU:HD13	3:C:229:ILE:HD12	1.79	0.63
1:A:74:LEU:HA	1:A:77:THR:HG22	1.81	0.63
2:B:15:PRO:HG3	2:B:273:GLU:HG3	1.81	0.62
3:C:67:ARG:HE	3:C:71:ASN:ND2	1.97	0.62
4:D:154:ARG:HG2	4:D:158:GLN:OE1	1.99	0.62
2:B:75:ILE:HD12	2:B:75:ILE:H	1.65	0.62
2:B:229:LEU:HD13	2:B:263:LEU:HD13	1.82	0.61
3:C:253:THR:HG22	3:C:280:LEU:HD13	1.81	0.61
4:D:148:GLU:O	4:D:152:GLN:HG2	2.01	0.61
3:C:16:PRO:HD2	3:C:19:LEU:HD22	1.82	0.60
1:A:202:ILE:HD13	1:A:242:PHE:CG	2.37	0.59
1:A:320:PHE:CZ	1:A:367:LEU:HD13	2.38	0.59
2:B:206:GLN:OE1	2:B:209:GLN:NE2	2.29	0.58
3:C:305:ILE:HG12	3:C:332:HIS:ND1	2.19	0.58
1:A:370:GLU:OE2	1:A:393:TYR:OH	2.12	0.58
4:D:18:VAL:O	4:D:22:LYS:HG3	2.04	0.58
3:C:93:ASP:HA	3:C:121:THR:HG22	1.86	0.57
3:C:218:CYS:O	3:C:221:ILE:HG12	2.04	0.57
3:C:344:ARG:HA	3:C:368:LEU:HA	1.86	0.57
3:C:291:MET:HE1	3:C:307:LEU:HG	1.87	0.57
1:A:202:ILE:CD1	1:A:242:PHE:CG	2.88	0.57
1:A:68:VAL:HG13	1:A:90:VAL:HG13	1.87	0.56
3:C:363:GLU:HG2	3:C:390:GLN:CG	2.31	0.56
2:B:274:GLU:OE2	6:B:501:HOH:O	2.18	0.56
1:A:42:TYR:CE2	1:A:44:ARG:HB2	2.41	0.56
4:D:16:VAL:HG13	4:D:20:ILE:HD11	1.87	0.56
4:D:35:GLY:HA2	4:D:36:MET:HB3	1.87	0.56
1:A:70:ASN:ND2	6:A:501:HOH:O	2.38	0.56
1:A:203:TRP:NE1	1:A:207:HIS:HD1	2.04	0.56
2:B:255:LEU:CD2	2:B:260:ARG:HD3	2.35	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:391:LEU:O	3:C:394:VAL:HG12	2.06	0.56
2:B:229:LEU:HD23	2:B:245:VAL:HB	1.87	0.56
1:A:317:ARG:NH2	2:B:221:GLU:OE1	2.30	0.55
2:B:295:ILE:HG23	2:B:306:ILE:HD13	1.89	0.55
3:C:32:LEU:HD21	3:C:51:ASN:HB3	1.87	0.55
3:C:117:ILE:HD12	3:C:121:THR:HG21	1.89	0.55
1:A:223:THR:HG23	1:A:240:ARG:HH21	1.71	0.55
4:D:136:ARG:NH2	4:D:150:GLU:OE2	2.40	0.55
3:C:70:GLU:O	3:C:73:SER:OG	2.24	0.55
2:B:229:LEU:HB2	2:B:242:SER:HA	1.88	0.54
3:C:304:LEU:HB3	3:C:333:LEU:CD2	2.37	0.54
2:B:213:ASP:CG	2:B:254:ARG:HH12	2.15	0.54
4:D:35:GLY:HA3	4:D:36:MET:C	2.33	0.54
3:C:78:GLY:O	3:C:105:ASN:ND2	2.32	0.54
1:A:190:ARG:HD2	6:A:510:HOH:O	2.06	0.53
2:B:321:GLN:O	2:B:325:VAL:HG22	2.09	0.53
3:C:148:LEU:HB3	3:C:177:LEU:HB2	1.89	0.53
4:D:26:THR:HG22	4:D:30:MET:CE	2.39	0.53
4:D:82:THR:HA	4:D:121:LYS:HG2	1.90	0.52
3:C:381:ARG:HG3	6:C:514:HOH:O	2.09	0.52
4:D:156:GLU:HG2	4:D:157:ASN:OD1	2.09	0.52
3:C:97:LYS:O	3:C:101:GLN:HG2	2.10	0.52
1:A:95:LEU:HD22	1:A:105:TRP:CE2	2.44	0.52
3:C:188:LEU:HD13	3:C:214:ASN:HB2	1.92	0.52
1:A:219:GLU:O	1:A:223:THR:HG22	2.11	0.51
4:D:155:LYS:HG2	4:D:156:GLU:N	2.25	0.51
4:D:98:GLY:O	4:D:102:GLU:HG2	2.10	0.51
3:C:67:ARG:HH22	3:C:74:LYS:NZ	2.07	0.51
3:C:290:LYS:HG3	3:C:316:ALA:HB3	1.92	0.51
3:C:326:THR:HG22	3:C:353:LEU:HD23	1.93	0.50
1:A:16:VAL:O	1:A:20:THR:HG23	2.11	0.50
1:A:181:MET:HE1	1:A:205:LEU:HD21	1.94	0.50
2:B:39:TYR:O	2:B:46:ARG:NH2	2.44	0.50
3:C:204:GLN:HB2	3:C:232:GLY:HA3	1.92	0.50
3:C:98:THR:HA	3:C:101:GLN:HE21	1.77	0.50
3:C:281:LEU:HD11	3:C:288:LEU:HD11	1.94	0.50
1:A:32:ILE:HG23	1:A:128:LEU:HD21	1.92	0.49
3:C:88:CYS:O	3:C:91:VAL:HG22	2.12	0.49
2:B:231:GLY:O	2:B:232:ARG:HG3	2.12	0.49
2:B:325:VAL:O	2:B:325:VAL:HG23	2.13	0.49
1:A:42:TYR:HD2	1:A:44:ARG:HB2	1.74	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:255:LEU:HD21	2:B:260:ARG:HD3	1.95	0.48
3:C:9:GLY:HA3	4:D:141:ILE:HD13	1.95	0.48
1:A:293:GLU:O	1:A:297:GLU:HG3	2.13	0.48
1:A:75:TYR:HB2	1:A:82:LEU:HD12	1.94	0.48
1:A:366:GLY:N	1:A:369:MET:SD	2.86	0.48
1:A:367:LEU:HD12	1:A:370:GLU:HB3	1.93	0.48
2:B:103:GLN:HG2	2:B:315:MET:HE2	1.96	0.48
1:A:367:LEU:HD11	1:A:397:LEU:CD2	2.41	0.48
1:A:195:TYR:CE1	2:B:188:GLU:HB3	2.49	0.48
4:D:24:SER:HB3	4:D:27:ILE:HB	1.96	0.48
1:A:101:PHE:CE1	2:B:91:HIS:CE1	3.02	0.48
2:B:129:GLU:H	2:B:129:GLU:CD	2.19	0.48
3:C:275:ASP:O	3:C:279:THR:HG23	2.14	0.47
4:D:26:THR:O	4:D:30:MET:HG3	2.14	0.47
2:B:51:TYR:OH	2:B:314:CYS:HB3	2.15	0.47
4:D:156:GLU:C	4:D:158:GLN:H	2.23	0.47
4:D:133:GLU:HG3	4:D:137:LYS:HE2	1.97	0.47
2:B:103:GLN:HG2	2:B:315:MET:CE	2.44	0.47
2:B:250:LYS:HD2	2:B:300:LEU:HD11	1.96	0.47
1:A:52:ASN:HB2	1:A:53:LYS:NZ	2.30	0.47
2:B:15:PRO:HG3	2:B:273:GLU:CG	2.45	0.47
3:C:92:GLY:HA3	3:C:116:LYS:HE2	1.97	0.47
3:C:350:ASN:O	3:C:352:LEU:HD12	2.15	0.47
4:D:26:THR:HG22	4:D:30:MET:HE3	1.96	0.47
1:A:291:LEU:HA	1:A:294:GLU:HG2	1.97	0.46
3:C:67:ARG:NE	3:C:71:ASN:HD22	2.12	0.46
1:A:292:LEU:HD13	1:A:323:GLN:HA	1.96	0.46
4:D:35:GLY:CA	4:D:36:MET:C	2.89	0.46
2:B:79:GLN:NE2	2:B:119:VAL:HG23	2.30	0.46
1:A:64:LEU:HD23	1:A:97:LEU:HD11	1.98	0.46
4:D:156:GLU:CG	4:D:157:ASN:N	2.73	0.46
4:D:159:TRP:O	4:D:160:CYS:HB2	2.15	0.46
2:B:49:GLY:HA2	2:B:52:TRP:CE3	2.51	0.45
2:B:75:ILE:HD12	2:B:75:ILE:N	2.30	0.45
3:C:17:LYS:CG	3:C:45:LEU:HD21	2.46	0.45
3:C:279:THR:HG22	3:C:306:GLN:CG	2.35	0.45
3:C:278:PHE:CE2	3:C:299:ILE:HD11	2.51	0.45
4:D:145:PHE:CB	4:D:150:GLU:HG2	2.46	0.45
2:B:124:LYS:HD2	2:B:161:ALA:O	2.16	0.45
2:B:155:LEU:HD11	2:B:315:MET:HE1	1.98	0.45
1:A:87:LEU:HD22	1:A:107:VAL:HG11	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:75:ILE:HG23	2:B:101:ALA:HB1	1.98	0.45
1:A:145:TRP:O	1:A:149:GLN:HG2	2.17	0.45
3:C:371:LEU:HD22	3:C:373:LEU:HG	1.99	0.45
3:C:118:THR:HG22	3:C:119:ASP:N	2.32	0.45
3:C:23:ILE:HD12	4:D:123:VAL:HB	2.00	0.44
1:A:320:PHE:HZ	1:A:367:LEU:HD13	1.81	0.44
3:C:305:ILE:O	3:C:308:SER:OG	2.27	0.44
3:C:67:ARG:NH2	3:C:74:LYS:HZ1	2.11	0.44
4:D:4:ILE:HG12	4:D:18:VAL:HG12	1.99	0.44
4:D:150:GLU:HA	4:D:150:GLU:OE1	2.18	0.44
1:A:87:LEU:HD23	1:A:87:LEU:HA	1.77	0.44
1:A:113:LEU:HD12	1:A:113:LEU:HA	1.77	0.44
4:D:49:ASN:HB2	4:D:109:TYR:CE2	2.53	0.44
3:C:17:LYS:HG3	3:C:45:LEU:HD21	2.00	0.44
3:C:133:LYS:HE3	3:C:157:ASN:O	2.17	0.44
4:D:146:THR:HG22	4:D:149:GLU:OE1	2.18	0.44
2:B:128:LYS:HE3	2:B:128:LYS:HB3	1.81	0.43
3:C:116:LYS:O	3:C:116:LYS:HG2	2.18	0.43
1:A:375:ASP:OD1	1:A:394:ARG:NH2	2.45	0.43
3:C:380:THR:O	3:C:384:ILE:HG12	2.19	0.43
3:C:201:LYS:O	3:C:204:GLN:HG3	2.18	0.43
4:D:93:LEU:HD23	4:D:93:LEU:HA	1.70	0.43
1:A:225:HIS:O	1:A:229:MET:HG3	2.19	0.43
2:B:300:LEU:HA	2:B:300:LEU:HD12	1.78	0.43
4:D:131:THR:O	4:D:135:ILE:HG12	2.19	0.43
1:A:368:GLU:O	1:A:372:ARG:HG2	2.18	0.43
1:A:65:LEU:O	1:A:65:LEU:HG	2.19	0.43
3:C:287:GLU:OE1	3:C:313:LYS:HD2	2.18	0.43
1:A:41:ARG:HA	3:C:381:ARG:HH12	1.84	0.42
3:C:99:PHE:HZ	3:C:106:ILE:CD1	2.31	0.42
4:D:8:SER:HB3	4:D:55:LEU:HD12	2.01	0.42
3:C:67:ARG:CZ	3:C:71:ASN:ND2	2.83	0.42
3:C:80:LEU:HD12	3:C:80:LEU:HA	1.64	0.42
3:C:305:ILE:HG12	3:C:332:HIS:CE1	2.54	0.42
1:A:75:TYR:CZ	1:A:87:LEU:HD11	2.53	0.42
1:A:92:CYS:O	1:A:96:LEU:HG	2.20	0.42
2:B:67:ASN:O	2:B:71:ILE:HG12	2.20	0.42
1:A:14:ARG:HG2	3:C:374:TYR:CG	2.55	0.42
1:A:14:ARG:HE	3:C:374:TYR:CB	2.33	0.42
1:A:179:GLU:O	1:A:183:VAL:HG23	2.20	0.42
3:C:119:ASP:HA	3:C:147:SER:OG	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:55:LEU:HA	4:D:55:LEU:HD23	1.86	0.41
1:A:32:ILE:HG23	1:A:128:LEU:HD11	2.03	0.41
3:C:283:ARG:O	3:C:286:HIS:CE1	2.72	0.41
2:B:38:GLU:O	2:B:38:GLU:HG2	2.19	0.41
1:A:96:LEU:HD23	1:A:96:LEU:HA	1.78	0.41
1:A:68:VAL:CG1	1:A:90:VAL:HG13	2.49	0.41
3:C:84:SER:O	3:C:85:LEU:HD23	2.20	0.41
1:A:99:PRO:HB2	1:A:139:THR:OG1	2.20	0.41
4:D:145:PHE:HB3	4:D:150:GLU:HG2	2.02	0.41
3:C:39:SER:OG	3:C:40:LYS:N	2.54	0.41
3:C:67:ARG:HE	3:C:71:ASN:HD21	1.67	0.41
3:C:174:ILE:O	3:C:178:VAL:HG22	2.21	0.41
3:C:204:GLN:HB3	3:C:229:ILE:HA	2.03	0.41
3:C:135:LEU:HD23	3:C:135:LEU:HA	1.95	0.41
1:A:12:VAL:HG11	1:A:67:TYR:CD1	2.57	0.40
1:A:374:ILE:HG22	1:A:394:ARG:NH2	2.37	0.40
3:C:329:GLY:O	3:C:333:LEU:HD23	2.22	0.40
4:D:16:VAL:CG1	4:D:20:ILE:HD11	2.51	0.40
3:C:73:SER:HB2	3:C:102:ASN:HB2	2.03	0.40
4:D:156:GLU:CG	4:D:157:ASN:OD1	2.70	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	292/407 (72%)	281 (96%)	11 (4%)	0	100	100
2	B	316/332 (95%)	310 (98%)	5 (2%)	1 (0%)	36	55
3	C	385/424 (91%)	366 (95%)	18 (5%)	1 (0%)	36	55
4	D	144/164 (88%)	136 (94%)	7 (5%)	1 (1%)	18	34
All	All	1137/1327 (86%)	1093 (96%)	41 (4%)	3 (0%)	36	55

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
4	D	157	ASN
3	C	362	LEU
2	B	326	GLN

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	277/370 (75%)	274 (99%)	3 (1%)	65	84
2	B	273/284 (96%)	273 (100%)	0	100	100
3	C	346/376 (92%)	341 (99%)	5 (1%)	59	81
4	D	135/151 (89%)	134 (99%)	1 (1%)	76	89
All	All	1031/1181 (87%)	1022 (99%)	9 (1%)	70	87

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	113	LEU
1	A	116	THR
1	A	125	LEU
3	C	109	LEU
3	C	135	LEU
3	C	161	LEU
3	C	265	LEU
3	C	396	VAL
4	D	156	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	148	GLN
1	A	149	GLN
1	A	241	GLN

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Mol	Chain	Res	Type
2	B	193	GLN
2	B	208	HIS
2	B	321	GLN
3	C	12	ASN
3	C	71	ASN
3	C	101	GLN
3	C	194	GLN
3	C	315	GLN
3	C	390	GLN
4	D	23	GLN
4	D	90	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 ⓘ

Of 1 ligands modelled in this entry, 1 is monoatomic - leaving 0 for Mogul analysis.

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 ⓘ

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 [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	300/407 (73%)	-0.11	1 (0%) 90 87	33, 51, 83, 96	0
2	B	320/332 (96%)	-0.29	3 (0%) 81 78	35, 44, 68, 98	0
3	C	389/424 (91%)	-0.03	6 (1%) 72 68	35, 56, 79, 113	0
4	D	148/164 (90%)	0.04	3 (2%) 65 61	38, 59, 103, 122	0
All	All	1157/1327 (87%)	-0.11	13 (1%) 78 75	33, 52, 82, 122	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	327	PRO	3.3
2	B	328	GLU	3.1
3	C	143	ILE	2.9
4	D	20	ILE	2.6
3	C	76	CYS	2.6
2	B	31	GLY	2.5
3	C	62	THR	2.4
1	A	42	TYR	2.3
4	D	21	ALA	2.1
3	C	336	SER	2.1
4	D	157	ASN	2.1
3	C	333	LEU	2.0
3	C	66	GLY	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($\text{\AA}^2$)	Q<0.9
5	ZN	B	401	1/1	0.99	0.03	41,41,41,41	0

6.5 Other polymers [i](#)

There are no such residues in this entry.