



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 03:58 AM UTC

PDB ID : 9JXX / pdb_00009jxx
Title : Crystal structure of SiRe_0806 in complex with cA4
Authors : Wang, F.; Zhao, P.; Bi, X.; She, Q.
Deposited on : 2024-10-12
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

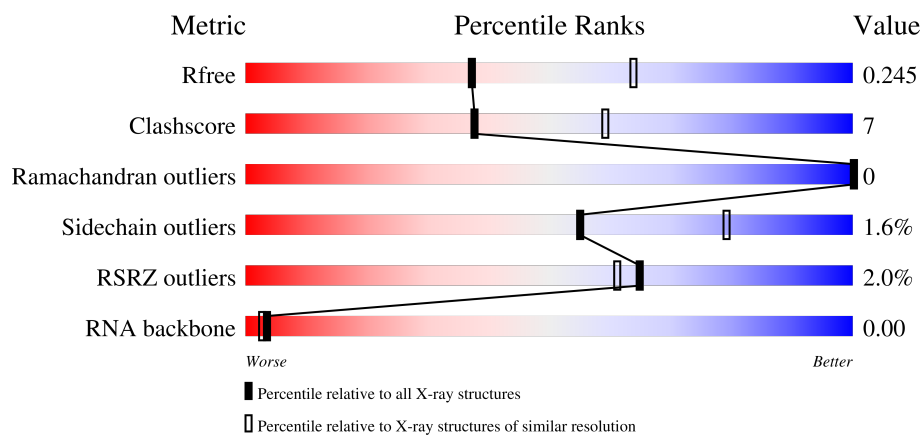
1 Overall quality at a glance

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)
RNA backbone	3983	1003 (2.78-2.22)

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	429	
1	B	429	
1	C	429	
1	D	429	

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Mol	Chain	Length	Quality of chain
2	X	4	 50%50%
2	Y	4	 75%25%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 13662 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 PIN domain-containing protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	412	Total	C	N	O	S	0	0	0
			3335	2132	566	630	7			
1	D	418	Total	C	N	O	S	0	0	0
			3394	2168	578	640	8			
1	A	407	Total	C	N	O	S	0	0	0
			3294	2105	560	622	7			
1	C	418	Total	C	N	O	S	0	0	0
			3387	2162	578	639	8			

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	418	ALA	-	expression tag	UNP F0NH84
B	419	ASP	-	expression tag	UNP F0NH84
B	420	GLN	-	expression tag	UNP F0NH84
B	421	ALA	-	expression tag	UNP F0NH84
B	422	ALA	-	expression tag	UNP F0NH84
B	423	ALA	-	expression tag	UNP F0NH84
B	424	HIS	-	expression tag	UNP F0NH84
B	425	HIS	-	expression tag	UNP F0NH84
B	426	HIS	-	expression tag	UNP F0NH84
B	427	HIS	-	expression tag	UNP F0NH84
B	428	HIS	-	expression tag	UNP F0NH84
B	429	HIS	-	expression tag	UNP F0NH84
D	418	ALA	-	expression tag	UNP F0NH84
D	419	ASP	-	expression tag	UNP F0NH84
D	420	GLN	-	expression tag	UNP F0NH84
D	421	ALA	-	expression tag	UNP F0NH84
D	422	ALA	-	expression tag	UNP F0NH84
D	423	ALA	-	expression tag	UNP F0NH84
D	424	HIS	-	expression tag	UNP F0NH84
D	425	HIS	-	expression tag	UNP F0NH84
D	426	HIS	-	expression tag	UNP F0NH84

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Chain	Residue	Modelled	Actual	Comment	Reference
D	427	HIS	-	expression tag	UNP F0NH84
D	428	HIS	-	expression tag	UNP F0NH84
D	429	HIS	-	expression tag	UNP F0NH84
A	418	ALA	-	expression tag	UNP F0NH84
A	419	ASP	-	expression tag	UNP F0NH84
A	420	GLN	-	expression tag	UNP F0NH84
A	421	ALA	-	expression tag	UNP F0NH84
A	422	ALA	-	expression tag	UNP F0NH84
A	423	ALA	-	expression tag	UNP F0NH84
A	424	HIS	-	expression tag	UNP F0NH84
A	425	HIS	-	expression tag	UNP F0NH84
A	426	HIS	-	expression tag	UNP F0NH84
A	427	HIS	-	expression tag	UNP F0NH84
A	428	HIS	-	expression tag	UNP F0NH84
A	429	HIS	-	expression tag	UNP F0NH84
C	418	ALA	-	expression tag	UNP F0NH84
C	419	ASP	-	expression tag	UNP F0NH84
C	420	GLN	-	expression tag	UNP F0NH84
C	421	ALA	-	expression tag	UNP F0NH84
C	422	ALA	-	expression tag	UNP F0NH84
C	423	ALA	-	expression tag	UNP F0NH84
C	424	HIS	-	expression tag	UNP F0NH84
C	425	HIS	-	expression tag	UNP F0NH84
C	426	HIS	-	expression tag	UNP F0NH84
C	427	HIS	-	expression tag	UNP F0NH84
C	428	HIS	-	expression tag	UNP F0NH84
C	429	HIS	-	expression tag	UNP F0NH84

- Molecule 2 is a RNA chain called RNA (5'-R(P*AP*AP*AP*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	Y	4	Total 88	C 40	N 20	O 24	P 4	0	0	0
2	X	4	Total 88	C 40	N 20	O 24	P 4	0	0	0

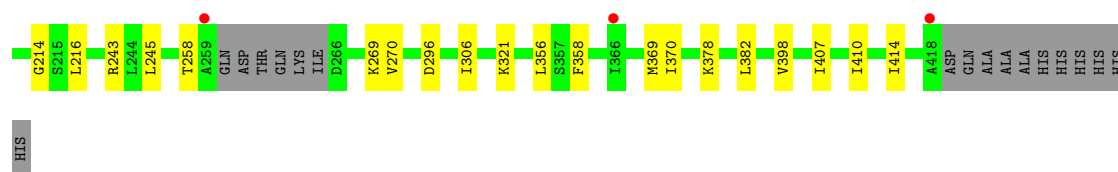
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	22	Total	O	0	0
			22	22		

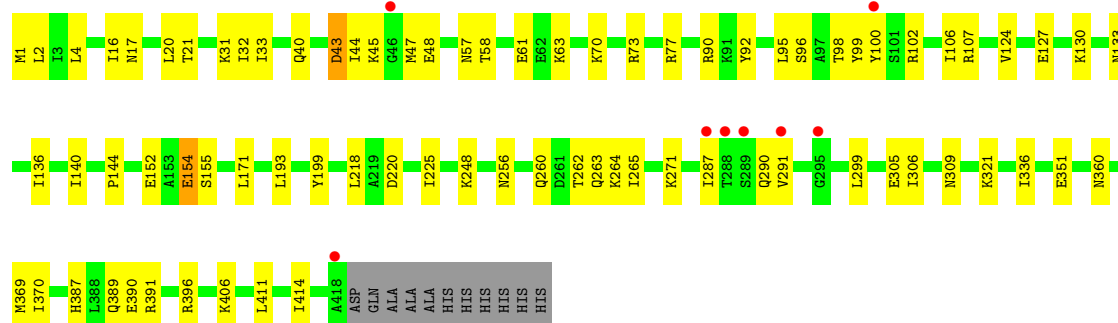
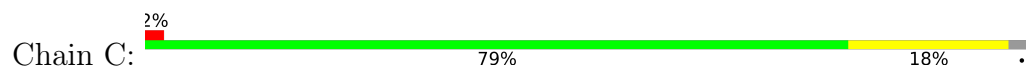
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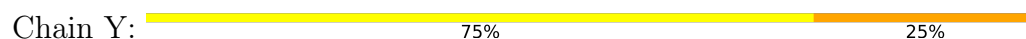
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	D	12	Total 12	O 12	0	0
3	Y	3	Total 3	O 3	0	0
3	A	17	Total 17	O 17	0	0
3	C	20	Total 20	O 20	0	0
3	X	2	Total 2	O 2	0	0



- Molecule 1: PIN domain-containing protein



- Molecule 2: RNA (5'-R(P*AP*AP*AP*A)-3')



- Molecule 2: RNA (5'-R(P*AP*AP*AP*A)-3')



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	79.47Å 81.84Å 86.22Å 75.47° 87.09° 61.09°	Depositor
Resolution (Å)	48.82 – 2.50 48.82 – 2.50	Depositor EDS
% Data completeness (in resolution range)	92.7 (48.82-2.50) 92.7 (48.82-2.50)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.01 (at 2.51Å)	Xtriage
Refinement program	PHENIX (1.19_4092: ???)	Depositor
R, R_{free}	0.206 , 0.244 0.207 , 0.245	Depositor DCC
R_{free} test set	2902 reflections (4.57%)	wwPDB-VP
Wilson B-factor (Å ²)	54.5	Xtriage
Anisotropy	0.438	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 41.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	13662	wwPDB-VP
Average B, all atoms (Å ²)	66.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.72% 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](#)

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	0.11	0/3346	0.28	0/4513
1	B	0.11	0/3388	0.33	0/4569
1	C	0.10	0/3440	0.28	0/4637
1	D	0.13	0/3448	0.32	0/4648
2	X	0.12	0/99	0.31	0/152
2	Y	0.13	0/99	0.34	0/152
All	All	0.12	0/13820	0.30	0/18671

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 [i](#)

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	3294	0	3346	34	0
1	B	3335	0	3393	48	0
1	C	3387	0	3464	61	0
1	D	3394	0	3471	61	0
2	X	88	0	44	3	0
2	Y	88	0	44	3	0
3	A	17	0	0	0	0
3	B	22	0	0	1	0
3	C	20	0	0	1	0
3	D	12	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	X	2	0	0	0	0
3	Y	3	0	0	0	0
All	All	13662	0	13762	191	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 (191) 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:6:THR:HG21	1:D:19:ILE:HD11	1.67	0.76
1:B:150:VAL:H	1:B:400:LYS:HE3	1.52	0.75
1:C:31:LYS:HD2	1:C:33:ILE:HD11	1.71	0.72
1:D:68:GLY:HA3	1:D:71:LEU:HD13	1.73	0.71
1:C:305:GLU:O	1:C:309:ASN:ND2	2.24	0.69
1:B:387:HIS:HB3	1:B:390:GLU:HB3	1.73	0.68
1:A:136:ILE:O	1:C:70:LYS:HE3	1.94	0.67
1:C:90:ARG:HD2	1:C:92:TYR:HE2	1.61	0.66
1:A:111:LEU:O	1:C:90:ARG:NH1	2.29	0.65
1:B:86:ILE:HD12	1:B:94:ALA:HB1	1.79	0.64
1:B:208:ILE:O	1:B:212:GLU:HG3	1.97	0.64
1:B:136:ILE:HG13	1:B:137:GLY:H	1.63	0.64
1:C:20:LEU:HD21	1:C:47:MET:HE1	1.80	0.63
1:A:98:THR:HG22	1:A:106:ILE:HG21	1.80	0.62
1:B:273:ILE:O	1:B:277:LEU:HG	2.00	0.61
1:B:193:LEU:HD13	1:B:199:TYR:HE2	1.66	0.60
1:A:48:GLU:HG3	1:A:60:ILE:HD12	1.83	0.60
1:C:16:ILE:O	1:C:20:LEU:HD23	2.01	0.60
1:D:362:TYR:HD2	1:A:321:LYS:HD3	1.67	0.60
1:C:271:LYS:NZ	3:C:501:HOH:O	2.33	0.59
1:C:411:LEU:HA	1:C:414:ILE:HD12	1.83	0.59
1:B:3:ILE:N	1:B:82:ASP:OD1	2.31	0.59
1:D:73:ARG:HG3	1:D:100:TYR:CD2	2.38	0.58
1:D:238:ARG:NH1	1:D:239:LYS:HG3	2.18	0.58
1:C:73:ARG:HG2	1:C:100:TYR:CE2	2.39	0.57
1:D:70:LYS:HA	1:D:73:ARG:HB2	1.84	0.57
1:C:77:ARG:O	1:C:102:ARG:NH1	2.35	0.57
1:C:57:ASN:HD22	1:C:58:THR:N	2.04	0.56
1:D:154:GLU:HG3	1:D:396:ARG:HG3	1.86	0.56
1:B:136:ILE:O	1:D:70:LYS:HE3	2.05	0.56
1:A:216:LEU:HD13	1:A:245:LEU:HD11	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:98:THR:HA	1:C:106:ILE:HD13	1.88	0.56
1:C:351:GLU:OE1	1:C:391:ARG:NH2	2.31	0.55
1:B:182:LEU:HD13	1:B:193:LEU:HD21	1.88	0.55
1:B:296:ASP:HB3	1:C:262:THR:HG21	1.88	0.55
1:A:133:ASN:HB3	1:A:136:ILE:HG12	1.89	0.55
1:C:264:LYS:H	1:C:264:LYS:CE	2.18	0.55
1:C:70:LYS:HA	1:C:73:ARG:HD3	1.88	0.55
1:D:77:ARG:HG2	1:D:100:TYR:HB3	1.89	0.55
1:C:17:ASN:O	1:C:21:THR:OG1	2.24	0.55
1:C:287:ILE:HD12	1:C:290:GLN:HB3	1.89	0.55
1:C:73:ARG:HG2	1:C:100:TYR:HE2	1.71	0.54
1:A:17:ASN:HD22	2:X:4:A:N6	2.06	0.54
1:A:136:ILE:HG13	1:A:137:GLY:N	2.22	0.54
1:A:180:ILE:HG21	1:A:356:LEU:HD13	1.89	0.54
1:A:378:LYS:HZ2	1:A:398:VAL:HG21	1.72	0.54
1:D:98:THR:HA	1:D:106:ILE:HD12	1.89	0.54
1:D:202:GLU:O	1:D:205:LYS:HG2	2.07	0.54
1:A:73:ARG:HD3	1:A:100:TYR:CZ	2.43	0.53
1:B:288:THR:HG21	1:D:246:SER:OG	2.08	0.53
1:B:193:LEU:HD13	1:B:199:TYR:CE2	2.44	0.53
1:B:163:ARG:NH1	3:B:502:HOH:O	2.35	0.53
1:B:37:GLU:HA	1:B:72:TRP:HH2	1.75	0.52
1:B:288:THR:OG1	1:D:248:LYS:HE2	2.10	0.52
1:B:187:GLN:HE22	1:B:271:LYS:NZ	2.08	0.52
1:D:17:ASN:HD22	2:Y:1:A:N6	2.07	0.52
1:D:115:ARG:NH1	3:D:503:HOH:O	2.42	0.52
1:C:193:LEU:HD13	1:C:199:TYR:HE2	1.74	0.52
1:B:223:VAL:HG12	1:B:231:LEU:HD11	1.91	0.52
1:D:262:THR:HG21	1:A:296:ASP:HB3	1.91	0.52
1:A:5:ALA:HB1	1:A:35:PHE:HE1	1.74	0.52
1:D:118:TYR:HE1	2:Y:1:A:H5"	1.75	0.52
1:B:84:PHE:HB3	1:B:106:ILE:HD12	1.92	0.52
1:C:20:LEU:HD21	1:C:47:MET:CE	2.41	0.51
1:C:360:ASN:ND2	1:C:370:ILE:HD11	2.25	0.51
1:B:360:ASN:HA	1:B:363:ARG:HG3	1.91	0.51
1:C:127:GLU:O	1:C:130:LYS:NZ	2.38	0.51
1:B:7:LEU:HD13	1:B:93:MET:HB2	1.93	0.51
1:C:248:LYS:HB3	1:C:287:ILE:HD11	1.91	0.51
1:A:98:THR:HG21	1:C:95:LEU:HD11	1.92	0.50
1:D:65:ILE:HG13	1:D:75:LYS:HD3	1.92	0.50
1:C:264:LYS:H	1:C:264:LYS:HE2	1.76	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:94:ALA:O	1:A:98:THR:HG23	2.12	0.50
1:D:81:ILE:O	1:D:103:ALA:HA	2.11	0.50
1:D:98:THR:HG22	1:D:106:ILE:HG21	1.93	0.49
1:A:270:VAL:HG13	1:A:358:PHE:HE1	1.77	0.49
1:C:144:PRO:HG3	1:C:171:LEU:HB3	1.95	0.49
1:A:136:ILE:HD12	1:A:138:ASP:HB2	1.93	0.49
1:D:45:LYS:HE3	1:C:44:ILE:HG22	1.94	0.49
1:D:387:HIS:HE1	1:D:389:GLN:HB3	1.78	0.49
1:B:149:ASN:N	1:B:400:LYS:HZ1	2.09	0.49
1:A:270:VAL:HG13	1:A:358:PHE:CE1	2.48	0.49
1:D:263:GLN:HG2	1:D:264:LYS:N	2.28	0.48
1:D:404:TYR:HA	1:D:407:ILE:HD12	1.95	0.48
1:D:252:ASN:ND2	1:D:291:VAL:HG12	2.29	0.48
1:D:22:GLU:O	1:D:25:THR:OG1	2.30	0.48
1:D:220:ASP:HB3	1:D:299:LEU:HD22	1.95	0.48
1:D:71:LEU:H	1:D:71:LEU:HD12	1.79	0.47
1:A:370:ILE:HG22	1:A:382:LEU:HD12	1.97	0.47
1:C:17:ASN:HD22	2:X:2:A:N6	2.13	0.47
1:A:53:TYR:CD1	1:A:150:VAL:HG22	2.49	0.47
1:D:166:ILE:HG23	1:D:175:VAL:HG11	1.97	0.47
1:D:248:LYS:O	1:D:287:ILE:HD11	2.15	0.46
1:D:252:ASN:HD21	1:D:291:VAL:HG12	1.78	0.46
1:A:136:ILE:HG13	1:A:137:GLY:H	1.80	0.46
1:C:220:ASP:HB3	1:C:299:LEU:HD22	1.95	0.46
1:C:133:ASN:HB3	1:C:136:ILE:HG12	1.96	0.46
1:B:387:HIS:O	1:B:390:GLU:HG2	2.16	0.46
1:C:218:LEU:HD11	1:C:306:ILE:HD12	1.96	0.46
1:C:369:MET:HE3	1:C:370:ILE:H	1.81	0.46
1:D:227:LEU:HD12	1:D:231:LEU:HD13	1.98	0.46
1:C:1:MET:SD	1:C:2:LEU:N	2.89	0.46
1:B:216:LEU:HD13	1:B:245:LEU:HD11	1.98	0.46
1:B:406:LYS:HA	1:B:406:LYS:HD3	1.60	0.45
1:D:69:ILE:HD12	1:D:69:ILE:N	2.31	0.45
1:D:129:LEU:HD23	1:D:142:TYR:OH	2.16	0.45
1:A:84:PHE:HB3	1:A:106:ILE:HD13	1.98	0.45
1:C:351:GLU:CD	1:C:391:ARG:HH22	2.23	0.45
1:D:144:PRO:HG3	1:D:171:LEU:HB3	1.99	0.45
1:B:81:ILE:O	1:B:103:ALA:HA	2.16	0.45
1:C:45:LYS:HG3	1:C:48:GLU:HB2	1.98	0.45
1:D:73:ARG:HG3	1:D:100:TYR:CE2	2.52	0.45
1:D:249:ASN:O	1:D:253:GLU:HG3	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:50:ALA:HB1	1:A:407:ILE:HG23	1.97	0.45
1:D:174:LYS:HD3	1:D:174:LYS:HA	1.59	0.45
1:D:2:LEU:HB2	1:D:29:PRO:HA	1.98	0.45
1:D:190:GLU:HA	1:D:193:LEU:HD12	1.99	0.45
1:C:90:ARG:HD2	1:C:92:TYR:CE2	2.46	0.45
1:B:288:THR:HG21	1:D:246:SER:HG	1.82	0.45
1:D:205:LYS:NZ	1:D:206:ASP:OD1	2.50	0.44
1:C:152:GLU:HB2	1:C:406:LYS:HE3	1.99	0.44
1:D:98:THR:HA	1:D:106:ILE:CD1	2.47	0.44
1:C:4:LEU:O	1:C:32:ILE:HA	2.18	0.44
1:B:287:ILE:HD11	1:D:245:LEU:HD13	1.98	0.44
1:C:32:ILE:HG13	1:C:58:THR:HG21	2.00	0.44
1:C:154:GLU:OE2	1:C:396:ARG:NH2	2.49	0.44
1:C:225:ILE:HG21	1:C:271:LYS:HG3	1.99	0.44
1:B:417:THR:O	1:B:417:THR:OG1	2.36	0.44
1:D:58:THR:OG1	1:D:59:LEU:N	2.50	0.44
1:D:387:HIS:CE1	1:D:389:GLN:HB3	2.52	0.44
1:C:44:ILE:HD11	1:C:414:ILE:HG12	1.99	0.44
1:D:306:ILE:HG22	1:D:314:VAL:HG21	1.98	0.44
1:A:133:ASN:HB3	1:A:136:ILE:CG1	2.48	0.43
1:A:258:THR:HA	1:A:269:LYS:HE3	1.99	0.43
1:B:297:MET:HE1	1:B:330:LYS:HE3	2.00	0.43
1:C:73:ARG:HH11	1:C:100:TYR:HE2	1.66	0.43
1:D:218:LEU:HD22	1:D:247:SER:HB2	1.99	0.43
1:B:149:ASN:H	1:B:400:LYS:HZ1	1.67	0.43
1:A:114:GLU:OE1	2:X:1:A:H2'	2.19	0.43
1:C:193:LEU:HB3	1:C:199:TYR:HD2	1.83	0.43
1:B:65:ILE:HB	1:B:72:TRP:CZ3	2.54	0.43
1:C:263:GLN:HG2	1:C:264:LYS:N	2.33	0.43
1:B:224:TYR:OH	1:B:283:HIS:HE1	2.02	0.42
1:D:94:ALA:O	1:D:98:THR:HG23	2.19	0.42
1:D:390:GLU:O	1:D:390:GLU:HG3	2.19	0.42
1:B:17:ASN:HD22	2:Y:3:A:N6	2.17	0.42
1:D:411:LEU:HA	1:D:414:ILE:HD12	2.01	0.42
1:C:256:ASN:HA	1:C:260:GLN:HG3	1.99	0.42
1:D:374:LEU:HB3	1:D:379:ILE:HG13	2.00	0.42
1:A:369:MET:HE3	1:A:369:MET:HB3	1.84	0.42
1:C:61:GLU:OE1	1:C:63:LYS:NZ	2.49	0.42
1:C:387:HIS:CE1	1:C:389:GLN:HB3	2.55	0.42
1:A:2:LEU:HA	1:A:82:ASP:OD2	2.20	0.42
1:C:321:LYS:HD3	1:C:336:ILE:HG21	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:220:ASP:OD1	1:B:223:VAL:HG23	2.19	0.42
1:D:366:ILE:H	1:D:366:ILE:HD12	1.85	0.42
1:A:5:ALA:HB1	1:A:35:PHE:CE1	2.53	0.42
1:A:214:GLY:O	1:A:243:ARG:NH2	2.47	0.42
1:A:410:ILE:O	1:A:414:ILE:HG13	2.19	0.42
1:B:36:ARG:HD3	1:B:62:GLU:HG2	2.01	0.42
1:A:27:LEU:HD21	1:A:83:ILE:HD12	2.02	0.42
1:C:99:TYR:C	1:C:100:TYR:HD1	2.27	0.42
1:B:42:LYS:C	1:B:44:ILE:H	2.28	0.42
1:B:162:LEU:O	1:B:166:ILE:HG12	2.20	0.42
1:B:369:MET:SD	1:B:383:HIS:HE1	2.43	0.42
1:D:4:LEU:O	1:D:32:ILE:HA	2.20	0.42
1:C:40:GLN:HG2	1:C:43:ASP:CG	2.45	0.42
1:D:209:LYS:O	1:D:213:ARG:HG3	2.19	0.41
1:C:107:ARG:NH2	1:C:140:ILE:HG21	2.35	0.41
1:D:113:ASP:OD2	1:D:115:ARG:NH2	2.53	0.41
1:B:4:LEU:O	1:B:32:ILE:HA	2.20	0.41
1:D:253:GLU:OE1	1:D:295:GLY:N	2.44	0.41
1:B:136:ILE:HG13	1:B:137:GLY:N	2.32	0.41
1:B:180:ILE:HD13	1:B:180:ILE:HA	1.93	0.41
1:D:238:ARG:HH11	1:D:239:LYS:HG3	1.83	0.41
1:C:387:HIS:CE1	1:C:390:GLU:HG2	2.56	0.41
1:C:387:HIS:HE1	1:C:389:GLN:HB3	1.85	0.41
1:B:267:GLU:C	1:B:269:LYS:H	2.27	0.41
1:C:70:LYS:HA	1:C:73:ARG:HB2	2.02	0.41
1:B:325:LEU:HD22	1:C:265:ILE:HD13	2.03	0.41
1:A:47:MET:HE3	1:A:50:ALA:HB3	2.03	0.41
1:B:16:ILE:HG12	1:B:47:MET:HE2	2.03	0.41
1:C:264:LYS:H	1:C:264:LYS:CD	2.33	0.40
1:C:287:ILE:HD13	1:C:287:ILE:HA	1.92	0.40
1:B:31:LYS:HG2	1:B:33:ILE:HD11	2.03	0.40
1:D:180:ILE:HG12	1:D:182:LEU:HD23	2.03	0.40
1:B:199:TYR:HE1	1:B:342:LYS:HD2	1.86	0.40
1:D:328:GLN:HE21	1:D:328:GLN:HB2	1.75	0.40
1:B:202:GLU:HG3	1:B:339:LYS:HB2	2.02	0.40
1:D:27:LEU:HD11	1:D:107:ARG:NH2	2.36	0.40
1:C:57:ASN:HD22	1:C:57:ASN:C	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	A	401/429 (94%)	398 (99%)	3 (1%)	0	100	100
1	B	408/429 (95%)	395 (97%)	13 (3%)	0	100	100
1	C	416/429 (97%)	401 (96%)	15 (4%)	0	100	100
1	D	416/429 (97%)	402 (97%)	14 (3%)	0	100	100
All	All	1641/1716 (96%)	1596 (97%)	45 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	361/382 (94%)	358 (99%)	3 (1%)	73	88
1	B	366/382 (96%)	361 (99%)	5 (1%)	59	81
1	C	373/382 (98%)	367 (98%)	6 (2%)	55	79
1	D	374/382 (98%)	365 (98%)	9 (2%)	43	70
All	All	1474/1528 (96%)	1451 (98%)	23 (2%)	55	79

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	9	SER
1	B	69	ILE

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Mol	Chain	Res	Type
1	B	107	ARG
1	B	174	LYS
1	B	390	GLU
1	D	76	ILE
1	D	86	ILE
1	D	124	VAL
1	D	155	SER
1	D	233	SER
1	D	264	LYS
1	D	288	THR
1	D	310	VAL
1	D	329	SER
1	A	86	ILE
1	A	182	LEU
1	A	306	ILE
1	C	43	ASP
1	C	96	SER
1	C	124	VAL
1	C	154	GLU
1	C	155	SER
1	C	291	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (29) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	17	ASN
1	B	30	ASN
1	B	257	HIS
1	B	283	HIS
1	B	360	ASN
1	B	403	ASN
1	D	17	ASN
1	D	128	GLN
1	D	133	ASN
1	D	257	HIS
1	D	290	GLN
1	D	328	GLN
1	D	389	GLN
1	A	17	ASN
1	A	28	ASN
1	A	57	ASN
1	A	184	ASN

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Mol	Chain	Res	Type
1	A	328	GLN
1	A	364	ASN
1	C	17	ASN
1	C	30	ASN
1	C	40	GLN
1	C	57	ASN
1	C	187	GLN
1	C	229	ASN
1	C	237	ASN
1	C	260	GLN
1	C	283	HIS
1	C	389	GLN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	X	3/4 (75%)	3 (100%)	0
2	Y	3/4 (75%)	3 (100%)	0
All	All	6/8 (75%)	6 (100%)	0

All (6) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	Y	2	A
2	Y	3	A
2	Y	4	A
2	X	2	A
2	X	3	A
2	X	4	A

There are no RNA pucker outliers to report.

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 [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

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	407/429 (94%)	0.08	7 (1%) 69 65	33, 60, 108, 162	0
1	B	412/429 (96%)	0.12	9 (2%) 62 58	30, 59, 115, 159	0
1	C	418/429 (97%)	0.10	8 (1%) 66 62	34, 60, 99, 145	0
1	D	418/429 (97%)	0.26	9 (2%) 62 58	34, 67, 114, 142	0
2	X	4/4 (100%)	-0.39	0 100 100	50, 53, 56, 62	0
2	Y	4/4 (100%)	-0.39	0 100 100	53, 62, 62, 67	0
All	All	1663/1724 (96%)	0.14	33 (1%) 65 61	30, 62, 112, 162	0

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	418	ALA	4.0
1	B	418	ALA	3.8
1	C	291	VAL	3.8
1	D	295	GLY	3.6
1	D	291	VAL	3.5
1	D	418	ALA	3.3
1	C	418	ALA	3.3
1	C	295	GLY	3.1
1	C	288	THR	3.0
1	B	44	ILE	2.9
1	D	99	TYR	2.9
1	A	99	TYR	2.7
1	D	44	ILE	2.6
1	B	268	ASN	2.6
1	D	134	VAL	2.6
1	B	365	GLY	2.5
1	D	362	TYR	2.4
1	B	39	PRO	2.4
1	D	237	ASN	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	81	ILE	2.3
1	A	2	LEU	2.3
1	A	259	ALA	2.2
1	B	43	ASP	2.2
1	A	64	VAL	2.2
1	C	46	GLY	2.2
1	C	100	TYR	2.2
1	B	134	VAL	2.1
1	C	289	SER	2.1
1	B	366	ILE	2.1
1	D	65	ILE	2.1
1	C	287	ILE	2.1
1	A	366	ILE	2.0
1	A	182	LEU	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.