



wwPDB X-ray Structure Validation Summary Report ⓘ

Jan 31, 2016 – 08:26 PM GMT

PDB ID : 1K9A
Title : Crystal structure analysis of full-length carboxyl-terminal Src kinase at 2.5 Å resolution
Authors : Ogawa, A.; Takayama, Y.; Nagata, A.; Chong, K.T.; Takeuchi, S.; Sakai, H.; Nakagawa, A.; Nada, S.; Okada, M.; Tsukihara, T.
Deposited on : 2001-10-28
Resolution : 2.50 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

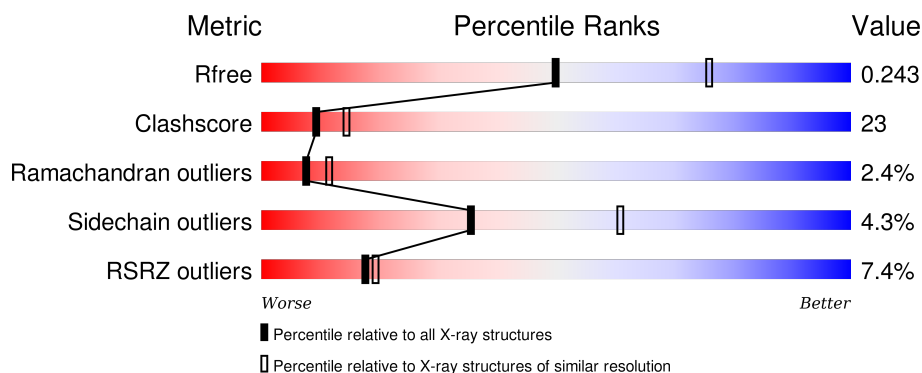
1 Overall quality at a glance ⓘ

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}	91344	3553 (2.50-2.50)
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	450	<div> <div>2%</div> <div>67% 28%</div> <div>..</div> </div>
1	B	450	<div> <div>2%</div> <div>66% 28%</div> <div>..</div> </div>
1	C	450	<div> <div>10%</div> <div>64% 30%</div> <div>..</div> </div>
1	D	450	<div> <div>9%</div> <div>60% 32%</div> <div>..</div> </div>
1	E	450	<div> <div>4%</div> <div>55% 38%</div> <div>..</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	450	<div><div></div><div>17%</div><div>47%</div><div>46%</div><div></div><div></div></div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 21313 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 Carboxyl-terminal Src kinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	439	Total	C	N	O	S	0	0	0
			3491	2231	595	644	21			
1	B	441	Total	C	N	O	S	0	0	0
			3506	2241	598	646	21			
1	C	438	Total	C	N	O	S	0	0	0
			3482	2225	593	643	21			
1	D	433	Total	C	N	O	S	0	0	0
			3443	2201	587	634	21			
1	E	440	Total	C	N	O	S	0	0	0
			3495	2232	595	647	21			
1	F	436	Total	C	N	O	S	0	0	0
			3468	2216	591	640	21			

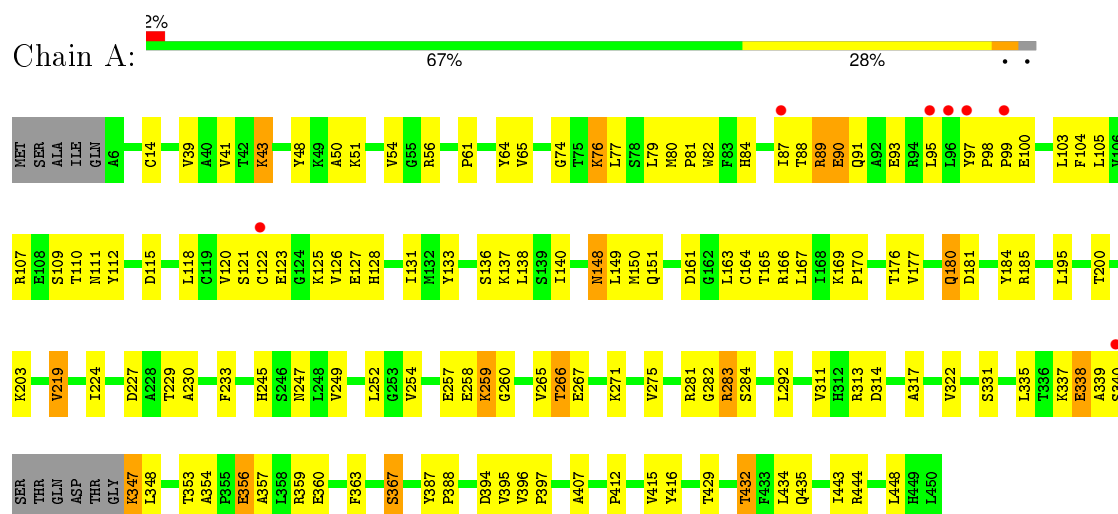
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	106	Total	O	0	0
			106	106		
2	B	89	Total	O	0	0
			89	89		
2	C	79	Total	O	0	0
			79	79		
2	D	85	Total	O	0	0
			85	85		
2	E	40	Total	O	0	0
			40	40		
2	F	29	Total	O	0	0
			29	29		

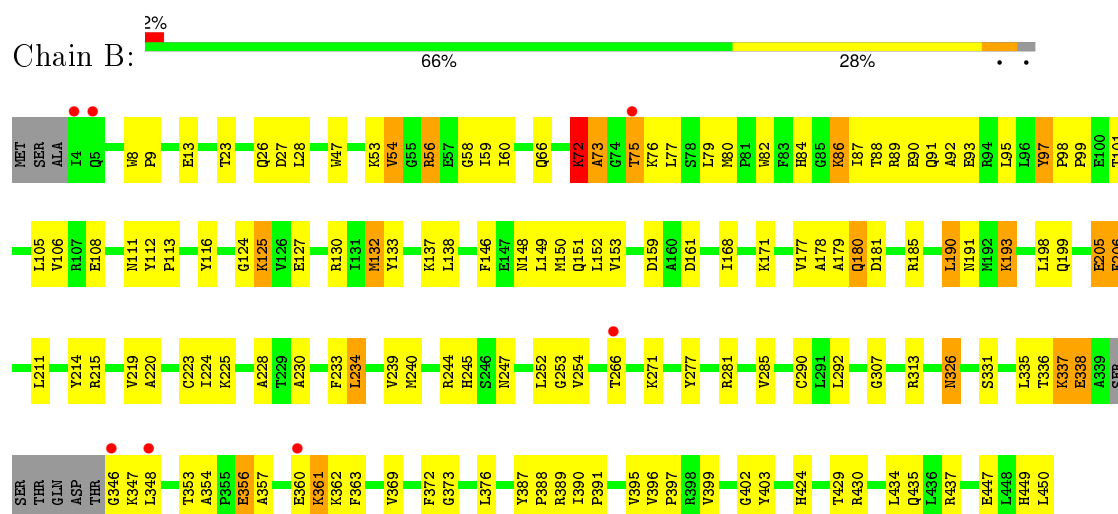
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: Carboxyl-terminal Src kinase

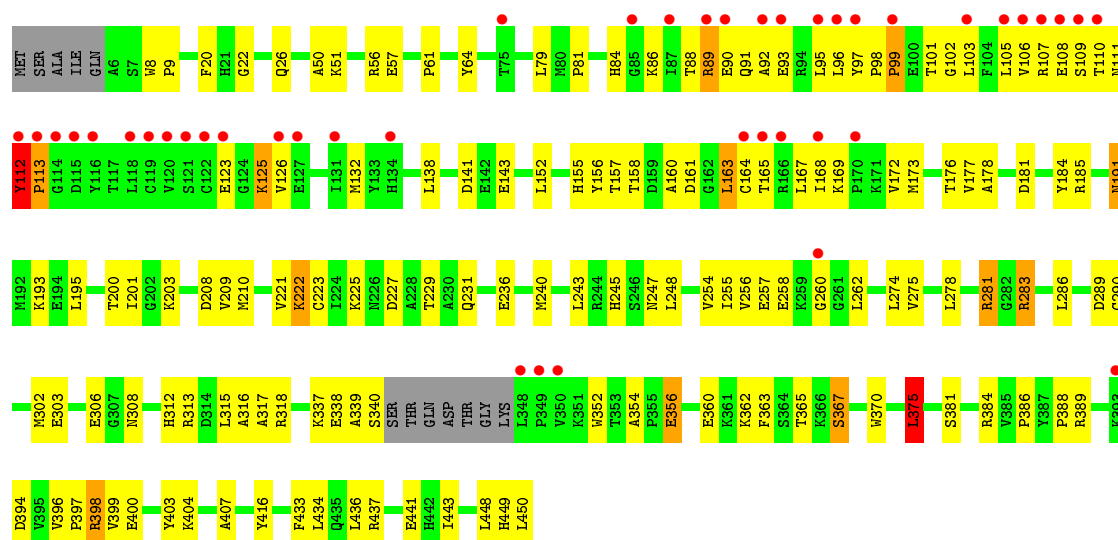


• Molecule 1: Carboxyl-terminal Src kinase

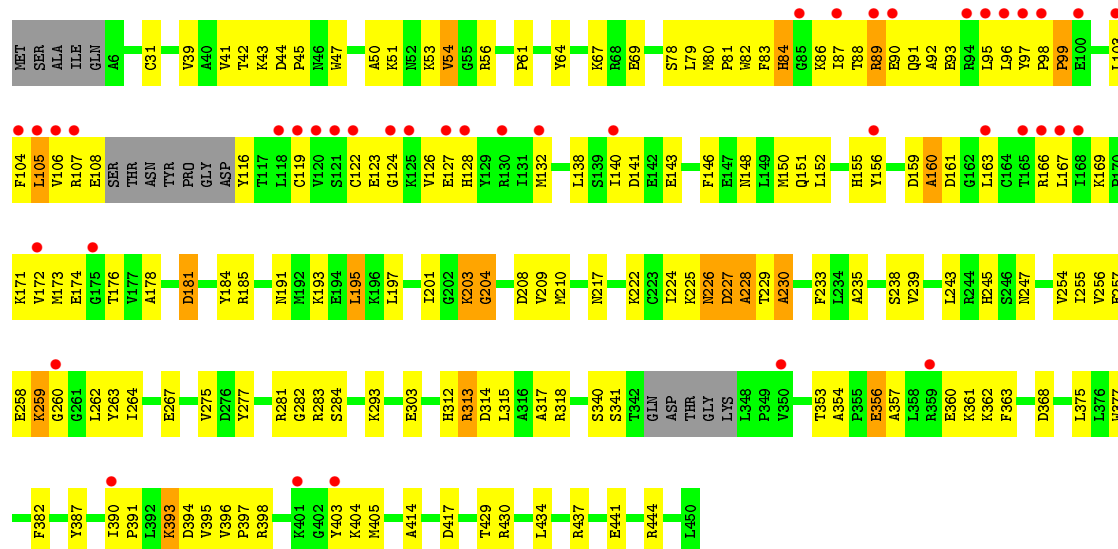


• Molecule 1: Carboxyl-terminal Src kinase

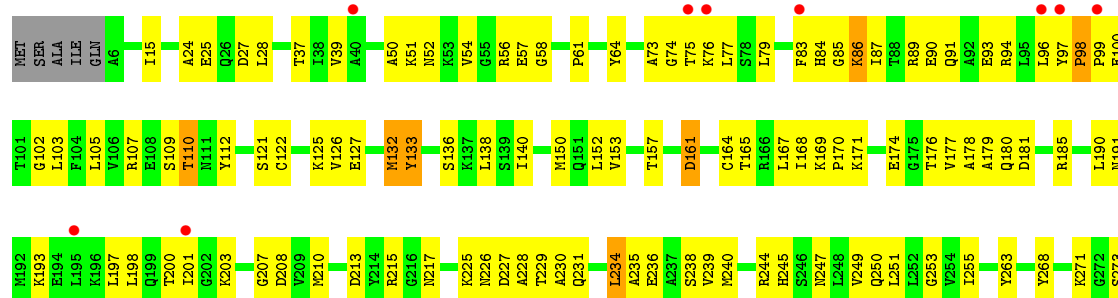


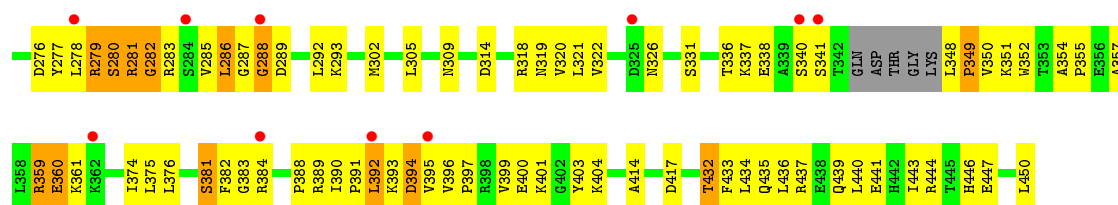


• Molecule 1: Carboxyl-terminal Src kinase

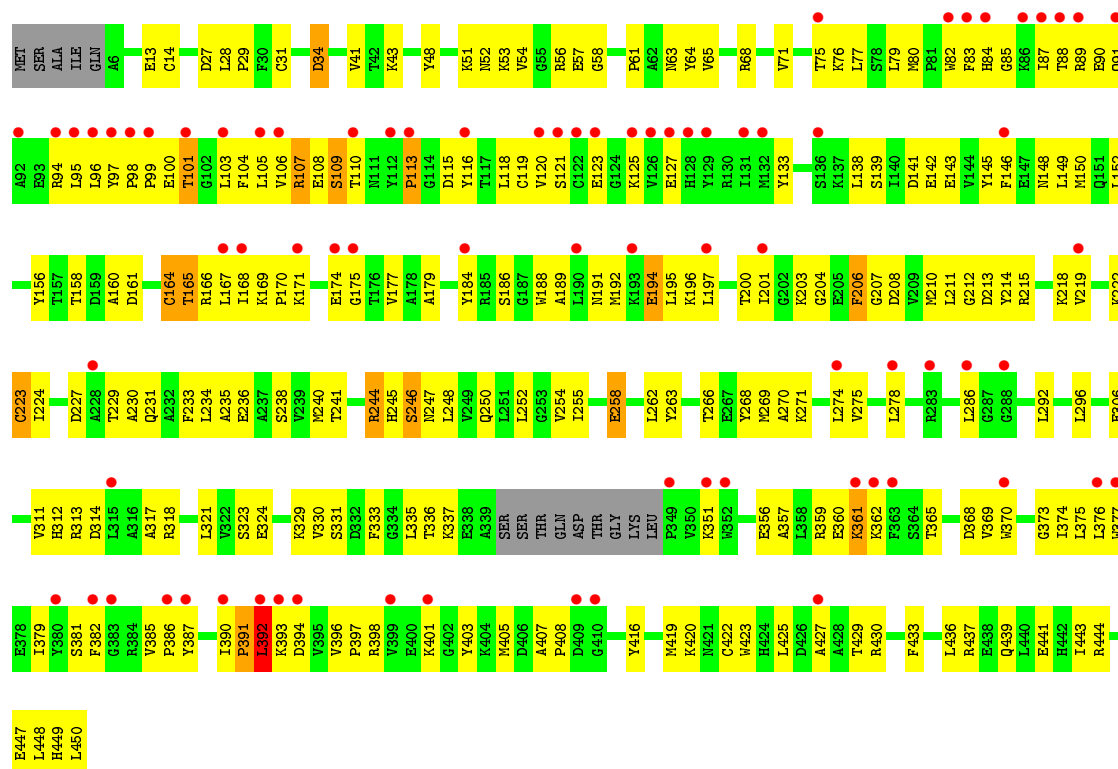


• Molecule 1: Carboxyl-terminal Src kinase





● Molecule 1: Carboxyl-terminal Src kinase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	115.00Å 162.60Å 232.40Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	73.50 – 2.50 73.03 – 2.48	Depositor EDS
% Data completeness (in resolution range)	(Not available) (73.50-2.50) 97.5 (73.03-2.48)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.35 (at 2.48Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.246 , 0.288 0.246 , 0.243	Depositor DCC
R_{free} test set	2959 reflections (2.00%)	DCC
Wilson B-factor (Å ²)	49.8	Xtriage
Anisotropy	0.650	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 54.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	1 of 150364 reflections (0.001%)	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	21313	wwPDB-VP
Average B, all atoms (Å ²)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.50% 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, 0.375 respectively for untwinned datasets, and 0.333, 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	0.41	0/3570	0.63	0/4828
1	B	0.39	0/3585	0.67	2/4848 (0.0%)
1	C	0.36	0/3561	0.61	1/4817 (0.0%)
1	D	0.38	0/3519	0.62	0/4757
1	E	0.37	0/3574	0.62	1/4835 (0.0%)
1	F	0.34	0/3547	0.58	0/4797
All	All	0.38	0/21356	0.62	4/28882 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	381	SER	N-CA-C	-6.43	93.63	111.00
1	B	72	LYS	N-CA-C	6.16	127.62	111.00
1	B	331	SER	N-CA-C	5.77	126.59	111.00
1	C	375	LEU	CA-CB-CG	5.18	127.20	115.30

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	3491	0	3484	137	0
1	B	3506	0	3501	127	0
1	C	3482	0	3471	147	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	3443	0	3441	164	0
1	E	3495	0	3483	168	0
1	F	3468	0	3456	222	0
2	A	106	0	0	3	0
2	B	89	0	0	4	0
2	C	79	0	0	6	0
2	D	85	0	0	5	0
2	E	40	0	0	1	0
2	F	29	0	0	3	0
All	All	21313	0	20836	952	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 23.

The worst 5 of 952 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:259:LYS:H	1:A:259:LYS:HD2	1.02	1.13
1:B:193:LYS:H	1:B:193:LYS:HD2	1.11	1.12
1:B:73:ALA:HB3	1:B:77:LEU:HB2	1.32	1.09
1:E:432:THR:HG22	1:E:435:GLN:H	1.18	1.06
1:A:432:THR:HG22	1:A:435:GLN:H	1.12	1.05

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	435/450 (97%)	403 (93%)	29 (7%)	3 (1%)	26	46
1	B	437/450 (97%)	405 (93%)	23 (5%)	9 (2%)	9	14

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	434/450 (96%)	386 (89%)	43 (10%)	5 (1%)	16	29
1	D	427/450 (95%)	376 (88%)	39 (9%)	12 (3%)	6	9
1	E	436/450 (97%)	388 (89%)	33 (8%)	15 (3%)	5	6
1	F	432/450 (96%)	351 (81%)	63 (15%)	18 (4%)	3	4
All	All	2601/2700 (96%)	2309 (89%)	230 (9%)	62 (2%)	7	11

5 of 62 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	112	TYR
1	D	123	GLU
1	D	204	GLY
1	D	228	ALA
1	D	230	ALA

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	376/385 (98%)	356 (95%)	20 (5%)	28	50
1	B	377/385 (98%)	359 (95%)	18 (5%)	31	55
1	C	375/385 (97%)	360 (96%)	15 (4%)	38	64
1	D	371/385 (96%)	357 (96%)	14 (4%)	40	67
1	E	377/385 (98%)	364 (97%)	13 (3%)	44	72
1	F	373/385 (97%)	357 (96%)	16 (4%)	35	61
All	All	2249/2310 (97%)	2153 (96%)	96 (4%)	35	61

5 of 96 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	191	ASN
1	D	54	VAL

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Mol	Chain	Res	Type
1	F	194	GLU
1	C	195	LEU
1	C	289	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 47 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	308	ASN
1	D	247	ASN
1	F	247	ASN
1	C	446	HIS
1	D	312	HIS

5.3.3 RNA [i](#)

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 carbohydrates 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 [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	439/450 (97%)	0.36	7 (1%) 74 78	23, 45, 79, 100	0
1	B	441/450 (98%)	0.43	7 (1%) 74 78	29, 48, 78, 95	0
1	C	438/450 (97%)	0.68	43 (9%) 10 10	31, 52, 111, 124	0
1	D	433/450 (96%)	0.63	41 (9%) 10 11	24, 50, 101, 118	0
1	E	440/450 (97%)	0.53	19 (4%) 39 44	31, 58, 91, 107	0
1	F	436/450 (96%)	1.12	78 (17%) 2 2	40, 74, 118, 136	0
All	All	2627/2700 (97%)	0.62	195 (7%) 17 19	23, 54, 103, 136	0

The worst 5 of 195 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	103	LEU	8.1
1	F	97	TYR	7.9
1	F	122	CYS	6.9
1	D	97	TYR	6.7
1	C	112	TYR	6.6

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 carbohydrates 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.