



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 07:03 pm GMT

PDB ID : 4J7B
Title : Crystal structure of polo-like kinase 1
Authors : Xu, J.; Shen, C.; Quan, J.; Wang, T.
Deposited on : 2013-02-13
Resolution : 2.30 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

<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
Xtriage (Phenix)	:	1.9-1692
EDS	:	trunk28620
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
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)	:	recalc28949

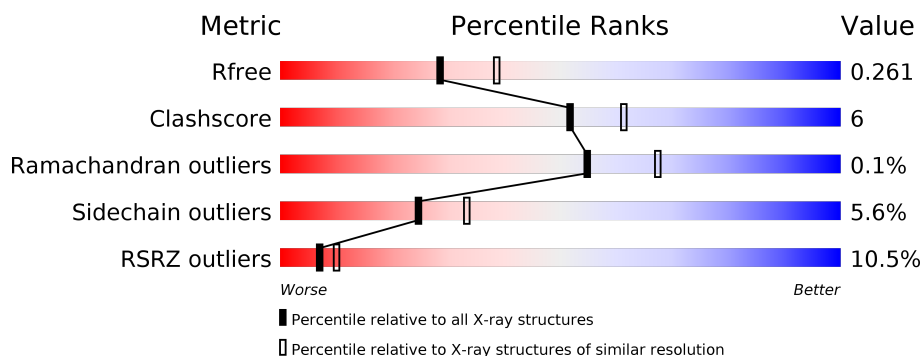
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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}	100719	4130 (2.30-2.30)
Clashscore	112137	4751 (2.30-2.30)
Ramachandran outliers	110173	4705 (2.30-2.30)
Sidechain outliers	110143	4704 (2.30-2.30)
RSRZ outliers	101464	4156 (2.30-2.30)

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	297	<div> <div>9%</div> <div>81% 13% . .</div> </div>
1	D	297	<div> <div>16%</div> <div>78% 15% . 5%</div> </div>
2	B	237	<div> <div>4%</div> <div>76% 13% . 10%</div> </div>
2	E	237	<div> <div>5%</div> <div>75% 14% . 10%</div> </div>
3	C	58	<div> <div>14%</div> <div>55% 28% . 16%</div> </div>
3	F	58	<div> <div>19%</div> <div>64% 19% . 16%</div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9170 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 Polo-like kinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	284	Total	C	N	O	S	0	0	0
			2287	1470	401	400	16			
1	D	281	Total	C	N	O	S	0	0	0
			2265	1456	397	397	15			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	16	MET	-	EXPRESSION TAG	UNP Q6DRK7
A	17	GLY	-	EXPRESSION TAG	UNP Q6DRK7
A	162	ASN	ASP	ENGINEERED MUTATION	UNP Q6DRK7
A	192	PHE	GLU	ENGINEERED MUTATION	UNP Q6DRK7
D	16	MET	-	EXPRESSION TAG	UNP Q6DRK7
D	17	GLY	-	EXPRESSION TAG	UNP Q6DRK7
D	162	ASN	ASP	ENGINEERED MUTATION	UNP Q6DRK7
D	192	PHE	GLU	ENGINEERED MUTATION	UNP Q6DRK7

- Molecule 2 is a protein called Polo-like kinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	213	Total	C	N	O	S	0	0	0
			1737	1109	295	320	13			
2	E	213	Total	C	N	O	S	0	0	0
			1737	1109	295	320	13			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	359	MET	-	EXPRESSION TAG	UNP Q6DRK7
B	384	GLY	ASP	ENGINEERED MUTATION	UNP Q6DRK7
E	359	MET	-	EXPRESSION TAG	UNP Q6DRK7
E	384	GLY	ASP	ENGINEERED MUTATION	UNP Q6DRK7

- Molecule 3 is a protein called 205 kDa microtubule-associated protein.

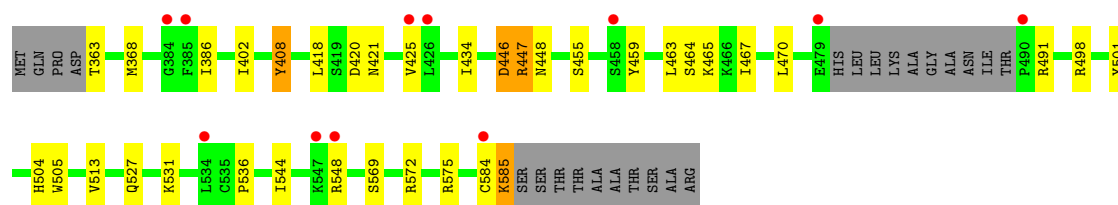
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	49	Total	C	N	O	S	0	0	0
			392	240	66	84	2			
3	F	49	Total	C	N	O	S	0	0	0
			392	240	66	84	2			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	268	MET	-	EXPRESSION TAG	UNP P23226
C	269	GLY	-	EXPRESSION TAG	UNP P23226
C	270	HIS	-	EXPRESSION TAG	UNP P23226
C	271	HIS	-	EXPRESSION TAG	UNP P23226
C	272	HIS	-	EXPRESSION TAG	UNP P23226
C	273	HIS	-	EXPRESSION TAG	UNP P23226
C	274	HIS	-	EXPRESSION TAG	UNP P23226
C	275	HIS	-	EXPRESSION TAG	UNP P23226
F	268	MET	-	EXPRESSION TAG	UNP P23226
F	269	GLY	-	EXPRESSION TAG	UNP P23226
F	270	HIS	-	EXPRESSION TAG	UNP P23226
F	271	HIS	-	EXPRESSION TAG	UNP P23226
F	272	HIS	-	EXPRESSION TAG	UNP P23226
F	273	HIS	-	EXPRESSION TAG	UNP P23226
F	274	HIS	-	EXPRESSION TAG	UNP P23226
F	275	HIS	-	EXPRESSION TAG	UNP P23226

- Molecule 4 is water.

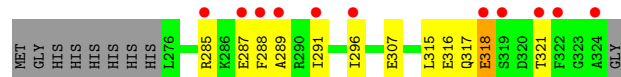
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	94	Total	O	0	0
			94	94		
4	B	128	Total	O	0	0
			128	128		
4	C	17	Total	O	0	0
			17	17		
4	D	54	Total	O	0	0
			54	54		
4	E	56	Total	O	0	0
			56	56		
4	F	11	Total	O	0	0
			11	11		



• Molecule 3: 205 kDa microtubule-associated protein



• Molecule 3: 205 kDa microtubule-associated protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	47.68Å 57.45Å 125.74Å 89.15° 89.84° 72.38°	Depositor
Resolution (Å)	25.00 – 2.30 24.99 – 2.30	Depositor EDS
% Data completeness (in resolution range)	93.9 (25.00-2.30) 93.7 (24.99-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.58 (at 2.31Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.205 , 0.264 0.204 , 0.261	Depositor DCC
R_{free} test set	2702 reflections (5.35%)	DCC
Wilson B-factor (Å ²)	33.6	Xtriage
Anisotropy	0.619	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 45.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.010 for -h,-k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9170	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.26% 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

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	0.78	0/2342	0.83	4/3161 (0.1%)
1	D	0.64	0/2319	0.77	1/3131 (0.0%)
2	B	0.99	2/1774 (0.1%)	1.03	8/2393 (0.3%)
2	E	0.80	0/1774	0.86	2/2393 (0.1%)
3	C	0.81	0/398	0.98	4/536 (0.7%)
3	F	0.67	0/398	0.84	0/536
All	All	0.79	2/9005 (0.0%)	0.87	19/12150 (0.2%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	422	SER	CB-OG	6.09	1.50	1.42
2	B	392	GLU	CD-OE1	5.37	1.31	1.25

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	310	ARG	NE-CZ-NH1	7.45	124.03	120.30
2	B	555	MET	CG-SD-CE	-7.37	88.40	100.20
3	C	277	ASP	CB-CG-OD1	-6.26	112.66	118.30
3	C	277	ASP	CB-CG-OD2	6.22	123.90	118.30
2	E	572	ARG	NE-CZ-NH2	-6.14	117.23	120.30
2	B	393	ASP	CB-CG-OD1	5.99	123.69	118.30
2	B	432	ARG	NE-CZ-NH2	-5.97	117.32	120.30
1	A	310	ARG	NE-CZ-NH2	-5.96	117.32	120.30
3	C	311	ARG	NE-CZ-NH2	-5.89	117.36	120.30
1	D	179	GLY	N-CA-C	5.74	127.45	113.10
2	B	432	ARG	NE-CZ-NH1	5.71	123.16	120.30
1	A	269	ARG	NE-CZ-NH1	5.63	123.12	120.30
1	A	157	ARG	NE-CZ-NH2	-5.59	117.50	120.30
2	B	383	ARG	NE-CZ-NH1	5.49	123.05	120.30
2	B	420	ASP	CB-CG-OD1	5.46	123.21	118.30
2	B	438	ASP	CB-CG-OD2	-5.38	113.45	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	290	ARG	NE-CZ-NH1	5.21	122.91	120.30
2	E	420	ASP	CB-CG-OD1	5.06	122.85	118.30
2	B	572	ARG	NE-CZ-NH1	5.04	122.82	120.30

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	2287	0	2341	27	0
1	D	2265	0	2315	25	0
2	B	1737	0	1728	17	0
2	E	1737	0	1728	26	0
3	C	392	0	359	15	0
3	F	392	0	359	13	0
4	A	94	0	0	3	0
4	B	128	0	0	5	0
4	C	17	0	0	0	0
4	D	54	0	0	2	0
4	E	56	0	0	3	0
4	F	11	0	0	0	0
All	All	9170	0	8830	112	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 6.

All (112) 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:368:MET:HG3	4:B:688:HOH:O	1.45	1.14
3:F:318:GLU:HA	3:F:318:GLU:OE2	1.58	1.03
3:C:316:GLU:HG3	3:C:316:GLU:O	1.75	0.85
3:C:287:GLU:HG2	3:C:288:PHE:CE2	2.17	0.80
1:D:23:LEU:N	1:D:23:LEU:CD2	2.44	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:23:LEU:HD23	1:D:23:LEU:N	1.98	0.77
1:A:269:ARG:HH11	1:A:269:ARG:HG2	1.46	0.77
1:D:275:ASP:OD2	1:D:277:THR:HG22	1.85	0.75
2:B:403:SER:HB2	3:C:316:GLU:OE2	1.86	0.75
1:D:97:PRO:HD3	4:D:439:HOH:O	1.87	0.75
2:E:425:VAL:HG23	2:E:470:LEU:HD23	1.70	0.73
3:F:318:GLU:CA	3:F:318:GLU:OE2	2.37	0.72
2:E:491:ARG:HG2	4:E:643:HOH:O	1.90	0.71
1:A:72:LYS:NZ	1:A:108:ASP:O	2.23	0.71
3:F:316:GLU:O	3:F:316:GLU:CG	2.40	0.69
1:A:201:PRO:HG2	1:A:204:ILE:HG22	1.74	0.68
1:A:204:ILE:HD12	1:A:209:LEU:HD21	1.75	0.67
3:F:316:GLU:O	3:F:316:GLU:HG2	1.93	0.67
2:E:425:VAL:HG23	2:E:470:LEU:CD2	2.27	0.64
3:C:287:GLU:CG	3:C:288:PHE:CE2	2.81	0.64
1:A:310:ARG:O	1:A:311:PHE:C	2.35	0.62
1:A:269:ARG:NH1	1:A:269:ARG:HG2	2.12	0.62
2:E:434:ILE:HD11	2:E:501:TYR:HB3	1.80	0.62
2:E:402:ILE:HD13	2:E:418:LEU:HD23	1.83	0.61
1:D:58:ASP:HB3	1:D:63:GLU:HG2	1.83	0.61
3:F:287:GLU:HG3	3:F:288:PHE:CE2	2.35	0.61
1:A:128:HIS:HD2	4:A:425:HOH:O	1.82	0.61
2:E:504:HIS:CD2	2:E:585:LYS:HE2	2.35	0.61
2:B:453:TYR:OH	2:B:498:ARG:NE	2.34	0.60
1:D:160:HIS:O	1:D:161:ARG:HB2	2.01	0.60
1:A:108:ASP:HB2	4:A:468:HOH:O	2.00	0.59
3:C:287:GLU:HG2	3:C:288:PHE:CD2	2.37	0.59
2:E:421:ASN:ND2	4:E:636:HOH:O	2.37	0.57
3:C:288:PHE:N	3:C:288:PHE:CD2	2.74	0.56
3:C:316:GLU:O	3:C:316:GLU:CG	2.45	0.55
1:A:261:ASN:CB	1:A:262:PRO:HD2	2.37	0.54
1:A:59:MET:O	1:A:62:LYS:NZ	2.30	0.54
1:A:164:LYS:HD3	1:A:202:ASN:HB2	1.89	0.54
1:A:282:VAL:HA	1:A:285:LEU:HD22	1.89	0.53
1:A:128:HIS:HE1	4:A:404:HOH:O	1.91	0.53
1:D:63:GLU:HG3	1:D:65:PHE:CE2	2.44	0.52
2:E:434:ILE:CD1	2:E:501:TYR:HB3	2.39	0.52
2:B:369:LEU:HD22	2:B:582:MET:CE	2.40	0.52
2:E:465:LYS:NZ	3:F:296:ILE:O	2.27	0.52
3:C:302:ARG:HG3	2:E:527:GLN:HG2	1.92	0.51
1:D:140:ARG:NH1	1:D:295:TYR:O	2.43	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:243:LYS:NZ	1:A:247:ILE:HD11	2.26	0.51
1:A:143:MET:HG3	1:A:227:LEU:HD13	1.93	0.51
1:A:51:ALA:HB2	1:A:70:VAL:HG22	1.92	0.51
1:D:51:ALA:HB2	1:D:70:VAL:HG22	1.92	0.51
3:F:287:GLU:HG3	3:F:288:PHE:CZ	2.47	0.50
4:B:712:HOH:O	3:C:294:ASP:HA	2.11	0.50
2:E:447:ARG:HG2	2:E:447:ARG:HH11	1.76	0.50
2:E:434:ILE:HD11	2:E:501:TYR:CB	2.40	0.50
2:E:585:LYS:N	2:E:585:LYS:HD2	2.26	0.50
2:B:371:GLN:OE1	2:B:539:GLY:HA2	2.12	0.49
1:D:125:LEU:HD11	1:D:129:LYS:NZ	2.27	0.49
1:D:308:PRO:O	2:E:498:ARG:HD3	2.11	0.49
2:B:465:LYS:HG3	4:B:669:HOH:O	2.13	0.49
2:B:448:ASN:HB2	4:B:690:HOH:O	2.12	0.48
3:C:313:HIS:CD2	3:C:315:LEU:HD22	2.49	0.48
1:A:261:ASN:HB2	1:A:262:PRO:HD2	1.96	0.47
2:E:446:ASP:OD1	2:E:448:ASN:N	2.48	0.47
2:B:575:ARG:NH1	4:B:627:HOH:O	2.46	0.47
1:A:32:ASP:O	1:A:36:MET:N	2.47	0.47
1:D:239:THR:HG22	1:D:245:THR:OG1	2.15	0.47
1:D:58:ASP:CB	1:D:63:GLU:HG2	2.44	0.47
1:A:24:LYS:HD2	1:A:110:ASP:CG	2.35	0.47
1:D:118:ILE:HG22	4:D:416:HOH:O	2.15	0.46
2:E:584:CYS:C	2:E:585:LYS:HD2	2.36	0.46
1:D:76:LEU:HD12	1:D:77:LYS:H	1.79	0.46
2:E:447:ARG:NH1	2:E:447:ARG:HG2	2.31	0.46
1:A:308:PRO:O	2:B:498:ARG:HG2	2.16	0.46
1:D:77:LYS:HD2	1:D:79:HIS:HE1	1.81	0.46
2:B:369:LEU:HD22	2:B:582:MET:HE1	1.99	0.45
1:A:58:ASP:O	1:A:62:LYS:N	2.46	0.45
2:B:445:ILE:HD12	2:B:445:ILE:N	2.32	0.45
3:F:317:GLN:O	3:F:318:GLU:C	2.52	0.45
1:D:30:LEU:HD22	1:D:113:TYR:CE1	2.53	0.44
2:E:446:ASP:OD1	2:E:446:ASP:C	2.55	0.44
3:F:285:ARG:HD2	3:F:285:ARG:HA	1.80	0.44
1:D:242:LEU:H	1:D:242:LEU:HD12	1.81	0.44
3:C:302:ARG:O	3:C:306:ILE:HG12	2.17	0.44
2:E:447:ARG:NH2	3:F:289:ALA:HB3	2.33	0.44
2:E:425:VAL:CG2	2:E:470:LEU:HD23	2.45	0.43
1:D:148:GLN:O	1:D:151:GLN:HG3	2.18	0.43
1:A:242:LEU:HG	1:A:246:TYR:CE2	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:76:LEU:HD12	1:D:77:LYS:HG3	2.00	0.42
1:A:68:LYS:NZ	1:A:87:GLU:OE1	2.53	0.42
1:D:161:ARG:NH1	1:D:183:LEU:O	2.52	0.42
1:D:299:ARG:HD2	1:D:299:ARG:HA	1.95	0.42
1:A:269:ARG:NH1	1:A:269:ARG:CG	2.81	0.42
2:B:397:ILE:HA	2:B:398:PRO:HD3	1.88	0.42
1:D:64:VAL:HG11	1:D:120:ARG:HH21	1.84	0.42
2:E:408:TYR:OH	3:F:307:GLU:HB2	2.20	0.42
1:A:58:ASP:C	1:A:58:ASP:OD1	2.58	0.41
2:E:491:ARG:CD	4:E:643:HOH:O	2.68	0.41
3:F:315:LEU:O	3:F:317:GLN:N	2.54	0.41
1:A:285:LEU:HD12	1:A:285:LEU:HA	1.97	0.41
2:B:506:PHE:CD1	3:C:279:LEU:HD22	2.56	0.41
2:B:546:GLU:N	2:B:546:GLU:OE1	2.37	0.41
1:D:77:LYS:HD2	1:D:79:HIS:CE1	2.55	0.41
1:A:80:GLN:OE1	1:A:80:GLN:HA	2.21	0.41
2:B:505:TRP:CD1	3:C:283:SER:HB3	2.56	0.41
3:F:291:ILE:HD12	3:F:291:ILE:H	1.86	0.41
2:B:506:PHE:HB2	3:C:279:LEU:HD13	2.03	0.41
2:E:505:TRP:HA	2:E:513:VAL:O	2.21	0.41
2:B:476:TYR:OH	3:C:309:ASP:OD1	2.30	0.41
2:E:459:TYR:CE1	2:E:464:SER:HA	2.55	0.41
2:E:368:MET:CE	2:E:536:PRO:HD3	2.52	0.40
2:E:531:LYS:HB2	2:E:544:ILE:HB	2.03	0.40
1:D:244:GLU:O	1:D:248:ARG:HG3	2.22	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	280/297 (94%)	269 (96%)	10 (4%)	1 (0%)	38 47

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	277/297 (93%)	266 (96%)	11 (4%)	0	100	100
2	B	209/237 (88%)	203 (97%)	6 (3%)	0	100	100
2	E	209/237 (88%)	204 (98%)	5 (2%)	0	100	100
3	C	47/58 (81%)	46 (98%)	1 (2%)	0	100	100
3	F	47/58 (81%)	44 (94%)	3 (6%)	0	100	100
All	All	1069/1184 (90%)	1032 (96%)	36 (3%)	1 (0%)	55	67

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	262	PRO

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	255/266 (96%)	243 (95%)	12 (5%)	30	41
1	D	252/266 (95%)	235 (93%)	17 (7%)	19	24
2	B	192/210 (91%)	184 (96%)	8 (4%)	34	47
2	E	192/210 (91%)	180 (94%)	12 (6%)	21	28
3	C	42/49 (86%)	38 (90%)	4 (10%)	10	12
3	F	42/49 (86%)	40 (95%)	2 (5%)	30	40
All	All	975/1050 (93%)	920 (94%)	55 (6%)	25	33

All (55) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	LYS
1	A	79	HIS
1	A	82	GLU
1	A	94	LEU
1	A	123	SER

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Mol	Chain	Res	Type
1	A	124	LEU
1	A	240	SER
1	A	241	CYS
1	A	243	LYS
1	A	277	THR
1	A	285	LEU
1	A	310	ARG
2	B	363	THR
2	B	408	TYR
2	B	426	LEU
2	B	455	SER
2	B	479	GLU
2	B	491	ARG
2	B	498	ARG
2	B	575	ARG
3	C	285	ARG
3	C	303	GLU
3	C	315	LEU
3	C	318	GLU
1	D	20	SER
1	D	23	LEU
1	D	38	ARG
1	D	50	PHE
1	D	76	LEU
1	D	85	SER
1	D	88	ILE
1	D	92	LYS
1	D	94	LEU
1	D	105	PHE
1	D	109	ASP
1	D	123	SER
1	D	124	LEU
1	D	180	ASP
1	D	240	SER
1	D	244	GLU
1	D	285	LEU
2	E	363	THR
2	E	386	ILE
2	E	408	TYR
2	E	446	ASP
2	E	447	ARG
2	E	455	SER

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Mol	Chain	Res	Type
2	E	463	LEU
2	E	467	ILE
2	E	548	ARG
2	E	569	SER
2	E	575	ARG
2	E	585	LYS
3	F	318	GLU
3	F	321	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	128	HIS
1	D	79	HIS
2	E	504	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 ⓘ

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	284/297 (95%)	0.40	26 (9%) 10 13	22, 51, 92, 105	0
1	D	281/297 (94%)	0.74	48 (17%) 2 2	36, 63, 104, 133	0
2	B	213/237 (89%)	0.11	10 (4%) 32 39	18, 31, 62, 82	0
2	E	213/237 (89%)	0.23	11 (5%) 28 35	25, 42, 77, 109	0
3	C	49/58 (84%)	0.78	8 (16%) 2 3	30, 51, 95, 114	0
3	F	49/58 (84%)	1.16	11 (22%) 1 1	30, 61, 127, 159	0
All	All	1089/1184 (91%)	0.45	114 (10%) 7 10	18, 48, 94, 159	0

All (114) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	78	PRO	6.8
3	C	319	SER	5.9
3	F	288	PHE	5.8
3	F	318	GLU	5.6
3	C	288	PHE	5.6
3	F	319	SER	5.4
3	F	289	ALA	5.3
3	C	318	GLU	5.2
1	D	199	GLY	5.1
1	D	50	PHE	4.8
1	A	200	THR	4.5
1	D	293	SER	4.5
1	D	242	LEU	4.5
1	D	77	LYS	4.4
2	E	490	PRO	4.4
1	A	198	CYS	4.3
1	D	35	THR	4.1
3	F	287	GLU	4.1
3	F	291	ILE	4.1

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Mol	Chain	Res	Type	RSRZ
1	D	200	THR	4.0
1	D	22	PRO	3.9
1	D	36	MET	3.9
1	D	295	TYR	3.9
1	A	246	TYR	3.8
1	D	223	LEU	3.8
3	C	287	GLU	3.7
1	D	76	LEU	3.6
1	D	241	CYS	3.6
1	A	18	PRO	3.5
3	C	289	ALA	3.4
1	A	213	GLY	3.4
2	B	584	CYS	3.3
1	D	20	SER	3.3
1	A	34	ARG	3.3
1	A	226	ILE	3.3
1	D	114	VAL	3.2
1	D	240	SER	3.1
1	D	21	ALA	3.1
1	D	226	ILE	3.1
1	D	79	HIS	3.0
1	A	223	LEU	3.0
1	A	255	SER	3.0
1	A	242	LEU	3.0
3	F	322	PHE	3.0
3	F	321	THR	3.0
3	F	324	ALA	3.0
1	D	262	PRO	2.9
2	B	548	ARG	2.9
1	D	311	PHE	2.8
1	A	222	SER	2.7
1	D	75	LEU	2.7
1	A	78	PRO	2.7
1	D	187	ILE	2.7
3	F	296	ILE	2.7
1	D	157	ARG	2.7
1	D	222	SER	2.7
1	A	244	GLU	2.6
2	E	458	SER	2.6
1	D	82	GLU	2.6
1	D	294	GLY	2.6
1	D	289	GLU	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	258	ARG	2.6
2	B	583	ALA	2.6
1	D	258	ARG	2.5
1	A	60	ASP	2.5
1	D	243	LYS	2.5
2	E	547	LYS	2.5
3	C	285	ARG	2.5
1	D	48	GLY	2.5
1	D	255	SER	2.5
2	B	490	PRO	2.5
2	E	426	LEU	2.5
3	F	285	ARG	2.4
1	A	77	LYS	2.4
1	D	74	MET	2.4
2	E	479	GLU	2.4
2	E	584	CYS	2.4
1	D	173	ASP	2.4
1	A	35	THR	2.4
2	E	384	GLY	2.3
2	E	534	LEU	2.3
1	D	59	MET	2.3
2	E	548	ARG	2.3
1	A	75	LEU	2.3
1	D	221	TRP	2.3
1	D	28	ASP	2.2
1	A	243	LYS	2.2
1	D	212	LYS	2.2
1	A	204	ILE	2.2
3	C	321	THR	2.2
2	B	534	LEU	2.2
1	A	199	GLY	2.2
1	D	211	LYS	2.2
2	B	547	LYS	2.2
1	D	225	CYS	2.2
1	A	82	GLU	2.2
1	A	36	MET	2.2
1	D	40	MET	2.2
1	D	60	ASP	2.2
1	D	186	LYS	2.1
2	B	532	LEU	2.1
1	A	293	SER	2.1
1	D	213	GLY	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	114	VAL	2.1
1	D	246	TYR	2.1
2	E	425	VAL	2.1
1	D	244	GLU	2.1
3	C	286	LYS	2.1
1	A	187	ILE	2.1
2	B	514	LEU	2.1
1	D	172	ASP	2.1
2	B	425	VAL	2.0
2	B	533	ILE	2.0
2	E	385	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 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.