



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 5, 2022 – 10:06 am BST

PDB ID : 7O1F
Title : PCNA from *Chaetomium thermophilum* in complex with PolD4 PIP peptide
Authors : Alphey, M.A.; MacNeill, S.; Yang, D.
Deposited on : 2021-03-29
Resolution : 2.45 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.27
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0267
CCP4 : 7.1.010 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.27

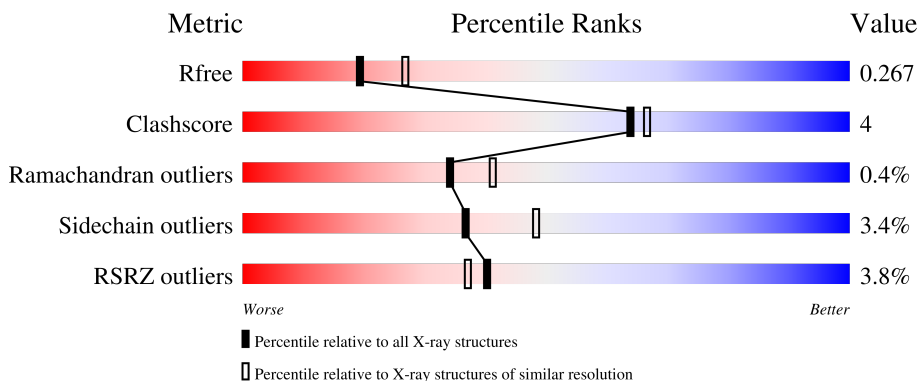
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.45 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.



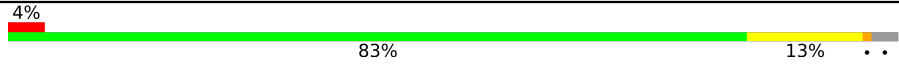
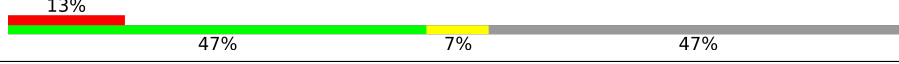

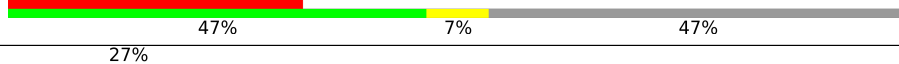
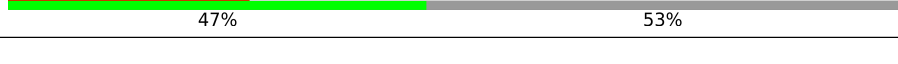
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1544 (2.48-2.44)
Clashscore	141614	1613 (2.48-2.44)
Ramachandran outliers	138981	1598 (2.48-2.44)
Sidechain outliers	138945	1598 (2.48-2.44)
RSRZ outliers	127900	1523 (2.48-2.44)

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	261	 2% 87% 10%
1	B	261	 % 83% 12%
1	C	261	 5% 77% 14% 9%
1	D	261	 3% 84% 12%
1	E	261	 3% 83% 10% 6%

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Mol	Chain	Length	Quality of chain
1	F	261	
2	J	15	
2	K	15	
2	M	15	
2	O	15	

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 11613 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 Proliferating cell nuclear antigen.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	253	Total 1889	C 1195	N 312	O 369	S 13	0	0	0
1	B	250	Total 1884	C 1187	N 309	O 375	S 13	0	0	0
1	C	238	Total 1774	C 1116	N 293	O 352	S 13	0	0	0
1	D	254	Total 1894	C 1195	N 314	O 372	S 13	0	0	0
1	E	246	Total 1852	C 1169	N 306	O 364	S 13	0	0	0
1	F	253	Total 1899	C 1197	N 315	O 374	S 13	0	0	0

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	ALA	-	expression tag	UNP G0SF70
A	0	MET	-	expression tag	UNP G0SF70
A	1	ALA	-	expression tag	UNP G0SF70
B	-1	ALA	-	expression tag	UNP G0SF70
B	0	MET	-	expression tag	UNP G0SF70
B	1	ALA	-	expression tag	UNP G0SF70
C	-1	ALA	-	expression tag	UNP G0SF70
C	0	MET	-	expression tag	UNP G0SF70
C	1	ALA	-	expression tag	UNP G0SF70
D	-1	ALA	-	expression tag	UNP G0SF70
D	0	MET	-	expression tag	UNP G0SF70
D	1	ALA	-	expression tag	UNP G0SF70
E	-1	ALA	-	expression tag	UNP G0SF70
E	0	MET	-	expression tag	UNP G0SF70
E	1	ALA	-	expression tag	UNP G0SF70
F	-1	ALA	-	expression tag	UNP G0SF70
F	0	MET	-	expression tag	UNP G0SF70

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Chain	Residue	Modelled	Actual	Comment	Reference
F	1	ALA	-	expression tag	UNP G0SF70

- Molecule 2 is a protein called Peptide fragment from PolD4.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	J	8	Total	C	N	O	0	0	0
			64	40	12	12			
2	K	8	Total	C	N	O	0	0	0
			64	40	12	12			
2	M	8	Total	C	N	O	0	0	0
			64	40	12	12			
2	O	7	Total	C	N	O	0	0	0
			51	33	8	10			

- Molecule 3 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	3	Total	Na	0	0
			3	3		
3	B	2	Total	Na	0	0
			2	2		
3	C	3	Total	Na	0	0
			3	3		
3	D	4	Total	Na	0	0
			4	4		
3	E	4	Total	Na	0	0
			4	4		
3	F	2	Total	Na	0	0
			2	2		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	24	Total	O	0	0
			24	24		
4	B	32	Total	O	0	0
			32	32		
4	C	27	Total	O	0	0
			27	27		
4	D	34	Total	O	0	0
			34	34		

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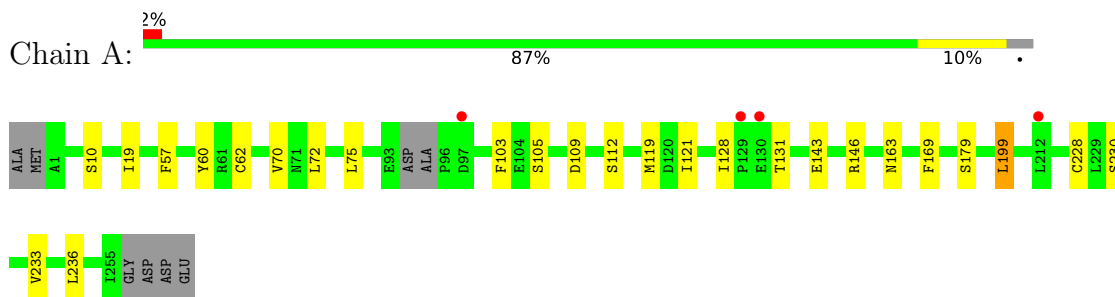
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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	E	20	Total O 20 20	0	0
4	F	20	Total O 20 20	0	0
4	M	3	Total O 3 3	0	0

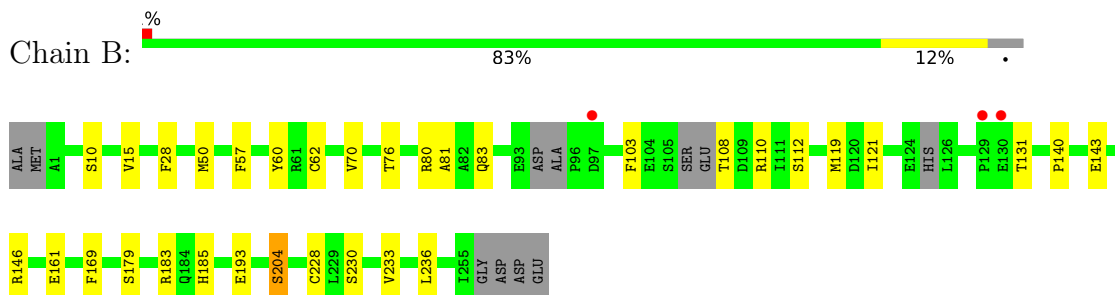
3 Residue-property plots [i](#)

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.

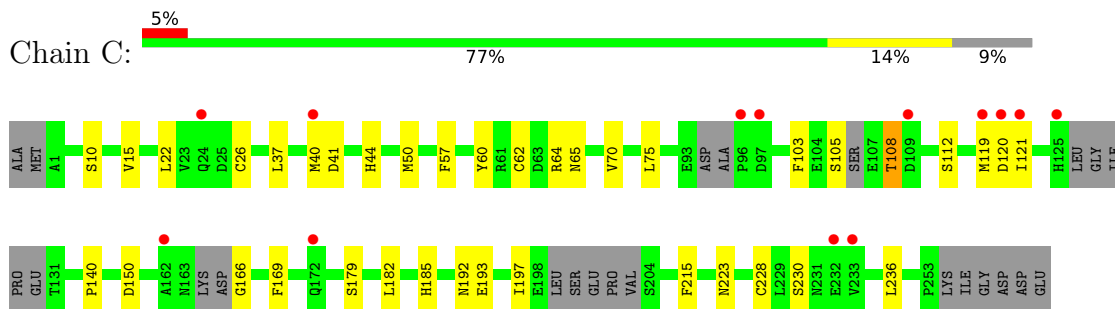
- Molecule 1: Proliferating cell nuclear antigen



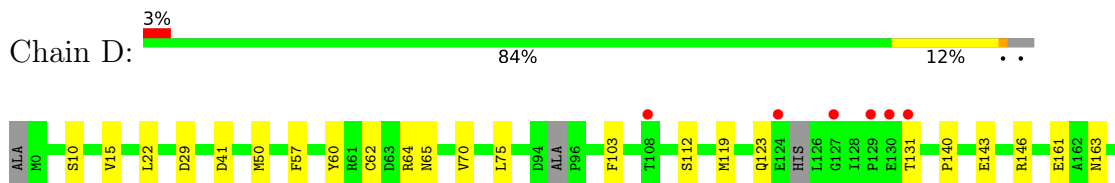
- Molecule 1: Proliferating cell nuclear antigen

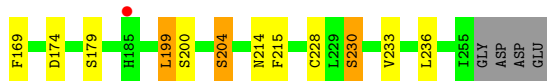


- Molecule 1: Proliferating cell nuclear antigen

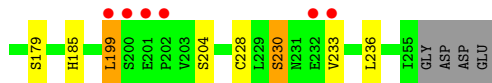
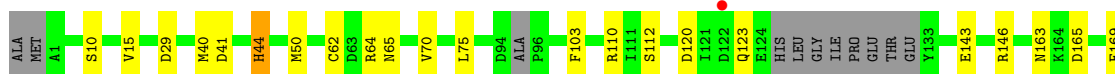
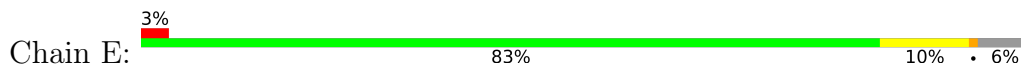


- Molecule 1: Proliferating cell nuclear antigen

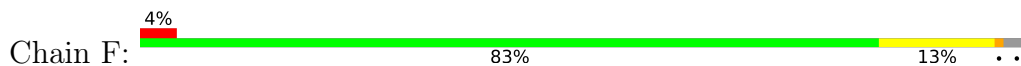




- Molecule 1: Proliferating cell nuclear antigen



- Molecule 1: Proliferating cell nuclear antigen



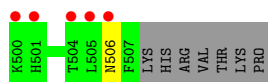
- Molecule 2: Peptide fragment from PolD4



- Molecule 2: Peptide fragment from PolD4



- Molecule 2: Peptide fragment from PolD4



- Molecule 2: Peptide fragment from PolD4



LYS	HS01	●
HS02	●	
HS03	●	
TS04	●	
LS05	●	
HS06	●	
FS07	●	
LYS		
HIS		
ARG		
VAL		
THR		
LYS		
PRO		

4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	71.89Å 84.71Å 84.89Å 60.94° 89.60° 81.22°	Depositor
Resolution (Å)	30.39 – 2.45 30.39 – 2.45	Depositor EDS
% Data completeness (in resolution range)	94.2 (30.39-2.45) 94.2 (30.39-2.45)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.35 (at 2.45Å)	Xtrriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.220 , 0.262 0.226 , 0.267	Depositor DCC
R_{free} test set	2829 reflections (4.74%)	wwPDB-VP
Wilson B-factor (Å ²)	45.7	Xtrriage
Anisotropy	0.148	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	(Not available) , (Not available)	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	0.013 for -h,-k+l,l	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	11613	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.35% 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: NA

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	0.68	0/1913	0.82	0/2594
1	B	0.68	0/1905	0.83	0/2579
1	C	0.69	0/1791	0.83	0/2422
1	D	0.70	0/1916	0.83	0/2597
1	E	0.69	0/1873	0.83	0/2535
1	F	0.68	0/1922	0.83	0/2605
2	J	0.62	0/65	0.70	0/87
2	K	0.68	0/65	0.70	0/87
2	M	0.67	0/65	0.73	0/87
2	O	0.74	0/51	0.75	0/68
All	All	0.69	0/11566	0.83	0/15661

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	1889	0	1879	14	0
1	B	1884	0	1876	17	0
1	C	1774	0	1734	19	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	1894	0	1873	18	0
1	E	1852	0	1846	15	0
1	F	1899	0	1888	17	0
2	J	64	0	54	1	0
2	K	64	0	54	0	0
2	M	64	0	54	0	0
2	O	51	0	43	0	0
3	A	3	0	0	0	0
3	B	2	0	0	0	0
3	C	3	0	0	0	0
3	D	4	0	0	0	0
3	E	4	0	0	0	0
3	F	2	0	0	0	0
4	A	24	0	0	0	0
4	B	32	0	0	0	0
4	C	27	0	0	0	0
4	D	34	0	0	0	0
4	E	20	0	0	0	0
4	F	20	0	0	0	0
4	M	3	0	0	0	0
All	All	11613	0	11301	95	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 4.

All (95) 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:E:230:SER:OG	1:E:233:VAL:HG22	1.89	0.72
1:D:22:LEU:HG	1:D:214:ASN:ND2	2.08	0.69
1:E:29:ASP:OD1	1:E:123:GLN:NE2	2.28	0.67
1:D:29:ASP:OD1	1:D:123:GLN:NE2	2.27	0.64
1:D:131:THR:HG21	1:D:233:VAL:HG21	1.81	0.62
1:B:131:THR:HG21	1:B:233:VAL:HG21	1.83	0.60
1:F:131:THR:HG21	1:F:233:VAL:HG21	1.83	0.60
1:A:131:THR:HG21	1:A:233:VAL:HG21	1.84	0.60
1:E:64:ARG:HD3	1:E:65:ASN:O	2.05	0.56
1:A:109:ASP:OD2	1:B:183:ARG:NH1	2.39	0.56
1:D:64:ARG:HD3	1:D:65:ASN:O	2.06	0.56
1:D:22:LEU:HG	1:D:214:ASN:HD22	1.70	0.55
1:D:103:PHE:HB2	1:D:112:SER:HB2	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:64:ARG:HD3	1:C:65:ASN:O	2.06	0.55
1:C:103:PHE:HB2	1:C:112:SER:HB2	1.89	0.54
1:E:103:PHE:HB2	1:E:112:SER:HB2	1.89	0.54
1:A:103:PHE:HB2	1:A:112:SER:HB2	1.89	0.54
1:B:103:PHE:HB2	1:B:112:SER:HB2	1.88	0.54
1:F:103:PHE:HB2	1:F:112:SER:HB2	1.89	0.53
1:B:119:MET:CE	1:B:121:ILE:HD11	2.40	0.51
1:B:140:PRO:HG3	1:B:193:GLU:HA	1.93	0.51
1:E:40:MET:HE2	1:E:44:HIS:HA	1.92	0.51
1:C:119:MET:CE	1:C:121:ILE:HD11	2.41	0.51
1:B:15:VAL:HG11	1:B:50:MET:SD	2.52	0.50
1:A:119:MET:CE	1:A:121:ILE:HD11	2.42	0.49
1:C:40:MET:HE2	1:C:44:HIS:HA	1.95	0.49
1:D:161:GLU:HG2	1:D:204:SER:HB2	1.95	0.49
1:E:163:ASN:HA	1:E:199:LEU:HD11	1.95	0.49
1:A:70:VAL:HG11	1:A:75:LEU:HD22	1.94	0.48
1:E:15:VAL:HG11	1:E:50:MET:SD	2.54	0.48
1:F:28:PHE:CE1	1:F:70:VAL:HG21	2.48	0.47
1:B:28:PHE:CE1	1:B:70:VAL:HG21	2.50	0.47
1:E:70:VAL:HG11	1:E:75:LEU:HD22	1.97	0.47
1:B:119:MET:HE2	1:B:121:ILE:HD11	1.97	0.47
1:D:174:ASP:O	1:F:117:LYS:HD2	2.15	0.47
1:D:64:ARG:NH1	1:D:65:ASN:O	2.47	0.47
1:A:119:MET:HE2	1:A:121:ILE:HD11	1.97	0.46
1:F:40:MET:HG2	1:F:44:HIS:HA	1.97	0.46
1:E:169:PHE:O	1:E:179:SER:HA	2.16	0.46
1:A:163:ASN:HA	1:A:199:LEU:CD1	2.45	0.46
1:C:15:VAL:HG11	1:C:50:MET:SD	2.55	0.46
1:C:70:VAL:HG11	1:C:75:LEU:HD22	1.98	0.46
1:C:140:PRO:HG3	1:C:193:GLU:HA	1.98	0.46
1:C:119:MET:HE2	1:C:121:ILE:HD11	1.98	0.46
1:D:15:VAL:HG11	1:D:50:MET:SD	2.55	0.46
1:D:169:PHE:O	1:D:179:SER:HA	2.16	0.46
1:F:15:VAL:HG11	1:F:50:MET:SD	2.57	0.45
1:F:169:PHE:O	1:F:179:SER:HA	2.17	0.45
1:B:169:PHE:O	1:B:179:SER:HA	2.16	0.45
1:C:64:ARG:NH1	1:C:65:ASN:O	2.45	0.45
1:A:163:ASN:HA	1:A:199:LEU:HD11	1.99	0.45
1:A:169:PHE:O	1:A:179:SER:HA	2.17	0.45
1:E:64:ARG:NH1	1:E:65:ASN:O	2.48	0.44
1:F:163:ASN:HA	1:F:199:LEU:HD11	1.98	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:110:ARG:HG3	1:C:182:LEU:HD23	1.99	0.44
1:C:169:PHE:O	1:C:179:SER:HA	2.16	0.44
1:E:163:ASN:HA	1:E:199:LEU:CD1	2.48	0.44
1:D:228:CYS:HB2	1:D:236:LEU:HB3	2.00	0.44
1:A:57:PHE:HB2	1:A:60:TYR:HB2	2.00	0.44
1:F:28:PHE:HE1	1:F:70:VAL:HG21	1.83	0.43
1:C:22:LEU:HD11	1:C:215:PHE:CE1	2.53	0.43
1:E:110:ARG:HG3	1:F:182:LEU:HD23	2.00	0.43
1:F:143:GLU:OE1	1:F:146:ARG:NH1	2.52	0.43
1:F:163:ASN:HA	1:F:199:LEU:CD1	2.49	0.43
1:B:228:CYS:HB2	1:B:236:LEU:HB3	2.00	0.43
1:B:143:GLU:OE1	1:B:146:ARG:NH1	2.51	0.43
1:A:128:ILE:HG23	2:J:507:PHE:CD1	2.54	0.43
1:C:57:PHE:HB2	1:C:60:TYR:HB2	2.00	0.43
1:D:143:GLU:OE1	1:D:146:ARG:NH1	2.52	0.43
1:D:230:SER:HB2	1:D:233:VAL:HG22	2.01	0.43
1:C:192:ASN:HD22	1:C:223:ASN:HB3	1.85	0.42
1:C:228:CYS:HB2	1:C:236:LEU:HB3	2.01	0.42
1:D:70:VAL:HG11	1:D:75:LEU:HD22	2.00	0.42
1:A:143:GLU:OE1	1:A:146:ARG:NH1	2.52	0.42
1:B:161:GLU:HG2	1:B:204:SER:HB2	2.01	0.42
1:D:57:PHE:HB2	1:D:60:TYR:HB2	2.01	0.42
1:F:57:PHE:HB2	1:F:60:TYR:HB2	2.01	0.42
1:F:228:CYS:HB2	1:F:236:LEU:HB3	2.00	0.42
1:B:76:THR:O	1:B:80:ARG:HG3	2.19	0.42
1:D:163:ASN:HA	1:D:199:LEU:CD1	2.50	0.42
1:B:28:PHE:HE1	1:B:70:VAL:HG21	1.85	0.42
1:C:26:CYS:SG	1:C:37:LEU:HD11	2.60	0.42
1:A:228:CYS:HB2	1:A:236:LEU:HB3	2.01	0.42
1:F:22:LEU:HD11	1:F:215:PHE:CE1	2.55	0.42
1:E:143:GLU:OE1	1:E:146:ARG:NH1	2.52	0.42
1:B:81:ALA:HA	1:C:150:ASP:OD2	2.20	0.41
1:E:228:CYS:HB2	1:E:236:LEU:HB3	2.02	0.41
1:F:19:ILE:HG22	1:F:72:LEU:CD1	2.51	0.41
1:B:57:PHE:HB2	1:B:60:TYR:HB2	2.01	0.41
1:D:22:LEU:HD11	1:D:215:PHE:CE1	2.56	0.41
1:A:19:ILE:HG22	1:A:72:LEU:CD1	2.50	0.41
1:E:40:MET:HE3	1:E:40:MET:HB3	1.95	0.41
1:F:161:GLU:HG2	1:F:204:SER:OG	2.22	0.40
1:C:166:GLY:HA2	1:C:197:ILE:HD12	2.04	0.40
1:C:105:SER:OG	1:C:108:THR:N	2.55	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	249/261 (95%)	243 (98%)	5 (2%)	1 (0%)	34	41
1	B	242/261 (93%)	239 (99%)	2 (1%)	1 (0%)	34	41
1	C	226/261 (87%)	222 (98%)	3 (1%)	1 (0%)	34	41
1	D	248/261 (95%)	243 (98%)	4 (2%)	1 (0%)	34	41
1	E	240/261 (92%)	234 (98%)	5 (2%)	1 (0%)	34	41
1	F	249/261 (95%)	243 (98%)	5 (2%)	1 (0%)	34	41
2	J	6/15 (40%)	6 (100%)	0	0	100	100
2	K	6/15 (40%)	6 (100%)	0	0	100	100
2	M	6/15 (40%)	6 (100%)	0	0	100	100
2	O	5/15 (33%)	5 (100%)	0	0	100	100
All	All	1477/1626 (91%)	1447 (98%)	24 (2%)	6 (0%)	34	41

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	10	SER
1	E	10	SER
1	F	10	SER
1	A	10	SER
1	C	10	SER
1	D	10	SER

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	207/229 (90%)	203 (98%)	4 (2%)	57	69
1	B	210/229 (92%)	204 (97%)	6 (3%)	42	53
1	C	192/229 (84%)	186 (97%)	6 (3%)	40	52
1	D	206/229 (90%)	198 (96%)	8 (4%)	32	42
1	E	204/229 (89%)	195 (96%)	9 (4%)	28	37
1	F	209/229 (91%)	200 (96%)	9 (4%)	29	38
2	J	7/15 (47%)	7 (100%)	0	100	100
2	K	7/15 (47%)	7 (100%)	0	100	100
2	M	7/15 (47%)	6 (86%)	1 (14%)	3	2
2	O	5/15 (33%)	5 (100%)	0	100	100
All	All	1254/1434 (87%)	1211 (97%)	43 (3%)	37	48

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	62	CYS
1	A	105	SER
1	A	199	LEU
1	A	230	SER
1	B	62	CYS
1	B	83	GLN
1	B	108	THR
1	B	185	HIS
1	B	204	SER
1	B	230	SER
1	C	41	ASP
1	C	62	CYS
1	C	108	THR
1	C	120	ASP
1	C	185	HIS
1	C	230	SER
1	D	41	ASP
1	D	62	CYS
1	D	119	MET
1	D	140	PRO

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Mol	Chain	Res	Type
1	D	199	LEU
1	D	200	SER
1	D	204	SER
1	D	230	SER
1	E	41	ASP
1	E	44	HIS
1	E	62	CYS
1	E	120	ASP
1	E	165	ASP
1	E	185	HIS
1	E	199	LEU
1	E	204	SER
1	E	230	SER
1	F	38	GLN
1	F	40	MET
1	F	62	CYS
1	F	105	SER
1	F	108	THR
1	F	119	MET
1	F	185	HIS
1	F	199	LEU
1	F	230	SER
2	M	506	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 18 ligands modelled in this entry, 18 are 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 [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	253/261 (96%)	0.10	4 (1%) 72 69	30, 43, 67, 81	0
1	B	250/261 (95%)	0.11	3 (1%) 79 77	27, 44, 72, 95	0
1	C	238/261 (91%)	0.22	13 (5%) 25 22	36, 52, 78, 96	0
1	D	254/261 (97%)	0.11	7 (2%) 53 49	29, 43, 69, 100	0
1	E	246/261 (94%)	0.16	7 (2%) 53 49	29, 45, 69, 84	0
1	F	253/261 (96%)	0.24	11 (4%) 35 32	35, 51, 81, 98	0
2	J	8/15 (53%)	1.17	2 (25%) 0 0	37, 41, 57, 78	0
2	K	8/15 (53%)	1.02	2 (25%) 0 0	30, 47, 70, 78	0
2	M	8/15 (53%)	2.08	5 (62%) 0 0	41, 48, 62, 67	0
2	O	7/15 (46%)	2.62	4 (57%) 0 0	67, 68, 79, 80	0
All	All	1525/1626 (93%)	0.19	58 (3%) 40 37	27, 47, 75, 100	0

All (58) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	O	505	LEU	4.7
1	F	129	PRO	4.3
1	B	130	GLU	3.7
1	F	132	GLU	3.7
1	E	232	GLU	3.7
1	E	200	SER	3.6
2	O	503	SER	3.5
2	J	500	LYS	3.4
1	E	233	VAL	3.4
1	D	129	PRO	3.3
2	J	501	HIS	3.3
2	M	501	HIS	3.2
2	O	507	PHE	3.2

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Mol	Chain	Res	Type	RSRZ
1	F	128	ILE	3.1
1	A	130	GLU	3.0
1	A	129	PRO	2.9
1	C	233	VAL	2.9
2	M	506	ASN	2.9
1	F	97	ASP	2.8
1	C	96	PRO	2.8
1	C	120	ASP	2.8
2	K	506	ASN	2.8
1	F	130	GLU	2.8
2	M	504	THR	2.7
1	C	119	MET	2.7
1	F	23	VAL	2.7
1	D	127	GLY	2.7
2	M	505	LEU	2.6
1	E	202	PRO	2.6
1	F	120	ASP	2.6
1	E	201	GLU	2.5
1	C	40	MET	2.5
1	A	97	ASP	2.5
1	E	122	ASP	2.4
2	K	500	LYS	2.4
1	D	130	GLU	2.4
1	C	125	HIS	2.4
1	D	108	THR	2.4
1	D	124	GLU	2.3
1	C	24	GLN	2.3
1	D	185	HIS	2.3
2	O	504	THR	2.3
1	F	7	GLU	2.2
1	B	97	ASP	2.2
1	B	129	PRO	2.2
1	E	199	LEU	2.2
1	F	24	GLN	2.1
1	C	232	GLU	2.1
1	A	212	LEU	2.1
1	C	97	ASP	2.1
1	D	131	THR	2.1
1	F	131	THR	2.1
1	C	109	ASP	2.1
2	M	500	LYS	2.1
1	C	162	ALA	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	121	ILE	2.1
1	C	172	GLN	2.0
1	F	164	LYS	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 monosaccharides 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
3	NA	E	302	1/1	0.70	0.15	57,57,57,57	0
3	NA	D	302	1/1	0.74	0.14	48,48,48,48	0
3	NA	D	303	1/1	0.79	0.12	34,34,34,34	0
3	NA	C	302	1/1	0.79	0.20	59,59,59,59	0
3	NA	B	301	1/1	0.80	0.12	50,50,50,50	0
3	NA	C	303	1/1	0.81	0.14	41,41,41,41	0
3	NA	E	303	1/1	0.85	0.20	40,40,40,40	0
3	NA	E	304	1/1	0.87	0.14	38,38,38,38	0
3	NA	F	301	1/1	0.87	0.14	47,47,47,47	0
3	NA	B	302	1/1	0.88	0.26	45,45,45,45	0
3	NA	A	303	1/1	0.90	0.38	66,66,66,66	0
3	NA	D	304	1/1	0.90	0.15	41,41,41,41	0
3	NA	A	302	1/1	0.90	0.12	49,49,49,49	0
3	NA	C	301	1/1	0.92	0.05	37,37,37,37	0
3	NA	D	301	1/1	0.94	0.12	42,42,42,42	0
3	NA	F	302	1/1	0.94	0.12	56,56,56,56	0
3	NA	E	301	1/1	0.95	0.17	47,47,47,47	0
3	NA	A	301	1/1	0.96	0.20	39,39,39,39	0

6.5 Other polymers [i](#)

There are no such residues in this entry.