



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 04:06 PM UTC

PDB ID : 9GI4 / pdb_00009gi4
Title : TFIIC5 DNA binding domain
Authors : Leen, E.; Bayliss, R.
Deposited on : 2024-08-16
Resolution : 2.63 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
Mogul	:	2022.3.0, CSD as543be (2022)
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

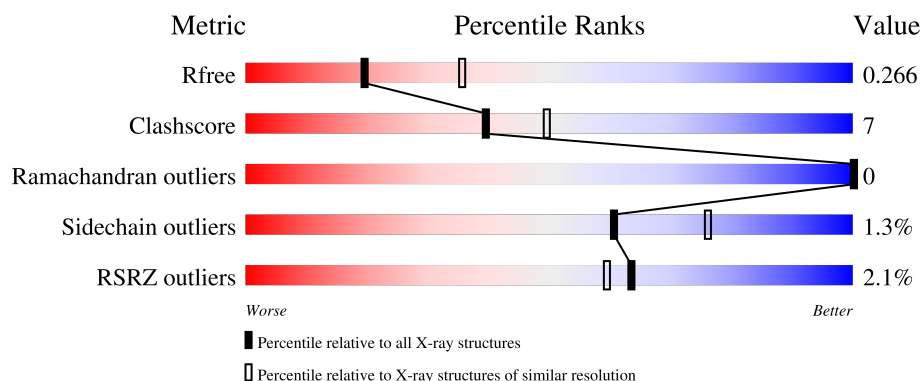
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.63 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	2053 (2.66-2.62)
Clashscore	190562	2097 (2.66-2.62)
Ramachandran outliers	187476	2066 (2.66-2.62)
Sidechain outliers	187428	2066 (2.66-2.62)
RSRZ outliers	180081	2052 (2.66-2.62)

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	242	<div> <div>4%</div> <div>64%</div> <div>14%</div> <div>22%</div> </div>
1	B	242	<div> <div>%</div> <div>66%</div> <div>13%</div> <div>21%</div> </div>
1	C	242	<div> <div>%</div> <div>69%</div> <div>8%</div> <div>22%</div> </div>
1	D	242	<div> <div>2%</div> <div>64%</div> <div>15%</div> <div>20%</div> </div>
1	E	242	<div> <div>%</div> <div>58%</div> <div>19%</div> <div>23%</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	242	<div><div>%</div><div><div></div><div>63%</div><div>15%</div><div>21%</div></div></div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9306 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 General transcription factor 3C polypeptide 5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	189	Total	C	N	O	S	0	0	0
			1532	984	269	271	8			
1	B	191	Total	C	N	O	S	0	1	0
			1560	1001	279	272	8			
1	C	188	Total	C	N	O	S	0	0	0
			1503	968	255	272	8			
1	D	193	Total	C	N	O	S	0	0	0
			1575	1009	281	277	8			
1	E	187	Total	C	N	O	S	0	1	0
			1531	981	269	273	8			
1	F	190	Total	C	N	O	S	0	0	0
			1540	988	272	272	8			

There are 138 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	GLY	-	expression tag	UNP Q9Y5Q8
A	?	-	TYR	deletion	UNP Q9Y5Q8
A	?	-	ASN	deletion	UNP Q9Y5Q8
A	?	-	TYR	deletion	UNP Q9Y5Q8
A	?	-	SER	deletion	UNP Q9Y5Q8
A	?	-	LEU	deletion	UNP Q9Y5Q8
A	?	-	PRO	deletion	UNP Q9Y5Q8
A	?	-	ILE	deletion	UNP Q9Y5Q8
A	?	-	THR	deletion	UNP Q9Y5Q8
A	?	-	VAL	deletion	UNP Q9Y5Q8
A	?	-	LYS	deletion	UNP Q9Y5Q8
A	?	-	LYS	deletion	UNP Q9Y5Q8
A	?	-	THR	deletion	UNP Q9Y5Q8
A	?	-	SER	deletion	UNP Q9Y5Q8
A	?	-	SER	deletion	UNP Q9Y5Q8
A	?	-	GLN	deletion	UNP Q9Y5Q8
A	?	-	LEU	deletion	UNP Q9Y5Q8

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Chain	Residue	Modelled	Actual	Comment	Reference
A	?	-	VAL	deletion	UNP Q9Y5Q8
A	?	-	THR	deletion	UNP Q9Y5Q8
A	?	-	MET	deletion	UNP Q9Y5Q8
A	?	-	HIS	deletion	UNP Q9Y5Q8
A	?	-	ASP	deletion	UNP Q9Y5Q8
A	?	-	LEU	deletion	UNP Q9Y5Q8
B	1	GLY	-	expression tag	UNP Q9Y5Q8
B	?	-	TYR	deletion	UNP Q9Y5Q8
B	?	-	ASN	deletion	UNP Q9Y5Q8
B	?	-	TYR	deletion	UNP Q9Y5Q8
B	?	-	SER	deletion	UNP Q9Y5Q8
B	?	-	LEU	deletion	UNP Q9Y5Q8
B	?	-	PRO	deletion	UNP Q9Y5Q8
B	?	-	ILE	deletion	UNP Q9Y5Q8
B	?	-	THR	deletion	UNP Q9Y5Q8
B	?	-	VAL	deletion	UNP Q9Y5Q8
B	?	-	LYS	deletion	UNP Q9Y5Q8
B	?	-	LYS	deletion	UNP Q9Y5Q8
B	?	-	THR	deletion	UNP Q9Y5Q8
B	?	-	SER	deletion	UNP Q9Y5Q8
B	?	-	SER	deletion	UNP Q9Y5Q8
B	?	-	GLN	deletion	UNP Q9Y5Q8
B	?	-	LEU	deletion	UNP Q9Y5Q8
B	?	-	VAL	deletion	UNP Q9Y5Q8
B	?	-	THR	deletion	UNP Q9Y5Q8
B	?	-	MET	deletion	UNP Q9Y5Q8
B	?	-	HIS	deletion	UNP Q9Y5Q8
B	?	-	ASP	deletion	UNP Q9Y5Q8
B	?	-	LEU	deletion	UNP Q9Y5Q8
C	1	GLY	-	expression tag	UNP Q9Y5Q8
C	?	-	TYR	deletion	UNP Q9Y5Q8
C	?	-	ASN	deletion	UNP Q9Y5Q8
C	?	-	TYR	deletion	UNP Q9Y5Q8
C	?	-	SER	deletion	UNP Q9Y5Q8
C	?	-	LEU	deletion	UNP Q9Y5Q8
C	?	-	PRO	deletion	UNP Q9Y5Q8
C	?	-	ILE	deletion	UNP Q9Y5Q8
C	?	-	THR	deletion	UNP Q9Y5Q8
C	?	-	VAL	deletion	UNP Q9Y5Q8
C	?	-	LYS	deletion	UNP Q9Y5Q8
C	?	-	LYS	deletion	UNP Q9Y5Q8
C	?	-	THR	deletion	UNP Q9Y5Q8

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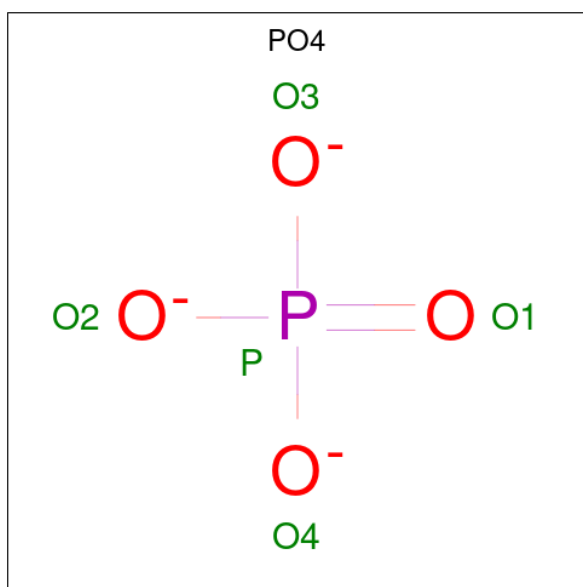
Chain	Residue	Modelled	Actual	Comment	Reference
C	?	-	SER	deletion	UNP Q9Y5Q8
C	?	-	SER	deletion	UNP Q9Y5Q8
C	?	-	GLN	deletion	UNP Q9Y5Q8
C	?	-	LEU	deletion	UNP Q9Y5Q8
C	?	-	VAL	deletion	UNP Q9Y5Q8
C	?	-	THR	deletion	UNP Q9Y5Q8
C	?	-	MET	deletion	UNP Q9Y5Q8
C	?	-	HIS	deletion	UNP Q9Y5Q8
C	?	-	ASP	deletion	UNP Q9Y5Q8
C	?	-	LEU	deletion	UNP Q9Y5Q8
D	1	GLY	-	expression tag	UNP Q9Y5Q8
D	?	-	TYR	deletion	UNP Q9Y5Q8
D	?	-	ASN	deletion	UNP Q9Y5Q8
D	?	-	TYR	deletion	UNP Q9Y5Q8
D	?	-	SER	deletion	UNP Q9Y5Q8
D	?	-	LEU	deletion	UNP Q9Y5Q8
D	?	-	PRO	deletion	UNP Q9Y5Q8
D	?	-	ILE	deletion	UNP Q9Y5Q8
D	?	-	THR	deletion	UNP Q9Y5Q8
D	?	-	VAL	deletion	UNP Q9Y5Q8
D	?	-	LYS	deletion	UNP Q9Y5Q8
D	?	-	LYS	deletion	UNP Q9Y5Q8
D	?	-	THR	deletion	UNP Q9Y5Q8
D	?	-	SER	deletion	UNP Q9Y5Q8
D	?	-	SER	deletion	UNP Q9Y5Q8
D	?	-	GLN	deletion	UNP Q9Y5Q8
D	?	-	LEU	deletion	UNP Q9Y5Q8
D	?	-	VAL	deletion	UNP Q9Y5Q8
D	?	-	THR	deletion	UNP Q9Y5Q8
D	?	-	MET	deletion	UNP Q9Y5Q8
D	?	-	HIS	deletion	UNP Q9Y5Q8
D	?	-	ASP	deletion	UNP Q9Y5Q8
D	?	-	LEU	deletion	UNP Q9Y5Q8
E	1	GLY	-	expression tag	UNP Q9Y5Q8
E	?	-	TYR	deletion	UNP Q9Y5Q8
E	?	-	ASN	deletion	UNP Q9Y5Q8
E	?	-	TYR	deletion	UNP Q9Y5Q8
E	?	-	SER	deletion	UNP Q9Y5Q8
E	?	-	LEU	deletion	UNP Q9Y5Q8
E	?	-	PRO	deletion	UNP Q9Y5Q8
E	?	-	ILE	deletion	UNP Q9Y5Q8
E	?	-	THR	deletion	UNP Q9Y5Q8

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Chain	Residue	Modelled	Actual	Comment	Reference
E	?	-	VAL	deletion	UNP Q9Y5Q8
E	?	-	LYS	deletion	UNP Q9Y5Q8
E	?	-	LYS	deletion	UNP Q9Y5Q8
E	?	-	THR	deletion	UNP Q9Y5Q8
E	?	-	SER	deletion	UNP Q9Y5Q8
E	?	-	SER	deletion	UNP Q9Y5Q8
E	?	-	GLN	deletion	UNP Q9Y5Q8
E	?	-	LEU	deletion	UNP Q9Y5Q8
E	?	-	VAL	deletion	UNP Q9Y5Q8
E	?	-	THR	deletion	UNP Q9Y5Q8
E	?	-	MET	deletion	UNP Q9Y5Q8
E	?	-	HIS	deletion	UNP Q9Y5Q8
E	?	-	ASP	deletion	UNP Q9Y5Q8
E	?	-	LEU	deletion	UNP Q9Y5Q8
F	1	GLY	-	expression tag	UNP Q9Y5Q8
F	?	-	TYR	deletion	UNP Q9Y5Q8
F	?	-	ASN	deletion	UNP Q9Y5Q8
F	?	-	TYR	deletion	UNP Q9Y5Q8
F	?	-	SER	deletion	UNP Q9Y5Q8
F	?	-	LEU	deletion	UNP Q9Y5Q8
F	?	-	PRO	deletion	UNP Q9Y5Q8
F	?	-	ILE	deletion	UNP Q9Y5Q8
F	?	-	THR	deletion	UNP Q9Y5Q8
F	?	-	VAL	deletion	UNP Q9Y5Q8
F	?	-	LYS	deletion	UNP Q9Y5Q8
F	?	-	LYS	deletion	UNP Q9Y5Q8
F	?	-	THR	deletion	UNP Q9Y5Q8
F	?	-	SER	deletion	UNP Q9Y5Q8
F	?	-	SER	deletion	UNP Q9Y5Q8
F	?	-	GLN	deletion	UNP Q9Y5Q8
F	?	-	LEU	deletion	UNP Q9Y5Q8
F	?	-	VAL	deletion	UNP Q9Y5Q8
F	?	-	THR	deletion	UNP Q9Y5Q8
F	?	-	MET	deletion	UNP Q9Y5Q8
F	?	-	HIS	deletion	UNP Q9Y5Q8
F	?	-	ASP	deletion	UNP Q9Y5Q8
F	?	-	LEU	deletion	UNP Q9Y5Q8

- Molecule 2 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	P	0	0
			5	4	1		
2	B	1	Total	O	P	0	0
			5	4	1		
2	C	1	Total	O	P	0	0
			5	4	1		
2	D	1	Total	O	P	0	0
			5	4	1		
2	D	1	Total	O	P	0	0
			5	4	1		
2	E	1	Total	O	P	0	0
			5	4	1		
2	F	1	Total	O	P	0	0
			5	4	1		

- Molecule 3 is water.

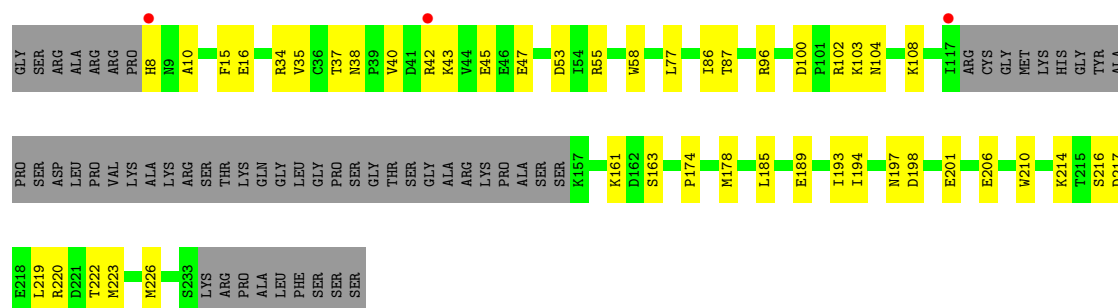
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	6	Total	O	0	0
			6	6		
3	B	6	Total	O	0	0
			6	6		
3	C	1	Total	O	0	0
			1	1		
3	D	4	Total	O	0	0
			4	4		
3	E	8	Total	O	0	0
			8	8		

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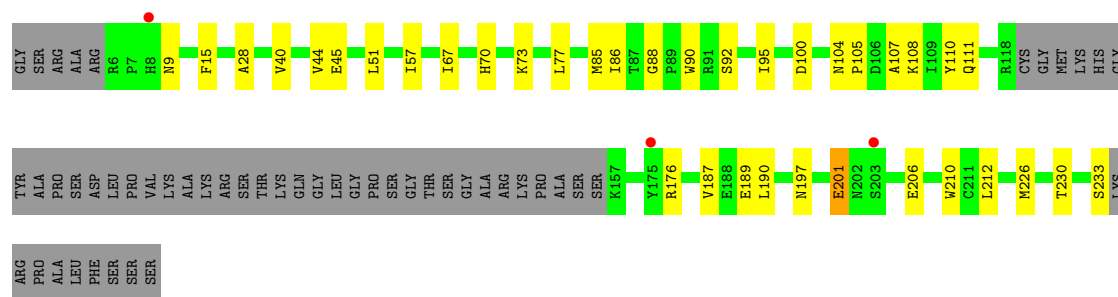
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	F	5	Total	O	0	0
			5	5		

- Molecule 1: General transcription factor 3C polypeptide 5



- Molecule 1: General transcription factor 3C polypeptide 5



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	37.52Å 37.64Å 276.68Å 90.01° 90.02° 59.90°	Depositor
Resolution (Å)	46.11 – 2.63 46.11 – 2.63	Depositor EDS
% Data completeness (in resolution range)	97.8 (46.11-2.63) 98.1 (46.11-2.63)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.72 (at 2.61Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487, PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.244 , 0.300 (Not available) , 0.266	Depositor DCC
R_{free} test set	1828 reflections (4.68%)	wwPDB-VP
Wilson B-factor (Å ²)	67.1	Xtriage
Anisotropy	0.361	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 70.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	0.009 for h-k,h,l 0.009 for k,-h+k,l 0.397 for -k,h-k,l 0.397 for -h+k,-h,l 0.417 for h,h-k,-l 0.407 for -k,-h,-l 0.013 for -h,-k,l 0.014 for h-k,-k,-l 0.418 for -h+k,k,-l 0.014 for -h,-h+k,-l 0.012 for k,h,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	9306	wwPDB-VP
Average B, all atoms (Å ²)	69.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.02% 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: PO4

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	0.09	0/1571	0.24	0/2133
1	B	0.09	0/1606	0.24	0/2179
1	C	0.09	0/1543	0.23	0/2100
1	D	0.10	0/1615	0.27	0/2189
1	E	0.09	0/1569	0.26	0/2129
1	F	0.10	0/1579	0.26	0/2144
All	All	0.09	0/9483	0.25	0/12874

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

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	1532	0	1494	20	0
1	B	1560	0	1522	20	0
1	C	1503	0	1436	15	0
1	D	1575	0	1545	21	0
1	E	1531	0	1486	27	0
1	F	1540	0	1496	23	0
2	A	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	5	0	0	0	0
2	C	5	0	0	0	0
2	D	10	0	0	0	0
2	E	5	0	0	0	0
2	F	5	0	0	0	0
3	A	6	0	0	0	0
3	B	6	0	0	0	0
3	C	1	0	0	0	0
3	D	4	0	0	0	0
3	E	8	0	0	0	0
3	F	5	0	0	0	0
All	All	9306	0	8979	123	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 (123) 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:F:176:ARG:HD3	1:F:176:ARG:H	1.39	0.85
1:D:18:GLU:HG3	1:D:19:GLU:HG2	1.64	0.79
1:F:187:VAL:HG21	1:F:226:MET:HE1	1.64	0.79
1:C:187:VAL:HG21	1:C:226:MET:HE1	1.68	0.75
1:A:173:PRO:HD2	1:A:227:ILE:HD11	1.68	0.75
1:E:161:LYS:HE3	1:E:163:SER:HB3	1.72	0.71
1:C:86:ILE:HG22	1:C:87:THR:HG23	1.73	0.70
1:B:26:GLU:OE2	1:E:42[B]:ARG:NE	2.26	0.67
1:E:16:GLU:N	1:E:16:GLU:OE1	2.29	0.65
1:A:86:ILE:HG22	1:A:87:THR:HG23	1.80	0.63
1:E:189:GLU:O	1:E:193:ILE:HD12	1.99	0.63
1:E:219:LEU:O	1:E:223:MET:HG3	1.98	0.62
1:F:176:ARG:H	1:F:176:ARG:CD	2.12	0.62
1:B:206:GLU:HG2	1:B:207:ARG:HD2	1.80	0.61
1:B:187:VAL:HG11	1:B:226:MET:HE1	1.82	0.61
1:D:57:ILE:HG13	1:D:95:ILE:HG23	1.80	0.61
1:E:43:LYS:O	1:E:47:GLU:HG3	1.99	0.61
1:B:108:LYS:HB2	1:B:210:TRP:HB2	1.81	0.61
1:C:166:ILE:HG12	1:C:184:ASP:HB3	1.83	0.60
1:D:85:MET:HE3	1:D:95:ILE:HD12	1.84	0.60
1:C:216:SER:HB3	1:C:220:ARG:HH21	1.67	0.59
1:B:40:VAL:O	1:B:44:VAL:HG23	2.02	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:115:PHE:HD1	1:B:220:ARG:HG2	1.66	0.59
1:D:69:VAL:HG11	1:D:74:LEU:HD13	1.86	0.57
1:C:45:GLU:HA	1:C:77:LEU:HD21	1.87	0.56
1:E:214:LYS:HE3	1:E:217:ASP:CG	2.30	0.56
1:A:219:LEU:O	1:A:223:MET:HG3	2.05	0.56
1:D:212:LEU:HD12	1:D:212:LEU:H	1.71	0.56
1:B:189:GLU:N	1:B:189:GLU:OE2	2.38	0.56
1:D:108:LYS:HB2	1:D:210:TRP:HB2	1.87	0.56
1:C:189:GLU:N	1:C:189:GLU:OE2	2.39	0.55
1:D:166:ILE:HG12	1:D:184:ASP:HB3	1.87	0.55
1:F:100:ASP:O	1:F:104:ASN:ND2	2.39	0.55
1:E:108:LYS:HB2	1:E:210:TRP:HB2	1.88	0.55
1:F:15:PHE:CD2	1:F:206:GLU:HG3	2.42	0.53
1:E:55:ARG:HD2	1:E:58:TRP:CE2	2.43	0.53
1:B:219:LEU:O	1:B:223:MET:HG3	2.09	0.53
1:F:88:GLY:HA2	1:F:206:GLU:HG2	1.90	0.52
1:C:85:MET:HG3	1:C:95:ILE:HD12	1.92	0.52
1:F:105:PRO:HB3	1:F:201:GLU:HG3	1.91	0.52
1:A:45:GLU:HA	1:A:77:LEU:HD21	1.92	0.52
1:F:57:ILE:HG23	1:F:95:ILE:HG13	1.92	0.52
1:B:201:GLU:HG2	1:B:210:TRP:HB3	1.92	0.52
1:A:115:PHE:HD1	1:A:220:ARG:HG2	1.74	0.51
1:E:197:ASN:ND2	1:E:210:TRP:O	2.44	0.51
1:C:115:PHE:HD2	1:C:220:ARG:HG3	1.76	0.51
1:D:17:ASP:O	1:D:102:ARG:NH1	2.38	0.50
1:B:85:MET:HE2	1:B:95:ILE:HD13	1.93	0.50
1:E:15:PHE:CD1	1:E:206:GLU:HG2	2.46	0.50
1:E:100:ASP:O	1:E:104:ASN:ND2	2.44	0.50
1:B:115:PHE:CD1	1:B:220:ARG:HG2	2.45	0.49
1:F:40:VAL:O	1:F:44:VAL:HG13	2.12	0.49
1:E:222:THR:O	1:E:226:MET:HG3	2.12	0.49
1:B:168:ARG:HG3	1:B:186:ASN:HB3	1.95	0.49
1:C:51:LEU:HD11	1:C:67:ILE:HG12	1.93	0.49
1:A:9:ASN:HB2	1:A:28:ALA:HB2	1.95	0.48
1:D:83:TYR:HE1	1:D:85:MET:HE2	1.78	0.48
1:E:53:ASP:HA	1:E:96:ARG:HH11	1.77	0.48
1:F:108:LYS:O	1:F:111:GLN:HG3	2.13	0.48
1:F:45:GLU:HA	1:F:77:LEU:HD21	1.95	0.48
1:C:45:GLU:HG3	1:C:77:LEU:HD11	1.96	0.48
1:A:216:SER:HA	1:A:219:LEU:HD12	1.95	0.48
1:B:44:VAL:HG13	1:B:67:ILE:HG21	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:107:ALA:HA	1:D:110:TYR:HD2	1.79	0.47
1:B:51:LEU:HD12	1:B:67:ILE:HD11	1.95	0.47
1:F:85:MET:HE3	1:F:90:TRP:HB2	1.96	0.47
1:F:70:HIS:HB3	1:F:73:LYS:HG3	1.96	0.47
1:A:173:PRO:HB3	1:A:178:MET:SD	2.54	0.47
1:D:207:ARG:HH22	1:D:214:LYS:HZ2	1.63	0.47
1:E:108:LYS:NZ	1:E:198:ASP:OD1	2.46	0.47
1:F:230:THR:O	1:F:233:SER:OG	2.31	0.47
1:F:9:ASN:HB2	1:F:28:ALA:HB2	1.97	0.47
1:E:38:ASN:HB3	1:E:40:VAL:HG12	1.96	0.46
1:E:86:ILE:HG23	1:E:87:THR:HG23	1.98	0.46
1:B:45:GLU:HA	1:B:77:LEU:HD21	1.97	0.46
1:F:187:VAL:HG12	1:F:190:LEU:H	1.79	0.46
1:B:157:LYS:HA	1:B:157:LYS:HD2	1.72	0.45
1:F:187:VAL:HB	1:F:190:LEU:HD12	1.98	0.45
1:B:55:ARG:O	1:B:58:TRP:NE1	2.49	0.45
1:C:187:VAL:HB	1:C:190:LEU:HD12	1.99	0.45
1:E:201:GLU:HG2	1:E:210:TRP:HB3	1.99	0.45
1:D:100:ASP:O	1:D:104:ASN:ND2	2.50	0.45
1:D:192:LYS:HD3	1:D:192:LYS:HA	1.61	0.44
1:E:174:PRO:HD2	1:E:178:MET:SD	2.57	0.44
1:A:114:ASP:OD1	1:A:114:ASP:N	2.45	0.44
1:A:69:VAL:HG21	1:A:74:LEU:HD13	2.00	0.44
1:C:216:SER:HB3	1:C:220:ARG:NH2	2.33	0.44
1:D:49:ARG:HG3	1:D:81:ILE:HD12	1.99	0.44
1:F:197:ASN:HB3	1:F:212:LEU:HG	2.00	0.44
1:F:108:LYS:HB2	1:F:210:TRP:HB2	2.00	0.44
1:D:219:LEU:O	1:D:223:MET:HG3	2.17	0.44
1:E:102:ARG:HE	1:E:102:ARG:HB2	1.62	0.44
1:A:18:GLU:O	1:A:102:ARG:NH2	2.50	0.43
1:A:212:LEU:O	1:A:215:THR:OG1	2.24	0.43
1:B:44:VAL:HG11	1:B:69:VAL:HB	1.99	0.43
1:B:173:PRO:HA	1:B:174:PRO:HD3	1.92	0.43
1:E:35:VAL:O	1:E:37:THR:HG23	2.18	0.43
1:D:84:TYR:CZ	1:D:92:SER:HA	2.53	0.43
1:D:109:ILE:HD12	1:D:183:CYS:HB3	2.01	0.43
1:B:26:GLU:HB2	1:E:42[B]:ARG:HG2	2.00	0.43
1:F:51:LEU:HD12	1:F:67:ILE:HD11	2.00	0.43
1:D:45:GLU:HA	1:D:77:LEU:HD21	2.01	0.43
1:E:8:HIS:C	1:E:10:ALA:H	2.27	0.42
1:A:34:ARG:HB2	1:D:19:GLU:OE1	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:45:GLU:HA	1:E:77:LEU:HD21	2.01	0.42
1:E:185:LEU:HD12	1:E:194:ILE:HD11	2.01	0.42
1:A:7:PRO:HB2	1:A:8:HIS:H	1.71	0.42
1:D:40:VAL:O	1:D:44:VAL:HG13	2.19	0.42
1:C:212:LEU:HD12	1:C:212:LEU:HA	1.81	0.42
1:E:214:LYS:HD2	1:E:214:LYS:HA	1.81	0.42
1:E:216:SER:HB2	1:E:220:ARG:NH2	2.35	0.41
1:A:12:PHE:HD1	1:A:86:ILE:HD11	1.85	0.41
1:A:187:VAL:HG21	1:A:226:MET:HE1	2.01	0.41
1:A:215:THR:O	1:A:219:LEU:HD12	2.21	0.41
1:A:38:ASN:HB2	1:A:40:VAL:HG12	2.03	0.41
1:C:86:ILE:HD13	1:C:86:ILE:HA	1.98	0.41
1:F:107:ALA:HA	1:F:110:TYR:HD1	1.86	0.41
1:A:173:PRO:HA	1:A:174:PRO:HD3	1.85	0.41
1:A:201:GLU:HG2	1:A:210:TRP:HB3	2.02	0.41
1:D:114:ASP:OD1	1:D:115:PHE:N	2.54	0.40
1:F:197:ASN:ND2	1:F:210:TRP:O	2.54	0.40
1:C:102:ARG:HE	1:C:102:ARG:HB2	1.60	0.40
1:F:189:GLU:OE1	1:F:189:GLU:N	2.55	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	185/242 (76%)	178 (96%)	7 (4%)	0	100	100
1	B	188/242 (78%)	181 (96%)	7 (4%)	0	100	100
1	C	184/242 (76%)	180 (98%)	4 (2%)	0	100	100
1	D	189/242 (78%)	184 (97%)	5 (3%)	0	100	100
1	E	184/242 (76%)	178 (97%)	6 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	F	186/242 (77%)	180 (97%)	6 (3%)	0	100	100
All	All	1116/1452 (77%)	1081 (97%)	35 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	162/214 (76%)	160 (99%)	2 (1%)	63	78
1	B	164/214 (77%)	160 (98%)	4 (2%)	43	64
1	C	157/214 (73%)	155 (99%)	2 (1%)	61	76
1	D	167/214 (78%)	166 (99%)	1 (1%)	78	87
1	E	161/214 (75%)	159 (99%)	2 (1%)	63	78
1	F	161/214 (75%)	158 (98%)	3 (2%)	50	70
All	All	972/1284 (76%)	958 (99%)	14 (1%)	61	75

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	GLU
1	A	112	VAL
1	B	42[A]	ARG
1	B	42[B]	ARG
1	B	81	ILE
1	B	164	VAL
1	C	112	VAL
1	C	212	LEU
1	D	204	CYS
1	E	34	ARG
1	E	103	LYS
1	F	86	ILE
1	F	92	SER

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Mol	Chain	Res	Type
1	F	201	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	38	ASN
1	A	177	GLN
1	A	191	GLN
1	B	181	GLN
1	C	66	ASN
1	D	66	ASN
1	E	14	ASN
1	F	66	ASN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

7 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	PO4	A	301	-	4,4,4	0.96	0	6,6,6	0.46	0
2	PO4	F	301	-	4,4,4	0.96	0	6,6,6	0.46	0
2	PO4	C	301	-	4,4,4	0.95	0	6,6,6	0.46	0
2	PO4	E	301	-	4,4,4	0.95	0	6,6,6	0.46	0
2	PO4	D	302	-	4,4,4	0.95	0	6,6,6	0.47	0
2	PO4	D	301	-	4,4,4	0.96	0	6,6,6	0.46	0
2	PO4	B	301	-	4,4,4	0.96	0	6,6,6	0.46	0

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	189/242 (78%)	0.46	9 (4%)	35 29	47, 68, 86, 105	0
1	B	191/242 (78%)	0.41	3 (1%)	70 67	42, 66, 88, 102	1 (0%)
1	C	188/242 (77%)	0.46	2 (1%)	78 75	51, 70, 89, 100	0
1	D	193/242 (79%)	0.40	4 (2%)	63 59	50, 66, 90, 97	0
1	E	187/242 (77%)	0.40	3 (1%)	70 67	32, 67, 88, 92	1 (0%)
1	F	190/242 (78%)	0.45	3 (1%)	70 67	52, 71, 89, 108	0
All	All	1138/1452 (78%)	0.43	24 (2%)	63 59	32, 68, 89, 108	2 (0%)

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	42[A]	ARG	4.4
1	E	8	HIS	3.9
1	A	7	PRO	3.7
1	A	37	THR	3.3
1	F	175	TYR	3.2
1	C	7	PRO	3.0
1	D	7	PRO	2.6
1	A	225	LEU	2.6
1	F	203	SER	2.5
1	E	117	ILE	2.5
1	A	229	GLN	2.5
1	A	175	TYR	2.3
1	A	51	LEU	2.3
1	B	204	CYS	2.2
1	D	110	TYR	2.2
1	C	117	ILE	2.2
1	F	8	HIS	2.1
1	A	230	THR	2.1
1	B	236	PRO	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	179	PHE	2.1
1	B	117	ILE	2.1
1	A	231	ILE	2.0
1	A	179	PHE	2.0
1	D	231	ILE	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(Å ²)	Q<0.9
2	PO4	D	302	5/5	0.65	0.15	86,91,100,110	0
2	PO4	B	301	5/5	0.82	0.14	79,83,96,115	0
2	PO4	A	301	5/5	0.83	0.17	90,91,92,114	0
2	PO4	E	301	5/5	0.83	0.15	74,86,86,113	0
2	PO4	C	301	5/5	0.85	0.14	82,88,104,123	0
2	PO4	D	301	5/5	0.85	0.13	93,100,111,125	0
2	PO4	F	301	5/5	0.87	0.12	80,87,96,118	0

6.5 Other polymers [i](#)

There are no such residues in this entry.