



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 30, 2022 – 04:14 pm BST

PDB ID : 7R08  
Title : Abortive infection DNA polymerase Abi-P2  
Authors : Gapinska, M.A.; Figiel, M.; Czarnocki Cieciora, M.; Nowotny, M.; Zajko, W.  
Deposited on : 2022-02-01  
Resolution : 3.10 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](mailto: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>.

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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 : 2.30  
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) : 2.30

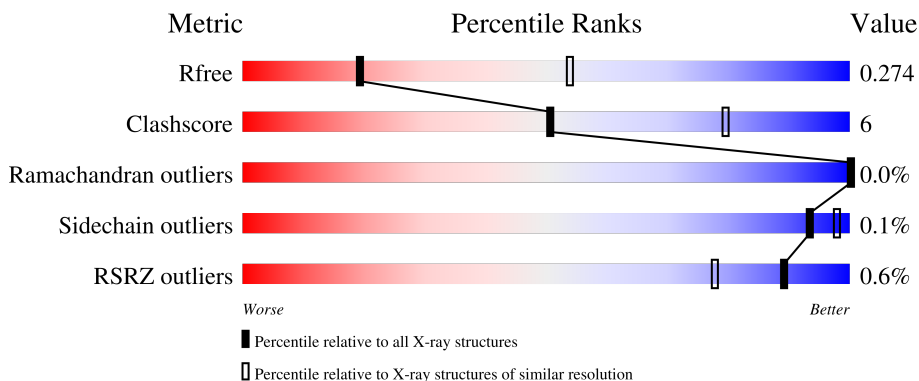
# 1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.10 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	1094 (3.10-3.10)
Clashscore	141614	1184 (3.10-3.10)
Ramachandran outliers	138981	1141 (3.10-3.10)
Sidechain outliers	138945	1141 (3.10-3.10)
RSRZ outliers	127900	1067 (3.10-3.10)

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  $\geq 3$ , 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	541	 75% 15% 10%
1	B	541	 76% 13% 11%
1	C	541	 77% 12% 11%
1	D	541	 75% 13% 11%
1	E	541	 76% 14% 10%

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Mol	Chain	Length	Quality of chain
1	F	541	 <p>A horizontal bar chart representing the quality of chain. The bar is divided into three segments: a green segment representing 79%, a yellow segment representing 10%, and a grey segment representing 11%. A small red square is located at the beginning of the bar, and a '%' symbol is positioned above it.</p>

## 2 Entry composition

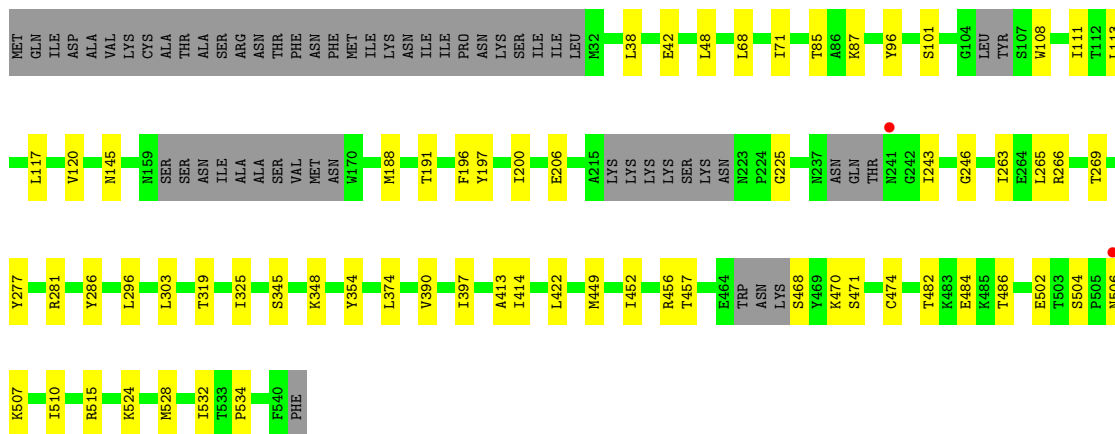
There is only 1 type of molecule in this entry. The entry contains 22205 atoms, of which 0 are hydrogens and 0 are deuteriums.

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.

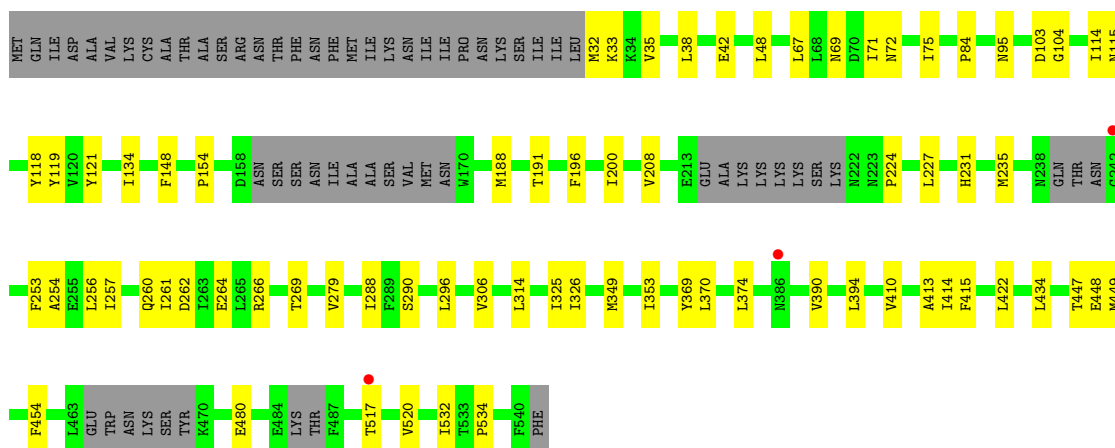
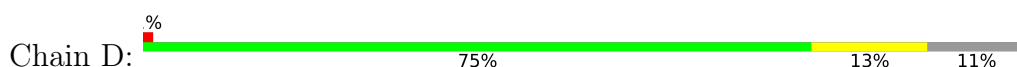
- Molecule 1 is a protein called Reverse transcriptase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	487	3794	2435	615	729	15	0	0	0
1	B	482	3757	2418	609	715	15	0	0	0
1	C	484	3668	2364	582	707	15	0	0	0
1	D	479	3599	2324	568	693	14	0	0	0
1	E	488	3729	2396	599	719	15	0	0	0
1	F	479	3658	2357	589	697	15	0	0	0

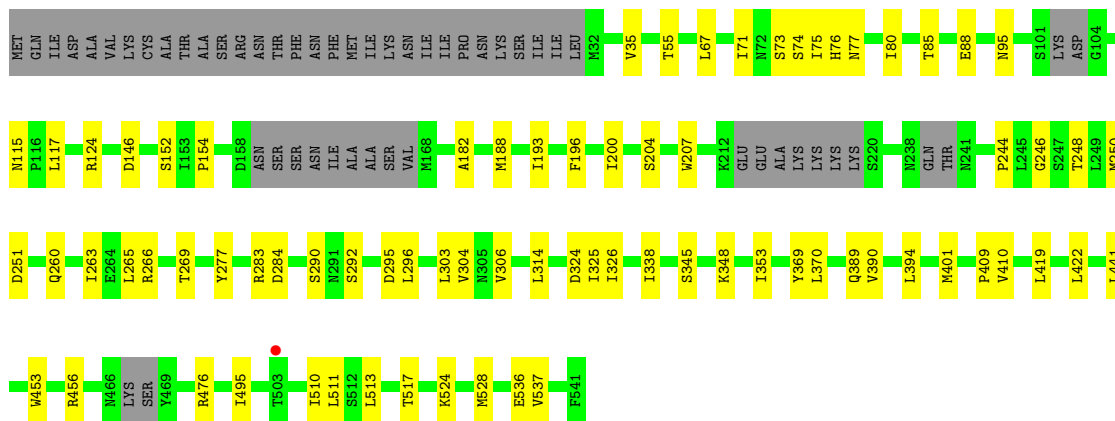
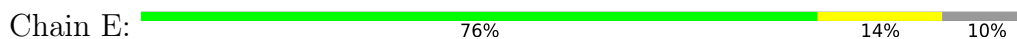




• Molecule 1: Reverse transcriptase

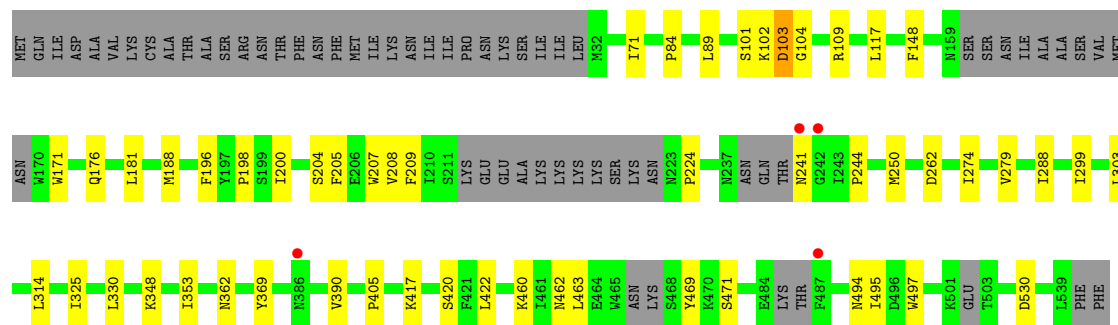


• Molecule 1: Reverse transcriptase



• Molecule 1: Reverse transcriptase

Chain F: %



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	134.33Å 88.36Å 161.09Å 90.00° 103.18° 90.00°	Depositor
Resolution (Å)	47.91 – 3.10 47.91 – 3.09	Depositor EDS
% Data completeness (in resolution range)	99.7 (47.91-3.10) 90.3 (47.91-3.09)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	-0.13 (at 3.07Å)	Xtriage
Refinement program	PHENIX 1.18.2_3874	Depositor
R, $R_{free}$	0.220 , 0.274 0.220 , 0.274	Depositor DCC
$R_{free}$ test set	2100 reflections (3.13%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	86.9	Xtriage
Anisotropy	0.518	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	(Not available) , (Not available)	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	22205	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	106.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.31% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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	0.24	0/3868	0.40	0/5256
1	B	0.23	0/3831	0.39	0/5199
1	C	0.23	0/3740	0.38	0/5104
1	D	0.23	0/3673	0.39	0/5016
1	E	0.23	0/3805	0.40	0/5192
1	F	0.23	0/3729	0.38	0/5077
All	All	0.23	0/22646	0.39	0/30844

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	3794	0	3588	47	0
1	B	3757	0	3532	45	0
1	C	3668	0	3367	37	0
1	D	3599	0	3258	43	0
1	E	3729	0	3411	45	0
1	F	3658	0	3398	32	0
All	All	22205	0	20554	242	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 (242) 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:38:LEU:HB3	1:D:42:GLU:HG3	1.79	0.65
1:F:102:LYS:O	1:F:104:GLY:N	2.30	0.65
1:A:189:PHE:HB2	1:A:296:LEU:HD23	1.79	0.63
1:F:102:LYS:HB2	1:F:109:ARG:HG3	1.82	0.61
1:A:89:LEU:HD12	1:A:94:ILE:HD11	1.82	0.61
1:A:125:LYS:HD3	1:A:227:LEU:HD11	1.83	0.61
1:D:196:PHE:HE2	1:D:254:ALA:HB1	1.66	0.60
1:D:390:VAL:HG11	1:D:422:LEU:HB2	1.83	0.60
1:A:456:ARG:HD2	1:A:532:ILE:HG12	1.82	0.60
1:F:148:PHE:HD2	1:F:262:ASP:HB3	1.66	0.60
1:F:188:MET:HB2	1:F:325:ILE:HG12	1.84	0.60
1:E:290:SER:HB3	1:E:296:LEU:HG	1.84	0.59
1:B:355:LEU:HA	1:B:358:LEU:HD12	1.85	0.59
1:E:73:SER:O	1:E:77:ASN:ND2	2.36	0.59
1:D:447:THR:OG1	1:D:448:GLU:N	2.37	0.58
1:F:274:ILE:HG21	1:F:299:ILE:HD11	1.86	0.58
1:D:279:VAL:HG22	1:D:288:ILE:HG12	1.85	0.58
1:A:57:GLU:O	1:A:453:TRP:NE1	2.29	0.57
1:D:208:VAL:HG11	1:D:257:ILE:HG22	1.84	0.57
1:E:204:SER:HA	1:E:207:TRP:CD1	2.39	0.57
1:A:489:VAL:HG11	1:A:514:LEU:HB3	1.87	0.57
1:C:265:LEU:HD22	1:C:303:LEU:HD13	1.86	0.57
1:D:434:LEU:HB3	1:D:454:PHE:HZ	1.70	0.57
1:C:471:SER:HB3	1:C:474:CYS:HB2	1.86	0.56
1:A:40:SER:OG	1:A:127:THR:O	2.22	0.56
1:F:204:SER:HA	1:F:207:TRP:HD1	1.71	0.56
1:A:94:ILE:HG22	1:A:115:ASN:HB2	1.88	0.56
1:A:353:ILE:HG12	1:A:369:TYR:CE2	2.41	0.56
1:E:345:SER:HB3	1:E:348:LYS:HB2	1.86	0.56
1:B:410:VAL:HB	1:B:537:VAL:HG21	1.88	0.56
1:C:145:ASN:HD22	1:C:263:ILE:HD13	1.70	0.56
1:D:266:ARG:HA	1:D:269:THR:HG22	1.87	0.56
1:D:353:ILE:HG12	1:D:369:TYR:CE2	2.41	0.55
1:D:196:PHE:O	1:D:200:ILE:HG13	2.07	0.55
1:A:96:TYR:HD2	1:A:113:LEU:HD12	1.71	0.55
1:F:198:PRO:O	1:F:241:ASN:N	2.40	0.55
1:B:61:TYR:HD2	1:B:520:VAL:HG13	1.71	0.54
1:C:117:LEU:HD21	1:C:510:ILE:HG21	1.89	0.54
1:A:427:THR:HA	1:A:430:LYS:HD3	1.89	0.54
1:C:96:TYR:HD2	1:C:113:LEU:HD12	1.72	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:150:CYS:SG	1:B:281:ARG:NH2	2.81	0.53
1:A:195:ASN:HB3	1:A:198:PRO:HD2	1.91	0.53
1:B:44:LEU:HB2	1:B:127:THR:HG21	1.90	0.53
1:A:269:THR:HG21	1:A:277:TYR:HE2	1.74	0.53
1:D:374:LEU:HG	1:D:534:PRO:HG3	1.91	0.53
1:C:374:LEU:HG	1:C:534:PRO:HG3	1.90	0.53
1:A:231:HIS:HA	1:A:234:MET:HG3	1.91	0.53
1:B:49:ARG:HG2	1:B:51:ASP:H	1.74	0.53
1:E:117:LEU:HD21	1:E:510:ILE:HG21	1.91	0.53
1:E:292:SER:HB3	1:E:295:ASP:HB2	1.91	0.53
1:F:244:PRO:HG2	1:F:250:MET:HG3	1.91	0.53
1:B:390:VAL:HG11	1:B:422:LEU:HB2	1.91	0.52
1:A:357:SER:HB2	1:A:404:ASN:ND2	2.24	0.52
1:B:401:MET:HB3	1:B:441:LEU:HD21	1.91	0.52
1:B:460:LYS:HG3	1:B:530:ASP:HA	1.92	0.52
1:B:279:VAL:HG22	1:B:288:ILE:HG12	1.92	0.52
1:B:152:SER:OG	1:B:255:GLU:OE2	2.25	0.51
1:B:188:MET:HB2	1:B:325:ILE:HG12	1.91	0.51
1:C:390:VAL:HG11	1:C:422:LEU:HB2	1.92	0.51
1:A:495:ILE:O	1:A:507:LYS:NZ	2.43	0.51
1:B:101:SER:HA	1:B:108:TRP:HA	1.93	0.51
1:D:413:ALA:HB2	1:D:532:ILE:HB	1.92	0.51
1:E:146:ASP:OD1	1:E:146:ASP:N	2.43	0.51
1:F:417:LYS:O	1:F:420:SER:OG	2.26	0.50
1:E:265:LEU:HD22	1:E:303:LEU:HD12	1.93	0.50
1:D:134:ILE:HG21	1:D:224:PRO:HG3	1.93	0.50
1:B:56:LEU:HD12	1:B:113:LEU:HD21	1.94	0.50
1:A:85:THR:H	1:A:89:LEU:HD23	1.76	0.50
1:B:440:LYS:NZ	1:E:182:ALA:O	2.32	0.50
1:C:38:LEU:HB3	1:C:42:GLU:HG3	1.93	0.50
1:D:48:LEU:HD13	1:D:69:ASN:HA	1.92	0.50
1:F:279:VAL:HG22	1:F:288:ILE:HG23	1.94	0.50
1:C:468:SER:O	1:C:470:LYS:N	2.45	0.50
1:E:476:ARG:NH2	1:E:517:THR:O	2.45	0.50
1:B:188:MET:HE3	1:B:322:TYR:HD2	1.77	0.49
1:C:413:ALA:HA	1:C:532:ILE:HD13	1.95	0.49
1:D:84:PRO:HG3	1:D:121:TYR:CE2	2.47	0.49
1:D:188:MET:HB2	1:D:325:ILE:HG12	1.94	0.49
1:C:269:THR:HG21	1:C:277:TYR:OH	2.13	0.49
1:A:374:LEU:HG	1:A:534:PRO:HG3	1.94	0.48
1:B:48:LEU:HD21	1:B:120:VAL:HG12	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:353:ILE:HD13	1:E:370:LEU:HD23	1.96	0.48
1:A:357:SER:HB2	1:A:404:ASN:HD21	1.78	0.48
1:D:353:ILE:HG12	1:D:369:TYR:CD2	2.48	0.48
1:C:524:LYS:O	1:C:528:MET:HG3	2.13	0.48
1:F:101:SER:HB2	1:F:362:ASN:HD21	1.79	0.48
1:B:257:ILE:O	1:B:261:ILE:HG13	2.14	0.48
1:D:48:LEU:HD11	1:D:72:ASN:HD22	1.78	0.48
1:D:253:PHE:HA	1:D:256:LEU:HD12	1.96	0.48
1:C:206:GLU:HA	1:C:225:GLY:HA3	1.96	0.48
1:D:32:MET:HG3	1:D:33:LYS:H	1.78	0.48
1:A:99:LEU:HD21	1:A:110:ARG:HG2	1.96	0.48
1:E:456:ARG:NH2	1:E:536:GLU:OE2	2.47	0.48
1:C:281:ARG:HD3	1:C:286:TYR:CE1	2.49	0.47
1:E:244:PRO:HG2	1:E:250:MET:HG3	1.96	0.47
1:A:434:LEU:HB3	1:A:454:PHE:CZ	2.49	0.47
1:B:415:PHE:CZ	1:B:437:LEU:HB2	2.49	0.47
1:A:349:MET:O	1:A:353:ILE:HG13	2.14	0.47
1:C:111:ILE:HG23	1:C:246:GLY:HA3	1.97	0.47
1:C:101:SER:HA	1:C:108:TRP:HA	1.96	0.47
1:A:401:MET:SD	1:A:411:GLY:HA3	2.56	0.46
1:F:303:LEU:HD22	1:F:314:LEU:HD21	1.96	0.46
1:A:353:ILE:HG21	1:A:370:LEU:HG	1.97	0.46
1:E:353:ILE:HG12	1:E:369:TYR:CE2	2.50	0.46
1:A:103:ASP:OD1	1:A:104:GLY:N	2.48	0.46
1:A:410:VAL:O	1:A:414:ILE:HG12	2.16	0.46
1:A:410:VAL:CG1	1:A:537:VAL:HG11	2.46	0.46
1:C:196:PHE:O	1:C:200:ILE:HG13	2.16	0.46
1:D:67:LEU:HA	1:D:517:THR:HG21	1.96	0.46
1:C:506:ASN:OD1	1:C:507:LYS:N	2.46	0.46
1:D:71:ILE:O	1:D:75:ILE:HG12	2.15	0.46
1:A:208:VAL:HG21	1:A:257:ILE:HG22	1.97	0.46
1:D:35:VAL:HG11	1:D:154:PRO:HG2	1.98	0.46
1:D:447:THR:O	1:D:449:MET:N	2.49	0.46
1:F:196:PHE:O	1:F:200:ILE:HG13	2.15	0.46
1:D:103:ASP:OD1	1:D:104:GLY:N	2.45	0.46
1:C:345:SER:HB3	1:C:348:LYS:HB2	1.97	0.46
1:D:227:LEU:O	1:D:231:HIS:ND1	2.49	0.46
1:E:188:MET:HB2	1:E:325:ILE:HG12	1.98	0.46
1:E:410:VAL:HB	1:E:537:VAL:HG11	1.98	0.46
1:E:389:GLN:HA	1:F:176:GLN:NE2	2.31	0.45
1:C:374:LEU:HD13	1:C:414:ILE:HG22	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:67:LEU:HA	1:E:517:THR:HG21	1.99	0.45
1:F:101:SER:HB3	1:F:405:PRO:HB2	1.98	0.45
1:B:354:TYR:OH	1:E:324:ASP:OD1	2.20	0.45
1:E:524:LYS:O	1:E:528:MET:HG3	2.17	0.45
1:F:390:VAL:HG11	1:F:422:LEU:HB2	1.97	0.45
1:E:146:ASP:OD1	1:E:266:ARG:NH1	2.32	0.45
1:A:341:LYS:HB3	1:A:343:HIS:CE1	2.52	0.45
1:D:191:THR:HB	1:D:314:LEU:HD22	1.97	0.45
1:E:269:THR:HG21	1:E:277:TYR:CE1	2.51	0.45
1:C:266:ARG:HA	1:C:269:THR:HG22	1.99	0.45
1:C:482:THR:HG23	1:C:484:GLU:HG3	1.99	0.45
1:D:410:VAL:O	1:D:414:ILE:HG23	2.17	0.45
1:E:193:ILE:H	1:E:284:ASP:HB2	1.81	0.45
1:B:103:ASP:H	1:B:362:ASN:ND2	2.14	0.45
1:B:348:LYS:HG2	1:E:338:ILE:HD12	1.99	0.45
1:A:297:ASP:O	1:A:301:LYS:HG3	2.17	0.45
1:B:189:PHE:HB2	1:B:296:LEU:HD23	1.99	0.44
1:C:48:LEU:HD22	1:C:68:LEU:HB3	1.99	0.44
1:C:296:LEU:HD12	1:C:296:LEU:HA	1.87	0.44
1:E:303:LEU:HA	1:E:306:VAL:HG12	1.98	0.44
1:C:354:TYR:CZ	1:D:326:ILE:HB	2.52	0.44
1:E:196:PHE:O	1:E:200:ILE:HG13	2.17	0.44
1:E:401:MET:HB3	1:E:441:LEU:HD21	1.98	0.44
1:B:379:LYS:HB2	1:B:379:LYS:HE3	1.84	0.44
1:E:390:VAL:HG11	1:E:422:LEU:HB2	1.99	0.44
1:A:101:SER:HA	1:A:108:TRP:CD1	2.51	0.44
1:B:48:LEU:HD13	1:B:69:ASN:HA	1.98	0.44
1:D:257:ILE:O	1:D:261:ILE:HG12	2.18	0.44
1:E:71:ILE:HA	1:E:513:LEU:HD21	1.99	0.44
1:F:171:TRP:CZ3	1:F:330:LEU:HD21	2.53	0.44
1:E:74:SER:HB3	1:E:80:ILE:HG12	2.00	0.44
1:A:255:GLU:HA	1:A:258:LEU:HB2	2.00	0.44
1:D:148:PHE:CD1	1:D:279:VAL:HB	2.52	0.44
1:F:460:LYS:HG3	1:F:530:ASP:HA	2.00	0.44
1:B:228:ILE:O	1:B:232:ILE:HG12	2.17	0.44
1:B:456:ARG:NH2	1:B:529:ASP:O	2.50	0.44
1:D:434:LEU:HB3	1:D:454:PHE:CZ	2.51	0.44
1:A:150:CYS:SG	1:A:281:ARG:NE	2.91	0.43
1:B:253:PHE:HA	1:B:256:LEU:HD12	2.00	0.43
1:D:353:ILE:HG21	1:D:370:LEU:CD2	2.48	0.43
1:E:71:ILE:O	1:E:75:ILE:HG13	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:269:THR:HG23	1:B:274:ILE:HB	1.99	0.43
1:F:353:ILE:HG12	1:F:369:TYR:CD2	2.53	0.43
1:A:449:MET:HA	1:A:452:ILE:HD12	2.00	0.43
1:C:191:THR:HG22	1:C:319:THR:HA	2.00	0.43
1:E:152:SER:HB2	1:E:283:ARG:HG2	1.99	0.43
1:C:486:THR:HG23	1:C:515:ARG:HD3	2.00	0.43
1:D:480:GLU:OE1	1:D:520:VAL:N	2.45	0.43
1:A:413:ALA:HB2	1:A:532:ILE:HB	2.01	0.43
1:B:346:LEU:HB3	1:B:373:PHE:HE1	1.84	0.43
1:D:95:ASN:OD1	1:D:115:ASN:N	2.40	0.43
1:A:103:ASP:HA	1:A:362:ASN:HD22	1.82	0.43
1:A:138:LYS:HE2	1:A:208:VAL:HG12	2.00	0.43
1:A:196:PHE:O	1:A:200:ILE:HG13	2.19	0.43
1:B:49:ARG:HD3	1:B:51:ASP:HB2	2.01	0.43
1:B:283:ARG:HD3	1:B:283:ARG:HA	1.90	0.43
1:A:84:PRO:HG3	1:A:121:TYR:CE2	2.54	0.43
1:F:71:ILE:HD11	1:F:117:LEU:HG	2.01	0.43
1:F:102:LYS:O	1:F:102:LYS:HG2	2.19	0.43
1:F:469:TYR:C	1:F:471:SER:H	2.22	0.43
1:C:397:ILE:HG12	1:C:414:ILE:HD11	2.01	0.43
1:D:208:VAL:HG12	1:D:260:GLN:HG2	2.01	0.42
1:F:209:PHE:CE2	1:F:224:PRO:HG2	2.53	0.42
1:B:331:LYS:H	1:B:331:LYS:HG2	1.41	0.42
1:B:349:MET:O	1:B:353:ILE:HG13	2.19	0.42
1:D:114:ILE:HG21	1:D:119:TYR:CD1	2.54	0.42
1:D:148:PHE:HD2	1:D:262:ASP:HB3	1.83	0.42
1:C:502:GLU:C	1:C:504:SER:H	2.23	0.42
1:E:394:LEU:HD21	1:E:419:LEU:HD21	2.01	0.42
1:B:116:PRO:O	1:B:120:VAL:HG22	2.19	0.42
1:C:188:MET:HB2	1:C:325:ILE:HG12	2.00	0.42
1:E:248:THR:HA	1:E:251:ASP:HB2	2.00	0.42
1:A:290:SER:HB3	1:A:296:LEU:CD1	2.49	0.42
1:B:171:TRP:HZ2	1:F:348:LYS:HG2	1.83	0.42
1:B:117:LEU:HD11	1:B:510:ILE:HG21	2.01	0.42
1:D:264:GLU:HB3	1:D:306:VAL:HG21	2.02	0.42
1:E:296:LEU:HD23	1:E:296:LEU:HA	1.91	0.42
1:A:385:ASP:HB3	1:A:389:GLN:HG3	2.01	0.41
1:B:415:PHE:CG	1:B:437:LEU:HD22	2.55	0.41
1:A:44:LEU:O	1:A:48:LEU:HG	2.20	0.41
1:D:118:TYR:CE1	1:D:235:MET:HB3	2.55	0.41
1:F:84:PRO:HB2	1:F:89:LEU:HD11	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:71:ILE:HG23	1:C:120:VAL:HG21	2.02	0.41
1:A:71:ILE:HG23	1:A:120:VAL:HG21	2.01	0.41
1:B:438:HIS:HB2	1:B:454:PHE:HE2	1.85	0.41
1:B:116:PRO:HD3	1:B:493:TRP:CE2	2.56	0.41
1:B:529:ASP:OD1	1:B:530:ASP:N	2.54	0.41
1:C:197:TYR:O	1:C:243:ILE:HG22	2.20	0.41
1:F:462:ASN:OD1	1:F:463:LEU:N	2.54	0.41
1:E:76:HIS:HB3	1:E:124:ARG:NH1	2.35	0.41
1:E:95:ASN:OD1	1:E:115:ASN:N	2.49	0.41
1:F:84:PRO:HB3	1:F:497:TRP:CZ3	2.55	0.41
1:F:181:LEU:HD23	1:F:181:LEU:HA	1.95	0.41
1:F:205:PHE:O	1:F:208:VAL:HG22	2.21	0.41
1:A:419:LEU:HD23	1:A:419:LEU:HA	1.92	0.41
1:E:409:PRO:HG3	1:E:453:TRP:NE1	2.35	0.41
1:C:449:MET:HA	1:C:452:ILE:HD12	2.01	0.41
1:E:35:VAL:HG11	1:E:154:PRO:HG2	2.03	0.41
1:E:55:THR:N	1:E:246:GLY:O	2.52	0.41
1:F:103:ASP:HA	1:F:362:ASN:HD22	1.85	0.41
1:B:46:TYR:CE2	1:B:248:THR:HG21	2.56	0.41
1:B:415:PHE:CE2	1:B:437:LEU:HD13	2.56	0.41
1:C:85:THR:HG22	1:C:87:LYS:H	1.85	0.41
1:D:290:SER:HB3	1:D:296:LEU:HD21	2.03	0.41
1:E:260:GLN:HA	1:E:263:ILE:HG22	2.02	0.41
1:F:494:ASN:OD1	1:F:495:ILE:N	2.54	0.41
1:A:71:ILE:HG13	1:A:80:ILE:HD11	2.02	0.41
1:A:362:ASN:ND2	1:A:405:PRO:HD2	2.35	0.41
1:C:457:THR:OG1	1:C:532:ILE:HD11	2.21	0.41
1:C:532:ILE:H	1:C:532:ILE:HD12	1.86	0.41
1:D:394:LEU:HD12	1:D:415:PHE:CZ	2.56	0.40
1:E:495:ILE:HD13	1:E:511:LEU:HD21	2.03	0.40
1:B:396:ILE:HG12	1:E:326:ILE:HD12	2.02	0.40
1:C:71:ILE:HD11	1:C:117:LEU:HG	2.03	0.40
1:D:349:MET:O	1:D:353:ILE:HG13	2.21	0.40
1:F:463:LEU:HD23	1:F:463:LEU:HA	1.86	0.40
1:A:101:SER:O	1:A:103:ASP:N	2.52	0.40
1:B:274:ILE:HG21	1:B:299:ILE:HD11	2.02	0.40
1:E:85:THR:HG23	1:E:88:GLU:H	1.86	0.40
1:E:304:VAL:HG22	1:E:314:LEU:HD13	2.04	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	479/541 (88%)	459 (96%)	20 (4%)	0	100	100
1	B	468/541 (86%)	449 (96%)	19 (4%)	0	100	100
1	C	472/541 (87%)	452 (96%)	20 (4%)	0	100	100
1	D	467/541 (86%)	441 (94%)	26 (6%)	0	100	100
1	E	476/541 (88%)	456 (96%)	20 (4%)	0	100	100
1	F	465/541 (86%)	447 (96%)	17 (4%)	1 (0%)	47	79
All	All	2827/3246 (87%)	2704 (96%)	122 (4%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	F	103	ASP

### 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	398/506 (79%)	397 (100%)	1 (0%)	92	96
1	B	389/506 (77%)	388 (100%)	1 (0%)	92	96
1	C	371/506 (73%)	370 (100%)	1 (0%)	92	96
1	D	357/506 (71%)	357 (100%)	0	100	100
1	E	379/506 (75%)	379 (100%)	0	100	100
1	F	373/506 (74%)	373 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	2267/3036 (75%)	2264 (100%)	3 (0%)	93 98

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	342	ARG
1	B	456	ARG
1	C	456	ARG

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](#)

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	487/541 (90%)	-0.19	3 (0%) 89 78	67, 103, 146, 206	0
1	B	482/541 (89%)	-0.16	4 (0%) 86 72	62, 107, 152, 193	0
1	C	484/541 (89%)	-0.22	2 (0%) 92 84	62, 103, 141, 204	0
1	D	479/541 (88%)	-0.17	3 (0%) 89 78	63, 108, 157, 204	0
1	E	488/541 (90%)	-0.20	1 (0%) 95 90	60, 96, 146, 208	0
1	F	479/541 (88%)	-0.11	4 (0%) 86 72	63, 111, 166, 236	0
All	All	2899/3246 (89%)	-0.18	17 (0%) 89 78	60, 104, 154, 236	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	241	ASN	3.8
1	D	242	GLY	3.5
1	D	386	ASN	3.5
1	F	386	ASN	3.2
1	B	387	GLY	3.2
1	B	386	ASN	2.7
1	C	506	ASN	2.6
1	B	510	ILE	2.6
1	D	517	THR	2.5
1	E	503	THR	2.4
1	A	194	SER	2.4
1	C	241	ASN	2.3
1	A	241	ASN	2.2
1	F	242	GLY	2.2
1	B	465	TRP	2.2
1	A	386	ASN	2.2
1	F	487	PHE	2.1

## 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](#)

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.