



# Full wwPDB X-ray Structure Validation Report ⓘ

Apr 17, 2014 – 04:18 AM EDT

PDB ID : 4JY1  
Title : CRYSTAL STRUCTURE OF HCV NS5B POLYMERASE IN COMPLEX  
WITH COMPOUND 5  
Authors : Coulombe, R.  
Deposited on : 2013-03-28  
Resolution : 2.60 Å(reported)

This is a full wwPDB 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 <http://wwpdb.org/ValidationPDFNotes.html>

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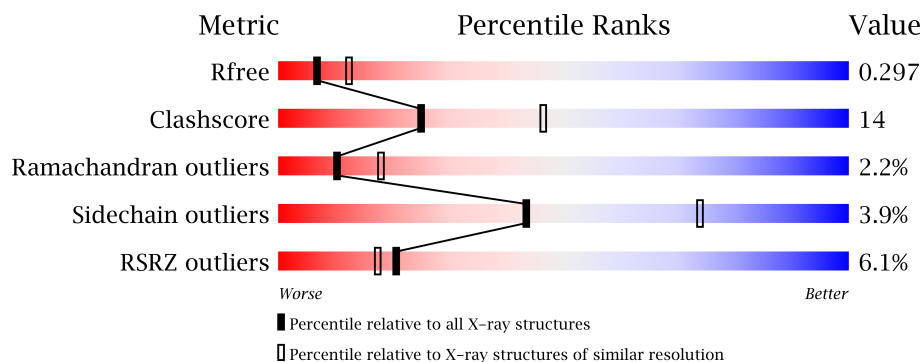
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.16 November 2013  
Xtriage (Phenix) : dev-1439  
EDS : stable22978  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22978

# 1 Overall quality at a glance

The reported resolution of this entry is 2.60 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))
$R_{free}$	66092	1718 (2.60-2.60)
Clashscore	79885	2154 (2.60-2.60)
Ramachandran outliers	78287	2113 (2.60-2.60)
Sidechain outliers	78261	2113 (2.60-2.60)
RSRZ outliers	66119	1718 (2.60-2.60)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	576	
1	B	576	

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8731 atoms, of which 0 are hydrogen and 0 are deuterium.

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.

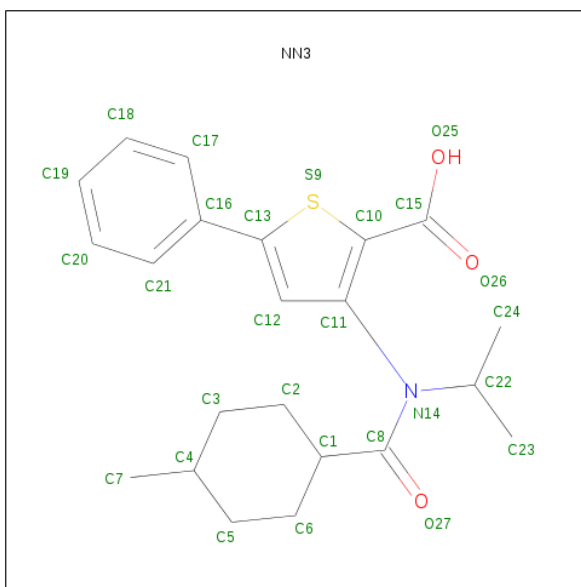
- Molecule 1 is a protein called Genome polyprotein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	559	Total	C	N	O	S	0	0	0
			4358	2745	770	811	32			
1	B	558	Total	C	N	O	S	0	0	0
			4346	2737	768	809	32			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	571	HIS	-	EXPRESSION TAG	UNP O92972
A	572	HIS	-	EXPRESSION TAG	UNP O92972
A	573	HIS	-	EXPRESSION TAG	UNP O92972
A	574	HIS	-	EXPRESSION TAG	UNP O92972
A	575	HIS	-	EXPRESSION TAG	UNP O92972
A	576	HIS	-	EXPRESSION TAG	UNP O92972
B	571	HIS	-	EXPRESSION TAG	UNP O92972
B	572	HIS	-	EXPRESSION TAG	UNP O92972
B	573	HIS	-	EXPRESSION TAG	UNP O92972
B	574	HIS	-	EXPRESSION TAG	UNP O92972
B	575	HIS	-	EXPRESSION TAG	UNP O92972
B	576	HIS	-	EXPRESSION TAG	UNP O92972

- Molecule 2 is 3-{ISOPROPYL[(TRANS-4-METHYLCYCLOHEXYL)CARBONYL]AMINO}-5-PHENYLTHIOPHENE-2-CARBOXYLICACID (three-letter code: NN3) (formula: C<sub>22</sub>H<sub>27</sub>NO<sub>3</sub>S).



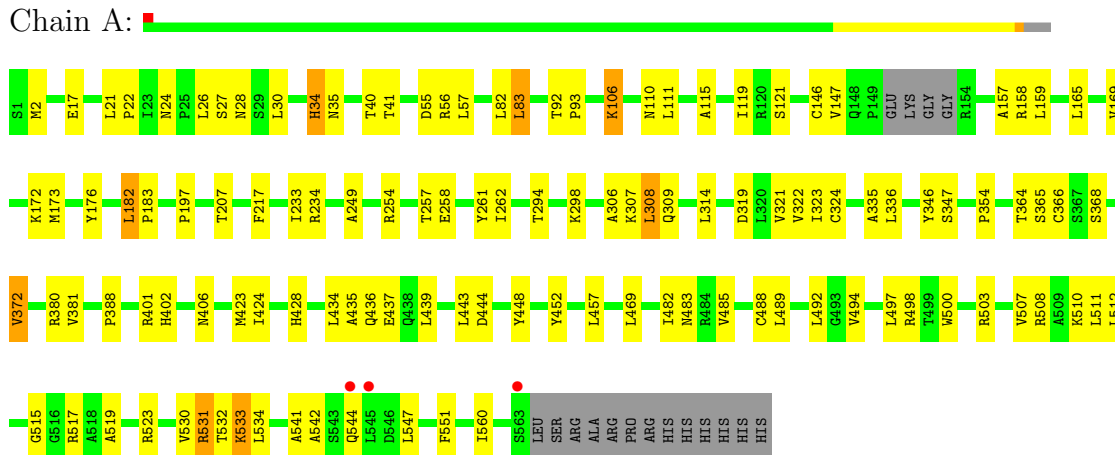
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			27	22	1	3	1		

### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

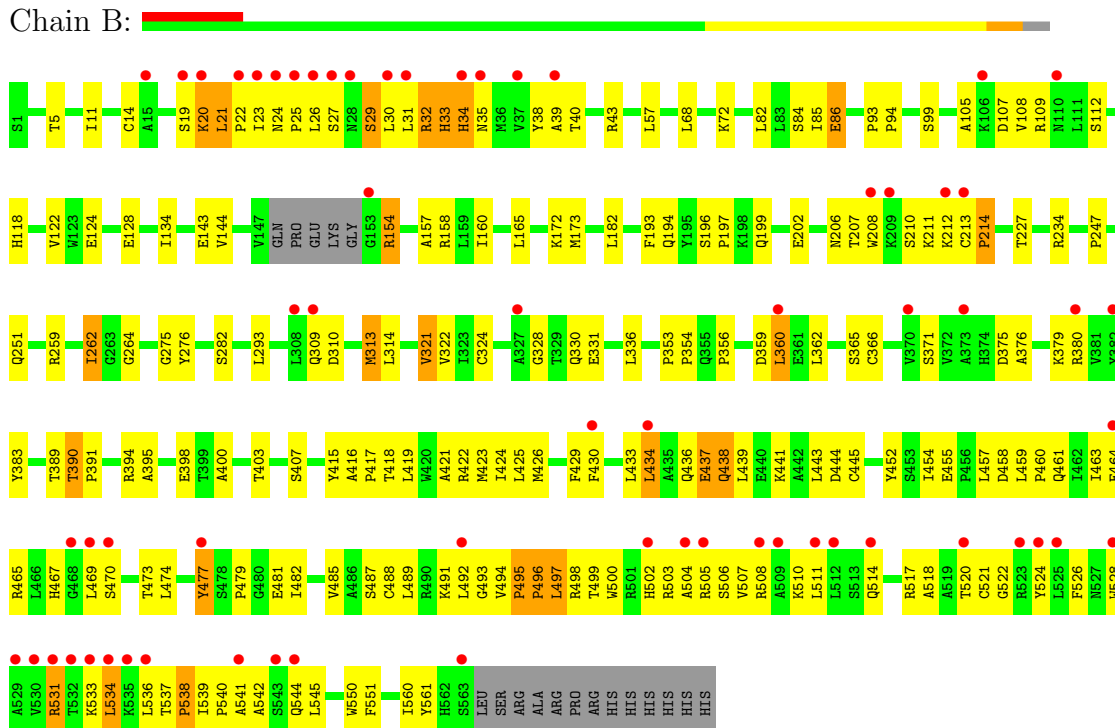
#### • Molecule 1: Genome polypeptide

Chain A:



#### • Molecule 1: Genome polypeptide

Chain B:



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	106.26Å 108.03Å 135.13Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.00 – 2.60 39.21 – 2.60	Depositor EDS
% Data completeness (in resolution range)	99.6 (40.00-2.60) 99.6 (39.21-2.60)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.92 (at 2.61Å)	Xtriage
Refinement program	CNX 2002	Depositor
R, $R_{free}$	0.252 , 0.299 0.251 , 0.297	Depositor DCC
$R_{free}$ test set	4875 reflections (11.21%)	DCC
Wilson B-factor (Å <sup>2</sup> )	48.0	Xtriage
Anisotropy	0.125	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 21.2	EDS
Estimated twinning fraction	0.018 for k,h,-l	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	1 of 48350 reflections (0.002%)	Xtriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	8731	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.38% of the height of the origin peak. No significant pseudotranslation is detected.*

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

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NN3

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	0.39	0/4453	0.65	0/6044
1	B	0.36	0/4440	0.65	1/6025 (0.0%)
All	All	0.37	0/8893	0.65	1/12069 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	29	SER	N-CA-C	-5.22	96.91	111.00

There are no chirality outliers.

There are no planarity outliers.

### 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4358	0	4369	82	0
1	B	4346	0	4357	161	0
2	A	27	0	26	0	0
All	All	8731	0	8752	241	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including

hydrogens) of the entry. The overall clashscore for this entry is 14.

All (241) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:531:ARG:H	1:B:531:ARG:HD3	1.28	0.98
1:B:461:GLN:HB3	1:B:545:LEU:HD11	1.60	0.84
1:B:31:LEU:HB3	1:B:494:VAL:HG22	1.61	0.83
1:A:485:VAL:O	1:A:489:LEU:HG	1.80	0.81
1:B:31:LEU:HD12	1:B:31:LEU:O	1.81	0.79
1:B:520:THR:HG21	1:B:539:ILE:HD12	1.63	0.79
1:A:40:THR:HB	1:A:157:ALA:HB2	1.65	0.78
1:B:30:LEU:O	1:B:494:VAL:HG22	1.84	0.77
1:A:314:LEU:HB3	1:A:321:VAL:HG12	1.64	0.76
1:B:489:LEU:HD22	1:B:494:VAL:HB	1.66	0.76
1:B:30:LEU:O	1:B:494:VAL:HG13	1.86	0.75
1:B:531:ARG:CD	1:B:531:ARG:H	2.00	0.75
1:B:419:LEU:HD13	1:B:477:TYR:CE1	2.23	0.74
1:B:309:GLN:O	1:B:324:CYS:HB2	1.89	0.73
1:B:531:ARG:HD3	1:B:531:ARG:N	2.04	0.72
1:B:314:LEU:HB3	1:B:321:VAL:HG13	1.72	0.72
1:B:336:LEU:HD23	1:B:356:PRO:HD3	1.72	0.71
1:A:82:LEU:HD13	1:A:249:ALA:HB2	1.73	0.71
1:B:29:SER:O	1:B:30:LEU:HB3	1.90	0.70
1:A:30:LEU:O	1:A:494:VAL:HG13	1.93	0.69
1:A:308:LEU:HD13	1:A:335:ALA:HB1	1.75	0.69
1:B:416:ALA:HB3	1:B:417:PRO:HD3	1.76	0.68
1:B:459:LEU:O	1:B:463:ILE:HG13	1.95	0.66
1:A:423:MET:HE1	1:A:497:LEU:HB2	1.78	0.64
1:A:34:HIS:HD2	1:A:35:ASN:H	1.44	0.64
1:B:489:LEU:CD2	1:B:494:VAL:HB	2.28	0.64
1:A:406:ASN:ND2	1:A:443:LEU:HB3	2.13	0.63
1:B:434:LEU:HG	1:B:439:LEU:HD11	1.79	0.63
1:B:436:GLN:O	1:B:437:GLU:HB3	1.98	0.63
1:B:21:LEU:HG	1:B:22:PRO:HD2	1.82	0.62
1:B:27:SER:C	1:B:29:SER:H	2.02	0.62
1:B:22:PRO:O	1:B:24:ASN:N	2.32	0.62
1:B:208:TRP:CE3	1:B:360:LEU:HD13	2.35	0.62
1:B:376:ALA:HB2	1:B:473:THR:HB	1.80	0.61
1:B:93:PRO:HG3	1:B:561:TYR:HB2	1.81	0.61
1:A:503:ARG:O	1:A:507:VAL:HG23	2.01	0.61
1:A:423:MET:CE	1:A:497:LEU:HD13	2.30	0.61
1:A:233:ILE:HD13	1:A:261:TYR:O	2.00	0.61
1:A:309:GLN:O	1:A:324:CYS:HB2	2.01	0.60
1:B:85:ILE:HD13	1:B:173:MET:SD	2.40	0.60

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:30:LEU:HB2	1:A:428:HIS:CE1	2.35	0.60
1:B:479:PRO:O	1:B:482:ILE:HG22	2.01	0.60
1:A:119:ILE:HD13	1:A:169:VAL:HG11	1.82	0.60
1:A:423:MET:HE2	1:A:497:LEU:HD13	1.83	0.59
1:B:124:GLU:O	1:B:128:GLU:HG2	2.01	0.59
1:B:30:LEU:HA	1:B:500:TRP:HZ2	1.66	0.59
1:B:40:THR:HB	1:B:157:ALA:HB2	1.83	0.59
1:B:505:ARG:HD3	1:B:531:ARG:HH12	1.66	0.59
1:B:196:SER:H	1:B:199:GLN:HB2	1.67	0.59
1:B:520:THR:HG21	1:B:539:ILE:CD1	2.32	0.59
1:B:375:ASP:CB	1:B:379:LYS:H	2.16	0.58
1:A:24:ASN:HB3	1:A:27:SER:OG	2.03	0.58
1:B:359:ASP:HB3	1:B:362:LEU:CD1	2.34	0.58
1:B:470:SER:HB2	1:B:474:LEU:HD21	1.85	0.58
1:B:434:LEU:CD1	1:B:507:VAL:HG13	2.34	0.58
1:A:257:THR:O	1:A:262:ILE:HG23	2.04	0.58
1:B:234:ARG:HG3	1:B:262:ILE:HD11	1.85	0.58
1:B:211:LYS:HB2	1:B:214:PRO:HB3	1.85	0.57
1:B:84:SER:OG	1:B:86:GLU:HG2	2.05	0.56
1:B:488:CYS:HA	1:B:491:LYS:HD3	1.87	0.56
1:B:439:LEU:HB3	1:B:457:LEU:HD21	1.88	0.56
1:B:86:GLU:CD	1:B:86:GLU:H	2.08	0.56
1:A:83:LEU:HD22	1:A:176:TYR:HB3	1.87	0.56
1:A:533:LYS:H	1:A:533:LYS:HD2	1.71	0.56
1:A:22:PRO:HG3	1:A:401:ARG:NH2	2.20	0.56
1:B:469:LEU:HD11	1:B:538:PRO:HG3	1.88	0.56
1:B:383:TYR:CE2	1:B:418:THR:HA	2.41	0.56
1:B:505:ARG:HD3	1:B:531:ARG:NH1	2.22	0.55
1:A:34:HIS:CD2	1:A:35:ASN:N	2.75	0.55
1:B:313:MET:SD	1:B:322:VAL:HG22	2.47	0.54
1:B:423:MET:HE3	1:B:500:TRP:HD1	1.72	0.54
1:B:20:LYS:O	1:B:21:LEU:HB2	2.07	0.54
1:B:359:ASP:HB3	1:B:362:LEU:HD12	1.89	0.54
1:B:380:ARG:HB3	1:B:380:ARG:HH11	1.71	0.54
1:A:34:HIS:HD2	1:A:35:ASN:N	2.06	0.54
1:A:541:ALA:O	1:A:544:GLN:HB3	2.08	0.54
1:B:437:GLU:HG2	1:B:438:GLN:N	2.20	0.54
1:B:434:LEU:HD12	1:B:507:VAL:HG13	1.89	0.54
1:A:336:LEU:HD11	1:A:354:PRO:HG2	1.89	0.53
1:A:531:ARG:HE	1:A:531:ARG:H	1.55	0.53
1:B:331:GLU:CD	1:B:331:GLU:H	2.12	0.53
1:B:197:PRO:HB2	1:B:467:HIS:HE1	1.73	0.53

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:390:THR:HB	1:B:391:PRO:HD3	1.90	0.53
1:A:336:LEU:HD11	1:A:354:PRO:HB2	1.91	0.53
1:A:457:LEU:HD12	1:A:517:ARG:HH21	1.74	0.52
1:A:182:LEU:HB3	1:A:183:PRO:HD3	1.89	0.52
1:B:526:PHE:HA	1:B:528:TRP:CD1	2.45	0.52
1:B:520:THR:CG2	1:B:539:ILE:HD12	2.39	0.52
1:B:107:ASP:HB3	1:B:112:SER:OG	2.09	0.52
1:B:160:ILE:HG13	1:B:160:ILE:O	2.08	0.52
1:A:388:PRO:HG2	1:A:488:CYS:SG	2.50	0.52
1:A:515:GLY:O	1:A:519:ALA:HB2	2.10	0.51
1:A:531:ARG:O	1:A:533:LYS:HD2	2.11	0.51
1:B:172:LYS:HE3	1:B:560:ILE:HD13	1.91	0.51
1:A:531:ARG:H	1:A:531:ARG:NE	2.08	0.51
1:A:533:LYS:H	1:A:533:LYS:CD	2.23	0.51
1:B:383:TYR:HE2	1:B:418:THR:HA	1.75	0.51
1:A:57:LEU:C	1:A:57:LEU:HD23	2.30	0.51
1:B:197:PRO:HB2	1:B:467:HIS:CE1	2.45	0.51
1:B:33:HIS:HB2	1:B:492:LEU:O	2.11	0.51
1:B:487:SER:O	1:B:491:LYS:HG3	2.11	0.51
1:B:105:ALA:O	1:B:109:ARG:HG3	2.11	0.51
1:B:371:SER:HB3	1:B:383:TYR:CE1	2.46	0.51
1:B:433:LEU:HB3	1:B:438:GLN:O	2.11	0.51
1:A:2:MET:SD	1:A:55:ASP:HB2	2.51	0.50
1:B:375:ASP:HB2	1:B:379:LYS:HB2	1.93	0.50
1:B:526:PHE:HA	1:B:528:TRP:NE1	2.26	0.50
1:B:506:SER:OG	1:B:510:LYS:HE3	2.12	0.50
1:A:56:ARG:HD3	1:A:56:ARG:N	2.26	0.50
1:B:455:GLU:HB2	1:B:458:ASP:OD2	2.12	0.50
1:A:24:ASN:OD1	1:A:26:LEU:N	2.45	0.49
1:B:11:ILE:O	1:B:11:ILE:HG22	2.12	0.49
1:B:26:LEU:O	1:B:27:SER:HB3	2.12	0.49
1:B:395:ALA:HB1	1:B:429:PHE:HZ	1.76	0.49
1:A:106:LYS:HA	1:A:106:LYS:HE2	1.95	0.49
1:B:496:PRO:O	1:B:498:ARG:N	2.43	0.49
1:B:202:GLU:O	1:B:206:ASN:ND2	2.46	0.49
1:A:234:ARG:HD3	1:B:247:PRO:HG3	1.93	0.49
1:B:365:SER:O	1:B:366:CYS:HB2	2.12	0.49
1:A:444:ASP:HA	1:A:452:TYR:O	2.12	0.49
1:B:144:VAL:HB	1:B:394:ARG:HG2	1.95	0.49
1:B:511:LEU:HD13	1:B:521:CYS:HB2	1.93	0.49
1:B:118:HIS:O	1:B:122:VAL:HG23	2.13	0.48
1:B:202:GLU:HG2	1:B:206:ASN:HD21	1.77	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:33:HIS:C	1:B:35:ASN:H	2.16	0.48
1:A:510:LYS:NZ	1:A:510:LYS:HB2	2.27	0.48
1:A:254:ARG:HG2	1:B:251:GLN:NE2	2.29	0.48
1:B:445:CYS:O	1:B:452:TYR:HB2	2.12	0.48
1:B:444:ASP:HA	1:B:452:TYR:O	2.13	0.48
1:B:464:GLU:OE1	1:B:539:ILE:HG22	2.12	0.48
1:A:531:ARG:HG2	1:A:532:THR:N	2.28	0.48
1:B:134:ILE:HG13	1:B:259:ARG:HB3	1.96	0.48
1:B:328:GLY:HA3	1:B:331:GLU:OE2	2.14	0.48
1:B:27:SER:C	1:B:29:SER:N	2.66	0.48
1:B:506:SER:O	1:B:510:LYS:HG3	2.13	0.48
1:A:336:LEU:CD1	1:A:354:PRO:HG2	2.43	0.48
1:A:482:ILE:HG13	1:A:483:ASN:N	2.28	0.48
1:A:294:THR:CG2	1:A:298:LYS:HE3	2.44	0.47
1:A:372:VAL:HG11	1:A:380:ARG:NH2	2.28	0.47
1:A:542:ALA:C	1:A:544:GLN:H	2.16	0.47
1:A:364:THR:HA	1:A:368:SER:O	2.14	0.47
1:A:523:ARG:HG3	1:A:534:LEU:HD12	1.95	0.47
1:B:383:TYR:HH	1:B:477:TYR:HD2	1.63	0.47
1:B:68:LEU:O	1:B:72:LYS:HG3	2.14	0.47
1:B:202:GLU:CG	1:B:206:ASN:HD21	2.28	0.47
1:B:517:ARG:HH11	1:B:517:ARG:HG3	1.80	0.47
1:B:202:GLU:HG2	1:B:206:ASN:ND2	2.30	0.47
1:A:336:LEU:HD11	1:A:354:PRO:CG	2.45	0.47
1:B:383:TYR:OH	1:B:481:GLU:HG2	2.15	0.46
1:B:160:ILE:HA	1:B:282:SER:OG	2.15	0.46
1:B:375:ASP:HB2	1:B:379:LYS:H	1.80	0.46
1:B:504:ALA:C	1:B:506:SER:N	2.68	0.46
1:B:94:PRO:HD3	1:B:561:TYR:CD2	2.51	0.46
1:A:314:LEU:HB3	1:A:321:VAL:CG1	2.39	0.46
1:B:31:LEU:HA	1:B:493:GLY:O	2.15	0.46
1:B:539:ILE:O	1:B:542:ALA:HB2	2.16	0.46
1:B:418:THR:OG1	1:B:421:ALA:HB2	2.16	0.46
1:A:115:ALA:O	1:A:119:ILE:HG13	2.16	0.45
1:A:83:LEU:HA	1:A:83:LEU:HD12	1.73	0.45
1:B:500:TRP:C	1:B:502:HIS:H	2.18	0.45
1:B:421:ALA:HA	1:B:425:LEU:HD12	1.99	0.45
1:B:30:LEU:O	1:B:494:VAL:CG2	2.61	0.45
1:B:144:VAL:HB	1:B:394:ARG:CG	2.47	0.45
1:A:439:LEU:O	1:A:457:LEU:HG	2.17	0.45
1:B:32:ARG:O	1:B:34:HIS:N	2.45	0.45
1:B:422:ARG:HA	1:B:426:MET:SD	2.57	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:522:GLY:HA2	1:B:526:PHE:HD2	1.82	0.45
1:B:5:THR:O	1:B:275:GLY:HA3	2.17	0.44
1:B:29:SER:O	1:B:30:LEU:CB	2.62	0.44
1:B:423:MET:CE	1:B:500:TRP:HD1	2.30	0.44
1:B:533:LYS:O	1:B:534:LEU:HB2	2.17	0.44
1:A:448:TYR:CE2	1:A:551:PHE:HD1	2.36	0.44
1:A:83:LEU:HB2	1:A:173:MET:HA	1.99	0.44
1:B:415:TYR:HB3	1:B:418:THR:HG21	1.99	0.44
1:A:17:GLU:OE1	1:A:41:THR:HB	2.18	0.44
1:B:540:PRO:O	1:B:541:ALA:HB3	2.17	0.44
1:A:217:PHE:CE1	1:A:322:VAL:HB	2.53	0.44
1:B:423:MET:CE	1:B:500:TRP:CD1	3.00	0.44
1:B:93:PRO:HA	1:B:94:PRO:HD3	1.85	0.44
1:A:172:LYS:HE3	1:A:560:ILE:HD13	2.00	0.43
1:A:336:LEU:HD11	1:A:354:PRO:CB	2.48	0.43
1:A:110:ASN:O	1:A:111:LEU:HB2	2.17	0.43
1:B:143:GLU:OE1	1:B:158:ARG:NH1	2.51	0.43
1:B:194:GLN:O	1:B:551:PHE:O	2.37	0.43
1:B:196:SER:O	1:B:197:PRO:C	2.57	0.43
1:B:353:PRO:HA	1:B:354:PRO:HD2	1.90	0.43
1:B:518:ALA:O	1:B:521:CYS:HB2	2.19	0.43
1:A:508:ARG:CZ	1:A:530:VAL:HG11	2.49	0.42
1:B:39:ALA:HA	1:B:143:GLU:O	2.18	0.42
1:B:196:SER:OG	1:B:199:GLN:HG3	2.18	0.42
1:A:434:LEU:C	1:A:436:GLN:H	2.23	0.42
1:B:172:LYS:HE3	1:B:560:ILE:CD1	2.50	0.42
1:B:524:TYR:CE2	1:B:536:LEU:HB3	2.53	0.42
1:B:465:ARG:NH1	1:B:545:LEU:HB2	2.34	0.42
1:B:22:PRO:HG2	1:B:400:ALA:HB1	2.01	0.42
1:B:439:LEU:O	1:B:457:LEU:HG	2.19	0.42
1:B:510:LYS:O	1:B:514:GLN:HG2	2.19	0.42
1:A:531:ARG:HE	1:A:531:ARG:N	2.17	0.42
1:B:29:SER:C	1:B:31:LEU:H	2.22	0.42
1:A:254:ARG:HH12	1:A:258:GLU:CG	2.32	0.42
1:B:398:GLU:OE2	1:B:407:SER:HB3	2.19	0.42
1:B:423:MET:HE1	1:B:424:ILE:HG12	2.02	0.42
1:B:19:SER:CB	1:B:43:ARG:HH22	2.33	0.42
1:A:207:THR:HG22	1:A:323:ILE:HG21	2.02	0.42
1:A:434:LEU:HD22	1:A:510:LYS:HD2	2.02	0.41
1:A:508:ARG:O	1:A:511:LEU:HB2	2.19	0.41
1:B:108:VAL:HG21	1:B:165:LEU:HD21	2.02	0.41
1:B:445:CYS:HB3	1:B:454:ILE:HD12	2.02	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:82:LEU:HD12	1:B:173:MET:O	2.21	0.41
1:A:92:THR:HA	1:A:93:PRO:HD3	1.98	0.41
1:A:346:TYR:O	1:A:347:SER:HB3	2.19	0.41
1:A:319:ASP:CG	1:A:366:CYS:H	2.23	0.41
1:B:419:LEU:HB2	1:B:477:TYR:CZ	2.55	0.41
1:B:418:THR:OG1	1:B:421:ALA:CB	2.68	0.41
1:B:541:ALA:HA	1:B:544:GLN:OE1	2.20	0.41
1:B:264:GLY:HA2	1:B:276:TYR:CZ	2.55	0.41
1:B:182:LEU:CD2	1:B:293:LEU:HD11	2.51	0.41
1:B:504:ALA:C	1:B:506:SER:H	2.24	0.41
1:A:21:LEU:HA	1:A:22:PRO:HD3	1.86	0.41
1:B:30:LEU:O	1:B:494:VAL:CG1	2.62	0.41
1:B:99:SER:HB2	1:B:165:LEU:HB3	2.02	0.41
1:B:213:CYS:HA	1:B:214:PRO:HD2	1.78	0.41
1:B:459:LEU:N	1:B:460:PRO:CD	2.83	0.41
1:B:389:THR:HG23	1:B:492:LEU:HD11	2.02	0.41
1:B:497:LEU:HG	1:B:497:LEU:O	2.21	0.41
1:B:537:THR:O	1:B:539:ILE:N	2.54	0.41
1:A:106:LYS:HA	1:A:106:LYS:CE	2.51	0.41
1:B:504:ALA:O	1:B:507:VAL:N	2.53	0.41
1:A:512:LEU:HA	1:A:519:ALA:HA	2.03	0.41
1:B:416:ALA:N	1:B:417:PRO:CD	2.84	0.41
1:B:436:GLN:O	1:B:437:GLU:CB	2.67	0.41
1:B:499:THR:O	1:B:503:ARG:HG3	2.21	0.41
1:B:207:THR:O	1:B:210:SER:OG	2.23	0.40
1:A:146:CYS:SG	1:A:492:LEU:HD11	2.61	0.40
1:A:165:LEU:O	1:A:169:VAL:HG23	2.21	0.40
1:A:306:ALA:O	1:A:307:LYS:HB2	2.21	0.40
1:A:423:MET:HE1	1:A:497:LEU:HD13	2.01	0.40
1:A:365:SER:O	1:A:366:CYS:HB2	2.21	0.40
1:A:424:ILE:HG23	1:A:500:TRP:CZ2	2.56	0.40
1:B:423:MET:CE	1:B:424:ILE:HG12	2.52	0.40
1:B:38:TYR:CE2	1:B:154:ARG:HG3	2.57	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	555/576 (96%)	526 (95%)	25 (4%)	4 (1%)	30	58
1	B	554/576 (96%)	482 (87%)	52 (9%)	20 (4%)	5	7
All	All	1109/1152 (96%)	1008 (91%)	77 (7%)	24 (2%)	10	18

All (24) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	23	ILE
1	B	441	LYS
1	B	495	PRO
1	A	469	LEU
1	B	25	PRO
1	B	193	PHE
1	B	438	GLN
1	B	477	TYR
1	B	534	LEU
1	A	437	GLU
1	B	33	HIS
1	B	34	HIS
1	B	154	ARG
1	B	360	LEU
1	B	497	LEU
1	B	20	LYS
1	B	212	LYS
1	B	330	GLN
1	B	538	PRO
1	A	435	ALA
1	B	21	LEU
1	A	147	VAL
1	B	214	PRO
1	B	496	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	477/491 (97%)	460 (96%)	17 (4%)	47	76
1	B	475/491 (97%)	455 (96%)	20 (4%)	40	71
All	All	952/982 (97%)	915 (96%)	37 (4%)	43	74

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	ASN
1	A	34	HIS
1	A	83	LEU
1	A	106	LYS
1	A	121	SER
1	A	158	ARG
1	A	159	LEU
1	A	182	LEU
1	A	197	PRO
1	A	308	LEU
1	A	372	VAL
1	A	381	VAL
1	A	402	HIS
1	A	498	ARG
1	A	531	ARG
1	A	533	LYS
1	A	547	LEU
1	B	14	CYS
1	B	32	ARG
1	B	57	LEU
1	B	86	GLU
1	B	227	THR
1	B	262	ILE
1	B	310	ASP
1	B	313	MET
1	B	321	VAL
1	B	390	THR
1	B	403	THR
1	B	430	PHE
1	B	434	LEU
1	B	437	GLU
1	B	443	LEU
1	B	485	VAL
1	B	495	PRO
1	B	508	ARG

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Mol	Chain	Res	Type
1	B	531	ARG
1	B	550	TRP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	28	ASN
1	A	34	HIS
1	A	251	GLN
1	A	273	ASN
1	A	309	GLN
1	A	406	ASN
1	A	428	HIS
1	A	502	HIS
1	A	514	GLN
1	A	544	GLN
1	A	562	HIS
1	B	24	ASN
1	B	206	ASN
1	B	251	GLN
1	B	273	ASN
1	B	374	HIS
1	B	461	GLN
1	B	467	HIS

### 5.3.3 RNA ⓘ

There are no RNA chains in this entry.

### 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

### 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.



## 5.6 Ligand geometry ⓘ

1 ligand is 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	NN3	A	601	-	29,29,29	2.08	5 (17%)	39,41,41	1.54	5 (12%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NN3	A	601	-	-	0/17/34/34	0/3/3/3

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	601	NN3	C8-N14	6.58	1.44	1.36
2	A	601	NN3	C12-C11	5.19	1.44	1.39
2	A	601	NN3	C1-C8	2.56	1.55	1.51
2	A	601	NN3	C22-N14	2.12	1.52	1.49
2	A	601	NN3	C12-C13	2.03	1.40	1.37

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	601	NN3	C1-C8-N14	4.91	123.99	119.20
2	A	601	NN3	C11-N14-C8	3.86	124.12	122.25
2	A	601	NN3	O27-C8-C1	-3.45	115.37	120.82
2	A	601	NN3	C11-N14-C22	-3.31	116.44	119.23
2	A	601	NN3	C6-C1-C8	3.15	115.52	109.60

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 5.7 Other polymers ⓘ

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, 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	559/576 (97%)	-0.11	3 (0%) 88 90	21, 38, 64, 89	0
1	B	558/576 (96%)	0.57	64 (11%) 5 4	24, 57, 92, 106	0
All	All	1117/1152 (96%)	0.23	67 (5%) 21 18	21, 44, 88, 106	0

All (67) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	26	LEU	6.8
1	B	532	THR	6.0
1	B	213	CYS	5.7
1	B	25	PRO	5.5
1	B	511	LEU	5.5
1	B	37	VAL	5.1
1	B	153	GLY	4.9
1	B	24	ASN	4.7
1	B	541	ALA	4.4
1	B	533	LYS	4.3
1	B	23	ILE	4.2
1	B	563	SER	4.0
1	B	529	ALA	4.0
1	B	509	ALA	3.9
1	B	469	LEU	3.9
1	B	502	HIS	3.7
1	B	528	TRP	3.7
1	B	39	ALA	3.6
1	B	22	PRO	3.5
1	B	28	ASN	3.5
1	B	512	LEU	3.4
1	B	530	VAL	3.4
1	B	27	SER	3.4
1	B	35	ASN	3.3

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Mol	Chain	Res	Type	RSRZ
1	B	534	LEU	3.2
1	B	380	ARG	3.2
1	B	106	LYS	3.1
1	A	544	GLN	3.1
1	B	15	ALA	3.0
1	B	544	GLN	3.0
1	B	309	GLN	3.0
1	B	110	ASN	2.9
1	A	545	LEU	2.9
1	B	524	TYR	2.8
1	B	34	HIS	2.8
1	B	523	ARG	2.7
1	B	212	LYS	2.7
1	B	30	LEU	2.6
1	B	19	SER	2.6
1	B	20	LYS	2.6
1	B	531	ARG	2.6
1	B	504	ALA	2.6
1	B	525	LEU	2.5
1	B	360	LEU	2.5
1	B	468	GLY	2.5
1	B	31	LEU	2.5
1	B	308	LEU	2.4
1	B	543	SER	2.4
1	B	514	GLN	2.4
1	B	370	VAL	2.4
1	B	520	THR	2.4
1	B	536	LEU	2.4
1	B	535	LYS	2.4
1	B	505	ARG	2.3
1	B	327	ALA	2.2
1	B	209	LYS	2.2
1	B	470	SER	2.2
1	B	430	PHE	2.2
1	B	208	TRP	2.2
1	B	477	TYR	2.2
1	B	492	LEU	2.2
1	B	434	LEU	2.1
1	B	373	ALA	2.1
1	B	464	GLU	2.1
1	A	563	SER	2.1
1	B	382	TYR	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	508	ARG	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](#)

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> 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	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	NN3	A	601	27/27	0.20	0.88	41,47,50,51	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.