



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:59 PM GMT

PDB ID : 4B76
Title : Discovery of an allosteric mechanism for the regulation of HCV NS3 protein function
Authors : Saalau-Bethell, S.M.; Woodhead, A.J.; Chessari, G.; Carr, M.G.; Coyle, J.; Graham, B.; Hiscock, S.D.; Murray, C.W.; Pathuri, P.; Rich, S.J.; Richardson, C.J.; Williams, P.A.; Jhoti, H.
Deposited on : 2012-08-16
Resolution : 2.14 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

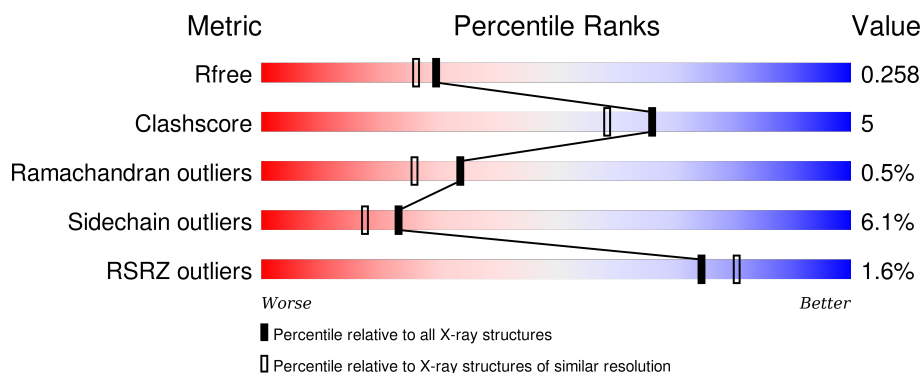
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION


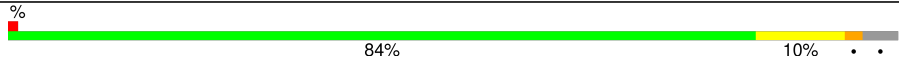
The reported resolution of this entry is 2.14 Å.

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}	91344	1693 (2.16-2.12)
Clashscore	102246	1824 (2.16-2.12)
Ramachandran outliers	100387	1798 (2.16-2.12)
Sidechain outliers	100360	1798 (2.16-2.12)
RSRZ outliers	91569	1699 (2.16-2.12)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	666	
1	B	666	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	PW1	A	1721	-	-	-	X
2	PW1	B	1721	-	-	-	X

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10215 atoms, of which 24 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 NON-STRUCTURAL PROTEIN 4A, SERINE PROTEASE NS3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	645	Total	C	N	O	S	0	0	0
			4808	3026	834	918	30			
1	B	642	Total	C	N	O	S	0	0	0
			4794	3019	831	914	30			

There are 56 discrepancies between the modelled and reference sequences:

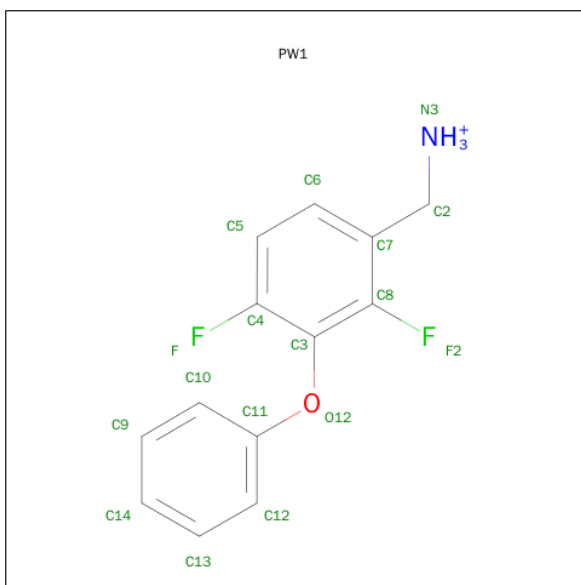
Chain	Residue	Modelled	Actual	Comment	Reference
A	684	MET	-	EXPRESSION TAG	UNP P26663
A	685	GLY	-	EXPRESSION TAG	UNP P26663
A	686	SER	-	EXPRESSION TAG	UNP P26663
A	687	SER	-	EXPRESSION TAG	UNP P26663
A	688	HIS	-	EXPRESSION TAG	UNP P26663
A	689	HIS	-	EXPRESSION TAG	UNP P26663
A	690	HIS	-	EXPRESSION TAG	UNP P26663
A	691	HIS	-	EXPRESSION TAG	UNP P26663
A	692	HIS	-	EXPRESSION TAG	UNP P26663
A	693	HIS	-	EXPRESSION TAG	UNP P26663
A	694	SER	-	EXPRESSION TAG	UNP P26663
A	695	SER	-	EXPRESSION TAG	UNP P26663
A	696	GLY	-	EXPRESSION TAG	UNP P26663
A	697	LEU	-	EXPRESSION TAG	UNP P26663
A	698	VAL	-	EXPRESSION TAG	UNP P26663
A	699	PRO	-	EXPRESSION TAG	UNP P26663
A	700	ARG	-	EXPRESSION TAG	UNP P26663
A	701	GLY	-	EXPRESSION TAG	UNP P26663
A	702	SER	-	EXPRESSION TAG	UNP P26663
A	703	HIS	-	EXPRESSION TAG	UNP P26663
A	704	MET	-	EXPRESSION TAG	UNP P26663
A	66	GLY	ALA	CONFLICT	UNP P26663
A	86	GLN	PRO	CONFLICT	UNP P26663
A	87	ALA	LYS	CONFLICT	UNP P26663

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Chain	Residue	Modelled	Actual	Comment	Reference
A	147	SER	PHE	CONFLICT	UNP P26663
A	718	SER	-	LINKER	UNP P26663
A	719	GLY	-	LINKER	UNP P26663
A	720	SER	-	LINKER	UNP P26663
B	684	MET	-	EXPRESSION TAG	UNP P26663
B	685	GLY	-	EXPRESSION TAG	UNP P26663
B	686	SER	-	EXPRESSION TAG	UNP P26663
B	687	SER	-	EXPRESSION TAG	UNP P26663
B	688	HIS	-	EXPRESSION TAG	UNP P26663
B	689	HIS	-	EXPRESSION TAG	UNP P26663
B	690	HIS	-	EXPRESSION TAG	UNP P26663
B	691	HIS	-	EXPRESSION TAG	UNP P26663
B	692	HIS	-	EXPRESSION TAG	UNP P26663
B	693	HIS	-	EXPRESSION TAG	UNP P26663
B	694	SER	-	EXPRESSION TAG	UNP P26663
B	695	SER	-	EXPRESSION TAG	UNP P26663
B	696	GLY	-	EXPRESSION TAG	UNP P26663
B	697	LEU	-	EXPRESSION TAG	UNP P26663
B	698	VAL	-	EXPRESSION TAG	UNP P26663
B	699	PRO	-	EXPRESSION TAG	UNP P26663
B	700	ARG	-	EXPRESSION TAG	UNP P26663
B	701	GLY	-	EXPRESSION TAG	UNP P26663
B	702	SER	-	EXPRESSION TAG	UNP P26663
B	703	HIS	-	EXPRESSION TAG	UNP P26663
B	704	MET	-	EXPRESSION TAG	UNP P26663
B	718	SER	-	LINKER	UNP P26663
B	719	GLY	-	LINKER	UNP P26663
B	720	SER	-	LINKER	UNP P26663
B	66	GLY	ALA	CONFLICT	UNP P26663
B	86	GLN	PRO	CONFLICT	UNP P26663
B	87	ALA	LYS	CONFLICT	UNP P26663
B	147	SER	PHE	CONFLICT	UNP P26663

- Molecule 2 is [2,4-BIS(FLUORANYL)-3-PHENOXY-PHENYL]METHYLAZANIUM (three-letter code: PW1) (formula: C₁₃H₁₂F₂NO).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	F	H	N	O	0	0
			29	13	2	12	1	1		
2	B	1	Total	C	F	H	N	O	0	0
			29	13	2	12	1	1		

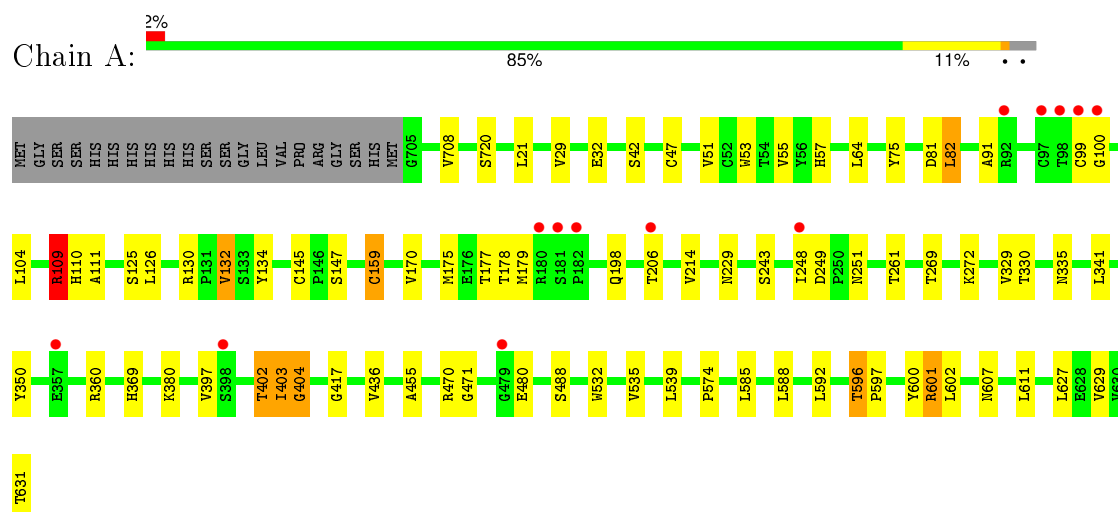
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	308	Total	O	0	0
			308	308		
3	B	247	Total	O	0	0
			247	247		

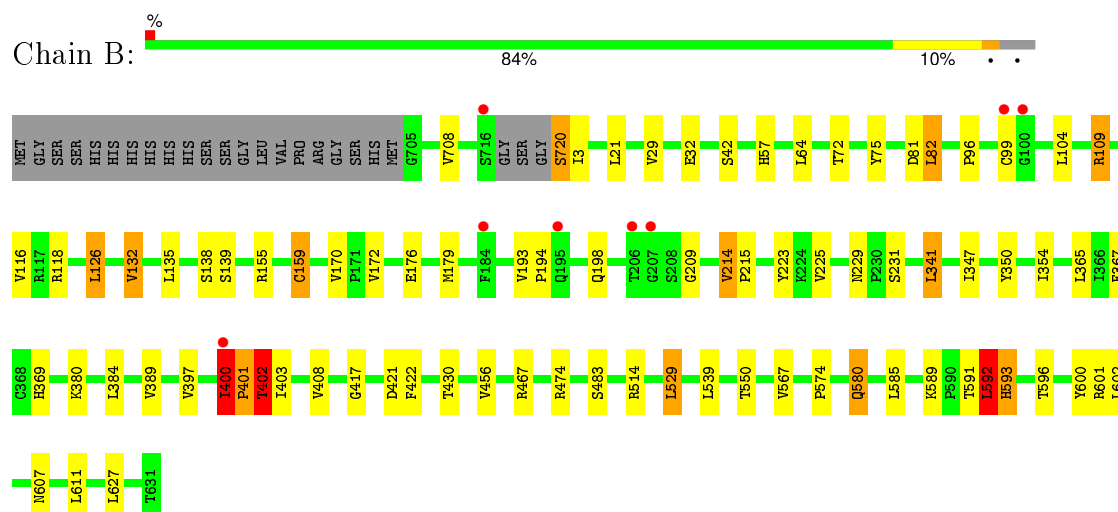
3 Residue-property plots [i](#)

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: NON-STRUCTURAL PROTEIN 4A, SERINE PROTEASE NS3



- Molecule 1: NON-STRUCTURAL PROTEIN 4A, SERINE PROTEASE NS3



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	91.35Å 110.31Å 142.15Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	87.15 – 2.14 87.15 – 2.14	Depositor EDS
% Data completeness (in resolution range)	99.2 (87.15-2.14) 99.2 (87.15-2.14)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.07 (at 2.14Å)	Xtriage
Refinement program	REFMAC 5.7.0025	Depositor
R, R_{free}	0.194 , 0.253 0.200 , 0.258	Depositor DCC
R_{free} test set	4034 reflections (5.38%)	DCC
Wilson B-factor (Å ²)	29.0	Xtriage
Anisotropy	0.063	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 40.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 79036 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	10215	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.15% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: PW1

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	0.62	2/4917 (0.0%)	0.74	2/6714 (0.0%)
1	B	0.58	1/4901 (0.0%)	0.74	5/6690 (0.1%)
All	All	0.60	3/9818 (0.0%)	0.74	7/13404 (0.1%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	720	SER	C-N	13.75	1.65	1.34
1	A	159	CYS	CB-SG	-6.72	1.70	1.82
1	B	159	CYS	CB-SG	-6.22	1.71	1.82

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	400	ILE	C-N-CD	-8.70	101.46	120.60
1	B	601	ARG	NE-CZ-NH2	-6.44	117.08	120.30
1	B	400	ILE	C-N-CA	6.42	148.96	122.00
1	A	109	ARG	NE-CZ-NH1	5.84	123.22	120.30
1	B	401	PRO	N-CA-C	5.79	127.15	112.10
1	B	402	THR	N-CA-CB	5.15	120.09	110.30
1	A	601	ARG	NE-CZ-NH2	-5.05	117.77	120.30

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	4808	0	4784	42	0
1	B	4794	0	4771	52	0
2	A	17	12	12	0	0
2	B	17	12	12	0	0
3	A	308	0	0	3	0
3	B	247	0	0	5	0
All	All	10191	24	9579	94	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 5.

All (94) 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:B:400:ILE:HB	1:B:401:PRO:CA	2.09	0.82
1:A:596:THR:HG22	1:A:607:ASN:HD22	1.44	0.81
1:A:132:VAL:HG11	1:A:159:CYS:SG	2.22	0.80
1:B:132:VAL:CG1	1:B:159:CYS:SG	2.70	0.79
1:A:132:VAL:CG1	1:A:159:CYS:SG	2.72	0.78
1:B:132:VAL:HG11	1:B:159:CYS:SG	2.24	0.77
1:A:269:THR:HG23	1:A:272:LYS:H	1.49	0.77
1:B:400:ILE:HB	1:B:401:PRO:HA	1.67	0.75
1:A:75:TYR:CD1	1:A:178:THR:HG22	2.23	0.74
1:B:32:GLU:OE2	3:B:2018:HOH:O	2.13	0.67
1:B:350:TYR:OH	1:B:369:HIS:HD2	1.77	0.67
1:B:550:THR:O	3:B:2224:HOH:O	2.13	0.66
1:A:269:THR:OG1	3:A:2123:HOH:O	2.13	0.66
1:B:596:THR:CG2	1:B:607:ASN:HD22	2.11	0.64
1:B:57:HIS:HD2	1:B:81:ASP:OD2	1.81	0.64
1:A:57:HIS:HD2	1:A:81:ASP:OD2	1.82	0.63
1:A:380:LYS:NZ	3:A:2184:HOH:O	2.32	0.63
1:A:132:VAL:HG13	1:A:159:CYS:SG	2.38	0.62
1:A:330:THR:CB	1:A:480:GLU:OE2	2.48	0.62
1:B:574:PRO:O	1:B:596:THR:HG22	1.99	0.61
1:B:397:VAL:HG11	1:B:417:GLY:O	2.01	0.61
1:B:596:THR:HG23	1:B:607:ASN:HD22	1.66	0.61
1:A:403:ILE:O	1:A:404:GLY:O	2.19	0.61
1:B:514:ARG:HG2	1:B:529:LEU:HD23	1.86	0.58
1:B:132:VAL:HG13	1:B:159:CYS:SG	2.45	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:455:ALA:HA	1:A:480:GLU:HG2	1.87	0.57
1:A:350:TYR:OH	1:A:369:HIS:HD2	1.88	0.57
1:A:330:THR:HB	1:A:480:GLU:OE2	2.06	0.56
1:B:389:VAL:HG21	1:B:400:ILE:O	2.06	0.55
1:B:397:VAL:O	1:B:400:ILE:HG12	2.07	0.55
1:B:720:SER:O	1:B:3:ILE:N	2.40	0.54
1:A:397:VAL:HG11	1:A:417:GLY:O	2.08	0.54
1:B:82:LEU:HG	1:B:170:VAL:HG11	1.90	0.54
1:A:532:TRP:O	1:A:535:VAL:HG22	2.07	0.54
1:B:580:GLN:HE21	1:B:580:GLN:HA	1.73	0.53
1:B:380:LYS:O	1:B:384:LEU:HD13	2.07	0.53
1:B:229:ASN:HD22	1:B:231:SER:H	1.56	0.52
1:B:400:ILE:HG21	1:B:402:THR:HG23	1.92	0.52
1:B:400:ILE:HG13	3:B:2158:HOH:O	2.09	0.52
1:B:408:VAL:HG11	1:B:422:PHE:CZ	2.45	0.52
1:A:403:ILE:HG22	1:A:404:GLY:N	2.25	0.51
1:B:592:LEU:HD23	1:B:593:HIS:CG	2.45	0.51
1:B:401:PRO:O	3:B:2159:HOH:O	2.19	0.51
1:B:214:VAL:HG22	1:B:215:PRO:HD3	1.91	0.51
1:B:589:LYS:O	1:B:592:LEU:HD13	2.11	0.51
1:A:53:TRP:CD1	1:A:175:MET:CE	2.94	0.51
1:B:421:ASP:HB3	1:B:467:ARG:HD2	1.93	0.50
1:A:402:THR:HG23	1:A:403:ILE:HB	1.94	0.50
1:A:601:ARG:NH2	3:A:2290:HOH:O	2.45	0.50
1:B:592:LEU:HG	1:B:593:HIS:H	1.77	0.49
1:B:96:PRO:HB3	1:B:172:VAL:HG21	1.94	0.49
1:A:42:SER:OG	1:A:109:ARG:NH2	2.38	0.49
1:B:400:ILE:HG21	1:B:402:THR:CG2	2.43	0.49
1:A:130:ARG:HD3	1:A:134:TYR:CD2	2.48	0.49
1:A:596:THR:CG2	1:A:607:ASN:HD22	2.20	0.48
1:B:42:SER:OG	1:B:109:ARG:NH2	2.41	0.48
1:A:53:TRP:NE1	1:A:175:MET:HE2	2.29	0.48
1:A:53:TRP:CD1	1:A:175:MET:HE2	2.49	0.48
1:B:116:VAL:HG22	1:B:126:LEU:HD12	1.97	0.47
1:A:330:THR:OG1	1:A:480:GLU:OE2	2.27	0.47
1:A:335:ASN:O	1:A:471:GLY:N	2.43	0.47
1:A:29:VAL:CG2	1:A:91:ALA:HB2	2.45	0.47
1:A:53:TRP:CE2	1:A:175:MET:HE2	2.50	0.46
1:A:99:CYS:SG	1:A:100:GLY:N	2.89	0.46
1:A:574:PRO:O	1:A:596:THR:HB	2.16	0.45
1:A:145:CYS:SG	1:A:147:SER:OG	2.71	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:110:HIS:O	1:A:111:ALA:HB3	2.17	0.45
1:A:82:LEU:HG	1:A:170:VAL:HG11	1.98	0.45
1:B:194:PRO:HG3	1:B:198:GLN:HB2	1.99	0.44
1:B:400:ILE:CG2	1:B:402:THR:N	2.80	0.44
1:A:229:ASN:O	1:A:269:THR:HA	2.18	0.44
1:A:251:ASN:HD22	1:A:261:THR:H	1.65	0.44
1:B:456:VAL:HG23	1:B:483:SER:HB2	1.99	0.44
1:B:229:ASN:ND2	1:B:231:SER:H	2.16	0.43
1:B:365:LEU:HD21	1:B:367:PHE:CZ	2.53	0.43
1:A:436:VAL:CG1	1:A:488:SER:HB3	2.48	0.43
1:B:135:LEU:O	1:B:138:SER:HB2	2.19	0.43
1:B:193:VAL:HG21	1:B:223:TYR:CE2	2.55	0.42
1:A:243:SER:HA	1:A:248:ILE:O	2.20	0.42
1:B:401:PRO:O	1:B:402:THR:CB	2.68	0.42
1:A:596:THR:HA	1:A:597:PRO:HD3	1.93	0.42
1:A:47:CYS:HA	1:A:51:VAL:O	2.20	0.42
1:B:341:LEU:HD13	1:B:474:ARG:HB3	2.03	0.41
1:B:209:GLY:O	1:B:214:VAL:HG13	2.21	0.41
1:B:720:SER:C	1:B:3:ILE:N	2.74	0.41
1:B:214:VAL:HG22	1:B:215:PRO:CD	2.51	0.41
1:B:430:THR:HG21	3:B:2172:HOH:O	2.20	0.41
1:B:347:ILE:HB	1:B:354:ILE:HB	2.03	0.41
1:B:400:ILE:HB	1:B:401:PRO:C	2.40	0.41
1:B:75:TYR:CE1	1:B:179:MET:HA	2.56	0.41
1:A:57:HIS:HE1	1:A:631:THR:O	2.04	0.40
1:A:29:VAL:HG23	1:A:91:ALA:HB2	2.03	0.40
1:B:592:LEU:CG	1:B:593:HIS:H	2.34	0.40
1:B:400:ILE:HG22	1:B:402:THR:N	2.37	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	643/666 (96%)	623 (97%)	18 (3%)	2 (0%)	46	41
1	B	637/666 (96%)	619 (97%)	14 (2%)	4 (1%)	30	21
All	All	1280/1332 (96%)	1242 (97%)	32 (2%)	6 (0%)	34	26

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	403	ILE
1	A	404	GLY
1	B	400	ILE
1	B	592	LEU
1	B	593	HIS
1	B	402	THR

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	526/544 (97%)	494 (94%)	32 (6%)	23	17
1	B	525/544 (96%)	493 (94%)	32 (6%)	23	17
All	All	1051/1088 (97%)	987 (94%)	64 (6%)	23	17

All (64) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	708	VAL
1	A	21	LEU
1	A	32	GLU
1	A	55	VAL
1	A	64	LEU
1	A	82	LEU
1	A	104	LEU
1	A	109	ARG
1	A	125	SER
1	A	126	LEU

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Mol	Chain	Res	Type
1	A	132	VAL
1	A	177	THR
1	A	179	MET
1	A	198	GLN
1	A	206	THR
1	A	214	VAL
1	A	249	ASP
1	A	329	VAL
1	A	341	LEU
1	A	360	ARG
1	A	402	THR
1	A	470	ARG
1	A	539	LEU
1	A	585	LEU
1	A	588	LEU
1	A	592	LEU
1	A	596	THR
1	A	600	TYR
1	A	602	LEU
1	A	611	LEU
1	A	627	LEU
1	A	629	VAL
1	B	708	VAL
1	B	720	SER
1	B	21	LEU
1	B	29	VAL
1	B	64	LEU
1	B	72	THR
1	B	82	LEU
1	B	99	CYS
1	B	104	LEU
1	B	109	ARG
1	B	118	ARG
1	B	126	LEU
1	B	132	VAL
1	B	139	SER
1	B	155	ARG
1	B	176	GLU
1	B	214	VAL
1	B	225	VAL
1	B	341	LEU
1	B	400	ILE

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Mol	Chain	Res	Type
1	B	403	ILE
1	B	529	LEU
1	B	539	LEU
1	B	567	VAL
1	B	580	GLN
1	B	585	LEU
1	B	591	THR
1	B	592	LEU
1	B	600	TYR
1	B	602	LEU
1	B	611	LEU
1	B	627	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	57	HIS
1	A	110	HIS
1	A	229	ASN
1	A	251	ASN
1	A	293	HIS
1	A	369	HIS
1	A	434	GLN
1	A	526	GLN
1	A	556	ASN
1	A	606	GLN
1	A	607	ASN
1	B	27	ASN
1	B	57	HIS
1	B	229	ASN
1	B	251	ASN
1	B	293	HIS
1	B	369	HIS
1	B	526	GLN
1	B	580	GLN
1	B	607	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

2 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	PW1	A	1721	-	18,18,18	0.42	0	22,24,24	1.16	1 (4%)
2	PW1	B	1721	-	18,18,18	0.44	0	22,24,24	1.12	3 (13%)

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	PW1	A	1721	-	-	0/6/6/6	0/2/2/2
2	PW1	B	1721	-	-	0/6/6/6	0/2/2/2

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1721	PW1	O12-C11-C12	2.04	126.74	119.42
2	B	1721	PW1	F2-C8-C7	2.29	119.98	117.75
2	B	1721	PW1	C6-C7-C8	3.33	118.40	116.42
2	A	1721	PW1	C6-C7-C8	3.81	118.68	116.42

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

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	645/666 (96%)	-0.19	13 (2%) 68 75	14, 32, 68, 95	0
1	B	642/666 (96%)	-0.29	8 (1%) 81 85	17, 32, 58, 103	0
All	All	1287/1332 (96%)	-0.24	21 (1%) 74 80	14, 32, 65, 103	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	98	THR	4.4
1	A	99	CYS	4.4
1	B	100	GLY	3.7
1	A	248	ILE	3.5
1	B	400	ILE	3.4
1	A	100	GLY	3.3
1	B	206	THR	3.2
1	A	182	PRO	3.1
1	A	357	GLU	2.7
1	A	181	SER	2.6
1	B	195	GLN	2.5
1	A	479	GLY	2.4
1	B	184	PHE	2.4
1	A	206	THR	2.3
1	A	180	ARG	2.3
1	B	99	CYS	2.2
1	A	398	SER	2.2
1	A	97	CYS	2.1
1	B	207	GLY	2.1
1	A	92	ARG	2.1
1	B	716	SER	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, 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	LLDF	B-factors(\AA^2)	Q<0.9
2	PW1	A	1721	17/17	0.96	0.15	4.13	18,23,32,47	29
2	PW1	B	1721	17/17	0.95	0.13	2.32	13,20,30,36	29

6.5 Other polymers [i](#)

There are no such residues in this entry.