



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 19, 2016 – 11:01 PM GMT

PDB ID : 5CYR
Title : Crystal structure of Thosea asigna virus RNA-dependent RNA polymerase (RdRP) complexed with ATP and ssRNA
Authors : Ferrero, D.S.; Buxaderas, M.; Rodriguez, J.F.; Verdaguer, N.
Deposited on : 2015-07-30
Resolution : 3.50 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 : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026982
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) : rb-20026982

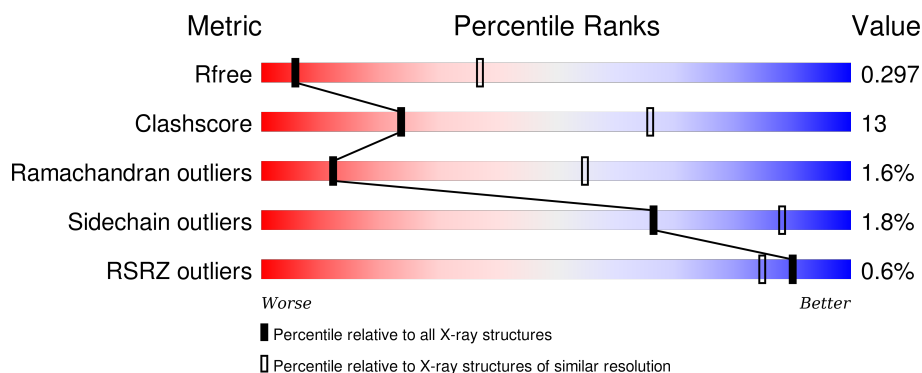
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.50 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	1051 (3.60-3.40)
Clashscore	102246	1157 (3.60-3.40)
Ramachandran outliers	100387	1120 (3.60-3.40)
Sidechain outliers	100360	1121 (3.60-3.40)
RSRZ outliers	91569	1058 (3.60-3.40)

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	705	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 24%, green 68%, grey 6%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> % 68% 24% •• 6% </div> </div>
1	B	705	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, green 64%, yellow 28%, orange 6%, grey 2%);"></div> <div style="display: flex; justify-content: space-between; width: 100%;"> 64% 28% • 6% </div> </div>

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	ATP	B	701	-	-	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 10528 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 RNA-dependent RNA polymerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	664	Total	C	N	O	S	43	0	0
			5243	3328	907	982	26			
1	A	662	Total	C	N	O	S	2	0	0
			5223	3314	903	980	26			

There are 62 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-30	MET	-	initiating methionine	UNP Q6A562
B	-29	SER	-	expression tag	UNP Q6A562
B	-28	TYR	-	expression tag	UNP Q6A562
B	-27	TYR	-	expression tag	UNP Q6A562
B	-26	HIS	-	expression tag	UNP Q6A562
B	-25	HIS	-	expression tag	UNP Q6A562
B	-24	HIS	-	expression tag	UNP Q6A562
B	-23	HIS	-	expression tag	UNP Q6A562
B	-22	HIS	-	expression tag	UNP Q6A562
B	-21	HIS	-	expression tag	UNP Q6A562
B	-20	ASP	-	expression tag	UNP Q6A562
B	-19	TYR	-	expression tag	UNP Q6A562
B	-18	ASP	-	expression tag	UNP Q6A562
B	-17	ILE	-	expression tag	UNP Q6A562
B	-16	PRO	-	expression tag	UNP Q6A562
B	-15	THR	-	expression tag	UNP Q6A562
B	-14	THR	-	expression tag	UNP Q6A562
B	-13	GLU	-	expression tag	UNP Q6A562
B	-12	ASN	-	expression tag	UNP Q6A562
B	-11	LEU	-	expression tag	UNP Q6A562
B	-10	TYR	-	expression tag	UNP Q6A562
B	-9	PHE	-	expression tag	UNP Q6A562
B	-8	GLN	-	expression tag	UNP Q6A562
B	-7	GLY	-	expression tag	UNP Q6A562
B	-6	ALA	-	expression tag	UNP Q6A562

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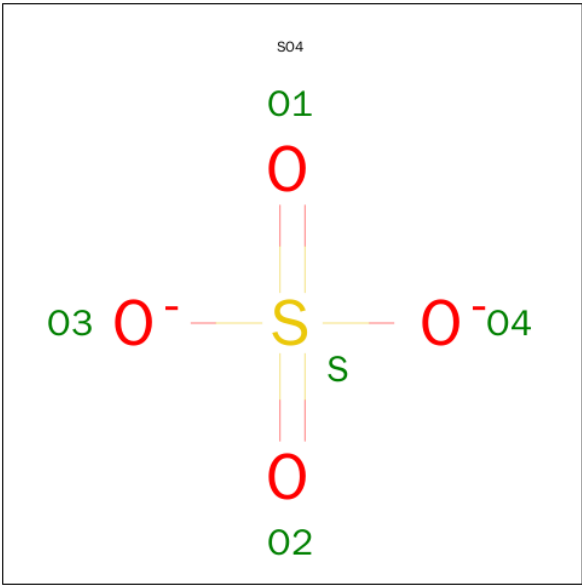
Chain	Residue	Modelled	Actual	Comment	Reference
B	-5	MET	-	expression tag	UNP Q6A562
B	-4	GLY	-	expression tag	UNP Q6A562
B	-3	ALA	-	expression tag	UNP Q6A562
B	-2	MET	-	expression tag	UNP Q6A562
B	-1	GLY	-	expression tag	UNP Q6A562
B	0	ILE	-	expression tag	UNP Q6A562
A	-30	MET	-	initiating methionine	UNP Q6A562
A	-29	SER	-	expression tag	UNP Q6A562
A	-28	TYR	-	expression tag	UNP Q6A562
A	-27	TYR	-	expression tag	UNP Q6A562
A	-26	HIS	-	expression tag	UNP Q6A562
A	-25	HIS	-	expression tag	UNP Q6A562
A	-24	HIS	-	expression tag	UNP Q6A562
A	-23	HIS	-	expression tag	UNP Q6A562
A	-22	HIS	-	expression tag	UNP Q6A562
A	-21	HIS	-	expression tag	UNP Q6A562
A	-20	ASP	-	expression tag	UNP Q6A562
A	-19	TYR	-	expression tag	UNP Q6A562
A	-18	ASP	-	expression tag	UNP Q6A562
A	-17	ILE	-	expression tag	UNP Q6A562
A	-16	PRO	-	expression tag	UNP Q6A562
A	-15	THR	-	expression tag	UNP Q6A562
A	-14	THR	-	expression tag	UNP Q6A562
A	-13	GLU	-	expression tag	UNP Q6A562
A	-12	ASN	-	expression tag	UNP Q6A562
A	-11	LEU	-	expression tag	UNP Q6A562
A	-10	TYR	-	expression tag	UNP Q6A562
A	-9	PHE	-	expression tag	UNP Q6A562
A	-8	GLN	-	expression tag	UNP Q6A562
A	-7	GLY	-	expression tag	UNP Q6A562
A	-6	ALA	-	expression tag	UNP Q6A562
A	-5	MET	-	expression tag	UNP Q6A562
A	-4	GLY	-	expression tag	UNP Q6A562
A	-3	ALA	-	expression tag	UNP Q6A562
A	-2	MET	-	expression tag	UNP Q6A562
A	-1	GLY	-	expression tag	UNP Q6A562
A	0	ILE	-	expression tag	UNP Q6A562

- Molecule 2 is ADENOSINE-5'-TRIPHOSPHATE (three-letter code: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$).



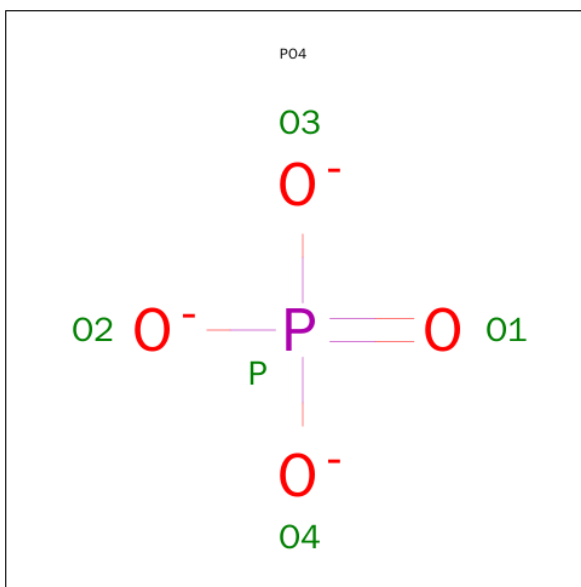
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	B	1	Total	C	N	O	P	0	0
			31	10	5	13	3		

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	O	P	0	0
			5	4	1		
4	A	1	Total	O	P	0	0
			5	4	1		
4	A	1	Total	O	P	0	0
			5	4	1		
4	A	1	Total	O	P	0	0
			5	4	1		

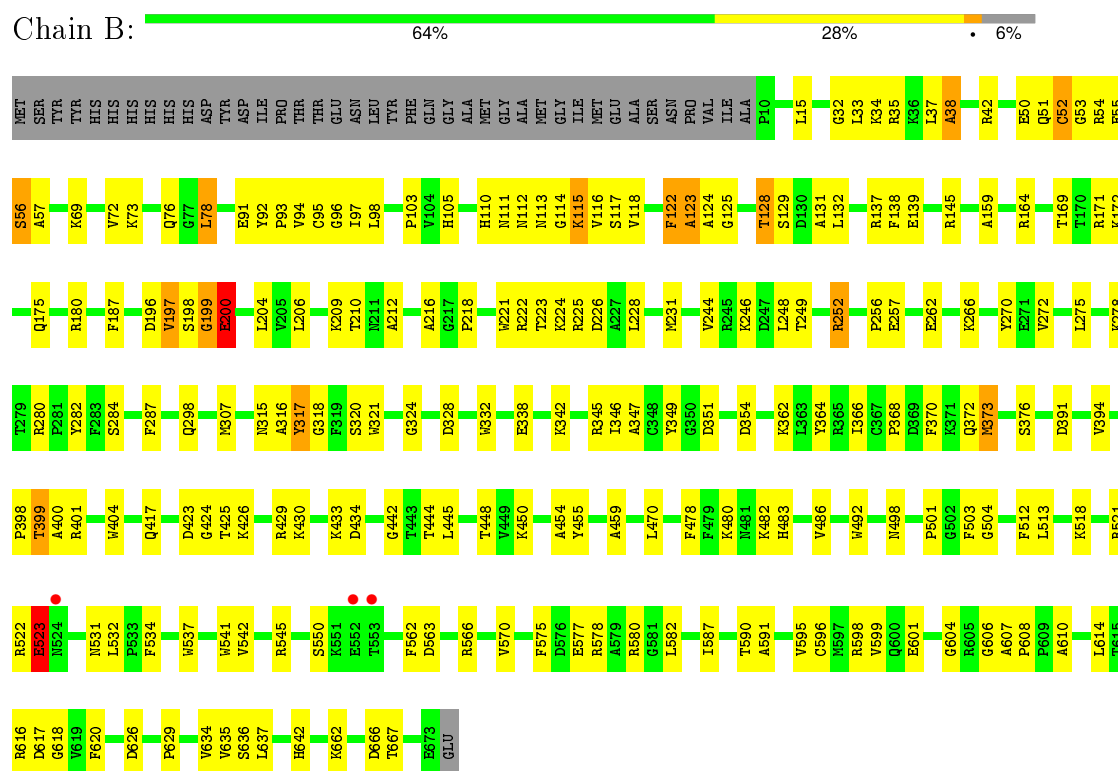
- Molecule 5 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Mg	0	0
			1	1		

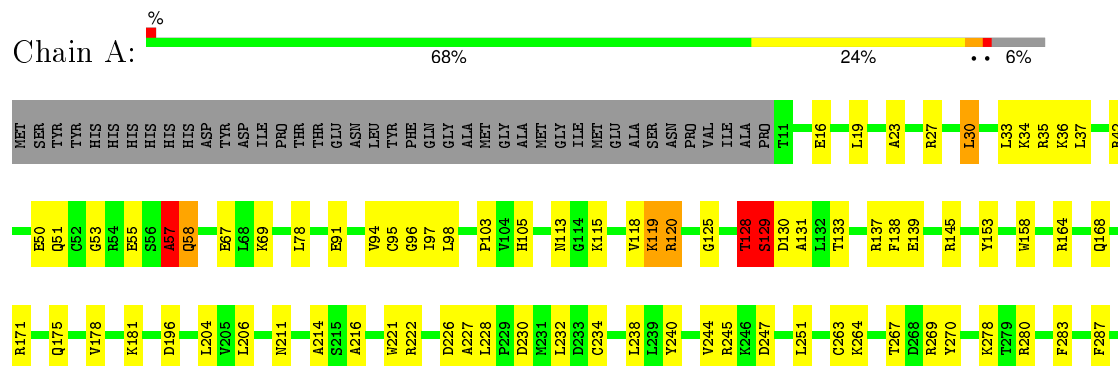
3 Residue-property plots

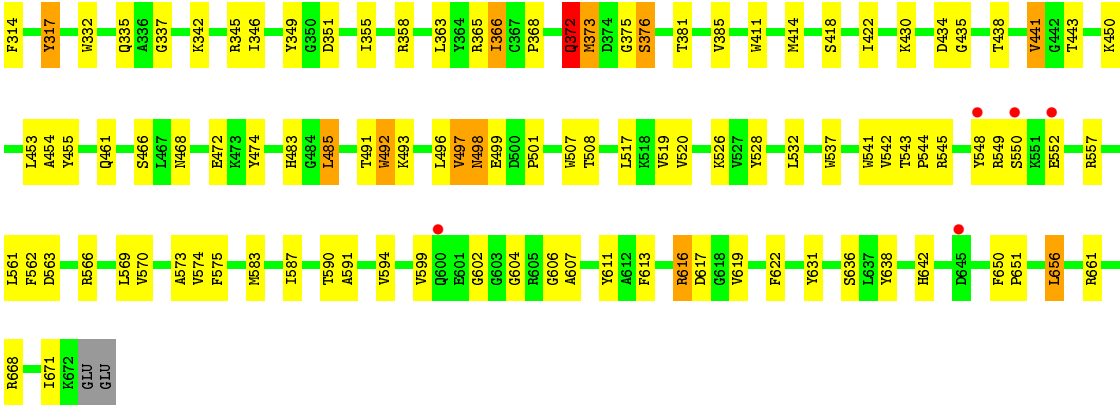
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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: RNA-dependent RNA polymerase



• Molecule 1: RNA-dependent RNA polymerase





4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	142.51Å 158.34Å 217.10Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.60 – 3.50 48.35 – 3.50	Depositor EDS
% Data completeness (in resolution range)	89.5 (47.60-3.50) 89.6 (48.35-3.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.15	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.47 (at 3.48Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.4_1496)	Depositor
R, R_{free}	0.252 , 0.296 0.257 , 0.297	Depositor DCC
R_{free} test set	1410 reflections (5.05%)	DCC
Wilson B-factor (Å ²)	25.0	Xtriage
Anisotropy	0.306	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 23.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.27$	Xtriage
Outliers	4 of 27974 reflections (0.014%)	Xtriage
F_o, F_c correlation	0.76	EDS
Total number of atoms	10528	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 23.50 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 4.6050e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MG, PO4, ATP, SO4

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	0.59	5/5341 (0.1%)	0.61	13/7238 (0.2%)
1	B	0.38	2/5363 (0.0%)	0.57	8/7267 (0.1%)
All	All	0.50	7/10704 (0.1%)	0.59	21/14505 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2
1	B	0	3
All	All	0	5

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	57	ALA	C-N	23.24	1.87	1.34
1	A	128	THR	C-N	-23.10	0.81	1.34
1	A	497	VAL	C-N	15.84	1.70	1.34
1	B	128	THR	C-N	-15.60	0.98	1.34
1	B	122	PHE	C-N	12.91	1.63	1.34
1	A	496	LEU	C-N	-12.77	1.04	1.34
1	A	129	SER	C-N	-9.96	1.11	1.34

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	523	GLU	CB-CA-C	16.99	144.38	110.40
1	A	372	GLN	CB-CA-C	-16.00	78.40	110.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	129	SER	O-C-N	-14.63	99.28	122.70
1	B	200	GLU	N-CA-CB	-10.53	91.65	110.60
1	A	373	MET	N-CA-CB	-10.52	91.66	110.60
1	B	373	MET	N-CA-CB	-10.43	91.83	110.60
1	A	57	ALA	O-C-N	-9.50	107.50	122.70
1	B	199	GLY	N-CA-C	-9.31	89.83	113.10
1	A	498	ASN	N-CA-C	-9.26	86.00	111.00
1	A	128	THR	O-C-N	9.22	137.45	122.70
1	A	128	THR	CA-C-N	-9.18	97.00	117.20
1	B	523	GLU	N-CA-C	-8.84	87.14	111.00
1	A	128	THR	C-N-CA	-7.81	102.17	121.70
1	A	129	SER	CA-C-N	7.49	133.69	117.20
1	A	57	ALA	C-N-CA	6.95	139.08	121.70
1	A	129	SER	C-N-CA	6.58	138.16	121.70
1	B	246	LYS	CB-CA-C	-6.51	97.37	110.40
1	A	498	ASN	CB-CA-C	-6.11	98.17	110.40
1	A	267	THR	N-CA-C	-5.88	95.12	111.00
1	B	122	PHE	O-C-N	-5.61	113.73	122.70
1	B	199	GLY	C-N-CA	5.26	134.84	121.70

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	128	THR	Mainchain
1	A	129	SER	Mainchain
1	B	128	THR	Mainchain
1	B	199	GLY	Peptide
1	B	523	GLU	Peptide

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	5223	0	5146	132	2
1	B	5243	0	5181	151	0
2	B	31	0	12	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	5	0	0	0	0
3	B	5	0	0	0	0
4	A	15	0	0	0	0
4	B	5	0	0	1	0
5	A	1	0	0	0	0
All	All	10528	0	10339	275	2

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 13.

All (275) 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:A:128:THR:C	1:A:129:SER:CA	1.81	1.50
1:A:497:VAL:C	1:A:498:ASN:N	1.70	1.45
1:A:128:THR:CA	1:A:129:SER:N	1.81	1.43
1:A:57:ALA:C	1:A:58:GLN:N	1.87	1.28
1:B:50:GLU:OE2	1:B:52:CYS:O	1.66	1.12
1:A:50:GLU:C	1:A:51:GLN:N	2.12	1.03
1:A:128:THR:O	1:A:129:SER:N	1.90	1.01
1:A:120:ARG:HG2	1:A:158:TRP:CZ2	1.96	1.00
1:B:398:PRO:O	1:B:400:ALA:N	1.95	1.00
1:B:164:ARG:HG2	1:B:372:GLN:NE2	1.77	0.98
1:B:164:ARG:HG2	1:B:372:GLN:HE22	1.25	0.96
1:B:50:GLU:HG3	1:B:137:ARG:O	1.68	0.94
1:A:372:GLN:HG3	1:A:372:GLN:O	1.68	0.92
1:A:128:THR:O	1:A:129:SER:CA	2.17	0.91
1:A:128:THR:O	1:A:129:SER:HA	1.72	0.89
1:B:398:PRO:C	1:B:400:ALA:H	1.78	0.87
1:A:128:THR:C	1:A:129:SER:N	0.80	0.86
1:A:128:THR:C	1:A:129:SER:HA	1.97	0.85
1:B:33:LEU:O	1:B:37:LEU:HB2	1.76	0.85
1:B:53:GLY:HA2	1:B:137:ARG:HG3	1.59	0.84
1:B:399:THR:O	1:A:668:ARG:HG3	1.78	0.83
1:B:282:TYR:OH	4:B:703:PO4:O3	1.97	0.83
1:A:372:GLN:CG	1:A:372:GLN:O	2.07	0.79
1:B:124:ALA:N	1:B:125:GLY:HA3	1.97	0.79
1:B:93:PRO:HA	1:B:115:LYS:HB3	1.65	0.78
1:A:355:ILE:HB	1:A:366:ILE:HG13	1.65	0.78
1:A:332:TRP:HA	1:A:335:GLN:HE21	1.50	0.75
1:B:54:ARG:O	1:B:56:SER:N	2.17	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:566:ARG:HH22	1:A:606:GLY:HA3	1.52	0.73
1:A:128:THR:CB	1:A:129:SER:N	2.52	0.72
1:B:34:LYS:O	1:B:38:ALA:HB2	1.90	0.71
1:B:601:GLU:O	1:B:604:GLY:N	2.22	0.71
1:B:198:SER:HB2	1:B:399:THR:HG23	1.72	0.71
1:B:175:GLN:NE2	1:B:434:ASP:OD1	2.24	0.70
1:A:443:THR:O	1:A:443:THR:HG22	1.89	0.69
1:A:120:ARG:NH1	1:A:422:ILE:O	2.26	0.69
1:A:351:ASP:HB3	1:A:373:MET:HG2	1.75	0.68
1:B:91:GLU:OE1	1:B:426:LYS:HA	1.94	0.68
1:B:278:LYS:NZ	2:B:701:ATP:O1G	2.21	0.68
1:A:206:LEU:HD11	1:A:232:LEU:HD21	1.75	0.67
1:A:120:ARG:HG2	1:A:158:TRP:HZ2	1.52	0.66
1:A:120:ARG:NH2	1:A:264:LYS:HA	2.10	0.66
1:B:231:MET:HG3	1:B:287:PHE:HD2	1.61	0.65
1:B:394:VAL:O	1:B:398:PRO:HB3	1.96	0.65
1:B:480:LYS:HE2	1:B:486:VAL:HG13	1.79	0.65
1:A:363:LEU:HD23	1:A:499:GLU:HG2	1.79	0.65
1:B:197:VAL:HA	1:B:204:LEU:HD21	1.79	0.65
1:A:358:ARG:NH2	1:A:499:GLU:OE2	2.24	0.65
1:A:119:LYS:O	1:A:120:ARG:HD2	1.98	0.64
1:B:42:ARG:NH2	1:B:587:ILE:O	2.31	0.64
1:B:616:ARG:HA	1:B:618:GLY:N	2.13	0.63
1:A:94:VAL:HG13	1:A:105:HIS:HD2	1.64	0.62
1:B:391:ASP:HA	1:B:394:VAL:HG22	1.79	0.62
1:A:450:LYS:HE2	1:A:485:LEU:HD13	1.80	0.62
1:A:454:ALA:HA	1:A:483:HIS:CD2	2.35	0.61
1:B:206:LEU:HB3	1:A:16:GLU:HB3	1.82	0.61
1:B:94:VAL:HG23	1:B:115:LYS:HA	1.82	0.61
1:B:417:GLN:HB2	1:B:433:LYS:HE2	1.82	0.61
1:B:52:CYS:HB3	1:B:57:ALA:HB2	1.83	0.60
1:B:69:LYS:NZ	1:B:73:LYS:HB2	2.17	0.60
1:A:507:TRP:HB2	1:A:517:LEU:HD22	1.84	0.60
1:B:169:THR:HG22	1:B:430:LYS:HE2	1.83	0.60
1:B:249:THR:HA	1:B:252:ARG:HG2	1.84	0.60
1:B:596:CYS:HA	1:B:626:ASP:HB3	1.84	0.59
1:A:461:GLN:O	1:A:466:SER:OG	2.21	0.59
1:B:111:ASN:HB3	1:B:116:VAL:HA	1.84	0.59
1:A:491:THR:O	1:A:493:LYS:N	2.36	0.58
1:A:131:ALA:O	1:A:541:TRP:NE1	2.37	0.58
1:B:563:ASP:HB3	1:B:607:ALA:HB3	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:175:GLN:NE2	1:A:434:ASP:OD1	2.36	0.58
1:B:398:PRO:C	1:B:400:ALA:N	2.47	0.58
1:B:76:GLN:NE2	1:B:122:PHE:O	2.34	0.58
1:A:636:SER:OG	1:A:642:HIS:O	2.21	0.58
1:A:91:GLU:HB3	1:A:115:LYS:HG2	1.84	0.57
1:B:222:ARG:NH1	1:B:226:ASP:O	2.37	0.57
1:B:351:ASP:HB3	1:B:373:MET:HG2	1.86	0.57
1:A:472:GLU:HB2	1:A:492:TRP:HB3	1.84	0.57
1:B:433:LYS:HA	1:B:433:LYS:CE	2.35	0.56
1:A:501:PRO:O	1:A:526:LYS:NZ	2.37	0.56
1:A:570:VAL:HG21	1:A:622:PHE:HE1	1.71	0.56
1:B:112:ASN:OD1	1:B:113:ASN:N	2.37	0.56
1:B:15:LEU:HD22	1:A:228:LEU:HD13	1.88	0.56
1:B:566:ARG:HG3	1:B:629:PRO:HD3	1.88	0.55
1:A:552:GLU:HG3	1:A:557:ARG:HE	1.72	0.55
1:B:598:ARG:NH1	1:B:599:VAL:O	2.40	0.55
1:B:53:GLY:O	1:B:57:ALA:N	2.38	0.55
1:A:368:PRO:HA	1:A:491:THR:HB	1.87	0.54
1:A:563:ASP:HB3	1:A:607:ALA:HB3	1.88	0.54
1:A:245:ARG:HG3	1:A:247:ASP:HB2	1.88	0.54
1:A:517:LEU:HD21	1:A:528:TYR:HD2	1.72	0.54
1:B:376:SER:O	1:B:450:LYS:NZ	2.40	0.54
1:B:171:ARG:NH1	1:B:372:GLN:O	2.41	0.54
1:B:616:ARG:HA	1:B:617:ASP:C	2.28	0.54
1:B:577:GLU:OE1	1:B:580:ARG:NH2	2.41	0.54
1:B:566:ARG:HH22	1:B:606:GLY:HA3	1.73	0.54
1:B:216:ALA:HB1	1:B:287:PHE:HB2	1.90	0.54
1:B:280:ARG:NH1	2:B:701:ATP:O1G	2.40	0.53
1:A:616:ARG:HG2	1:A:617:ASP:N	2.23	0.53
1:B:545:ARG:HH22	1:B:610:ALA:HB1	1.71	0.53
1:B:347:ALA:HB3	1:B:354:ASP:HB3	1.89	0.53
1:B:433:LYS:HE3	1:B:433:LYS:HA	1.91	0.53
1:A:430:LYS:NZ	1:A:435:GLY:O	2.31	0.53
1:A:570:VAL:O	1:A:638:TYR:OH	2.18	0.53
1:B:105:HIS:NE2	1:B:110:HIS:O	2.37	0.52
1:A:590:THR:HG22	1:A:594:VAL:HG23	1.91	0.52
1:B:198:SER:CB	1:B:399:THR:HG23	2.39	0.52
1:A:569:LEU:HD11	1:A:583:MET:HB2	1.92	0.52
1:B:566:ARG:NH2	1:B:606:GLY:HA3	2.25	0.52
1:A:98:LEU:HA	1:A:103:PRO:HA	1.91	0.52
1:B:98:LEU:HA	1:B:103:PRO:HA	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:128:THR:HA	1:A:129:SER:N	2.09	0.52
1:A:519:VAL:HG13	1:A:526:LYS:HB2	1.91	0.52
1:A:376:SER:HB3	1:A:485:LEU:HD12	1.91	0.52
1:B:32:GLY:HA2	1:B:35:ARG:HE	1.75	0.52
1:A:120:ARG:HH21	1:A:264:LYS:HA	1.75	0.52
1:B:634:VAL:HA	1:B:637:LEU:HD12	1.91	0.51
1:A:548:TYR:CE2	1:A:550:SER:HA	2.45	0.51
1:A:128:THR:HG22	1:A:129:SER:N	2.25	0.51
1:B:321:TRP:HZ2	1:B:513:LEU:HB2	1.76	0.51
1:A:118:VAL:HG22	1:A:119:LYS:N	2.26	0.51
1:B:417:GLN:OE1	1:B:433:LYS:HD3	2.10	0.51
1:B:187:PHE:HD1	1:B:315:ASN:HD22	1.59	0.51
1:A:499:GLU:O	1:A:501:PRO:HD3	2.11	0.51
1:B:366:ILE:HG22	1:B:368:PRO:HD3	1.93	0.50
1:B:575:PHE:HE1	1:B:635:VAL:HG22	1.77	0.50
1:B:317:TYR:CE1	1:B:444:THR:HG21	2.47	0.50
1:B:532:LEU:HB2	1:B:537:TRP:CE2	2.46	0.50
1:B:50:GLU:HG2	1:B:51:GLN:N	2.27	0.50
1:A:36:LYS:O	1:A:656:LEU:HD11	2.11	0.50
1:B:209:LYS:O	1:B:298:GLN:NE2	2.45	0.50
1:A:67:GLU:OE1	1:A:145:ARG:NH2	2.44	0.50
1:A:269:ARG:HG3	1:A:543:THR:HG21	1.94	0.50
1:A:153:TYR:HE2	1:A:545:ARG:HE	1.59	0.49
1:B:131:ALA:O	1:B:541:TRP:NE1	2.44	0.49
1:A:128:THR:CG2	1:A:129:SER:N	2.74	0.49
1:A:120:ARG:HD3	1:A:422:ILE:O	2.12	0.49
1:B:636:SER:OG	1:B:642:HIS:O	2.30	0.49
1:A:542:VAL:O	1:A:544:PRO:HD3	2.13	0.49
1:A:363:LEU:CD2	1:A:499:GLU:HG2	2.42	0.49
1:B:417:GLN:HG2	1:B:429:ARG:HG2	1.95	0.49
1:A:168:GLN:HA	1:A:375:GLY:HA3	1.95	0.49
1:A:497:VAL:C	1:A:498:ASN:CA	2.75	0.49
1:B:266:LYS:O	1:B:270:TYR:OH	2.23	0.49
1:B:95:CYS:HB3	1:B:97:ILE:H	1.78	0.49
1:A:95:CYS:SG	1:A:97:ILE:HG13	2.53	0.48
1:B:346:ILE:HD11	1:B:459:ALA:HB2	1.95	0.48
1:B:53:GLY:O	1:B:57:ALA:HB3	2.13	0.48
1:A:443:THR:O	1:A:443:THR:CG2	2.59	0.48
1:B:248:LEU:O	1:B:252:ARG:HD2	2.13	0.48
1:B:117:SER:OG	1:B:118:VAL:N	2.46	0.48
1:A:222:ARG:NH1	1:A:226:ASP:O	2.45	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:365:ARG:NH1	1:A:508:THR:OG1	2.46	0.48
1:A:337:GLY:HA2	1:A:342:LYS:HZ2	1.78	0.48
1:A:562:PHE:HE2	1:A:587:ILE:HG23	1.79	0.48
1:B:666:ASP:OD1	1:B:667:THR:N	2.46	0.48
1:A:164:ARG:CZ	1:A:278:LYS:HE3	2.44	0.48
1:B:262:GLU:HB3	1:B:284:SER:HB3	1.94	0.48
1:B:50:GLU:HA	1:B:138:PHE:HA	1.95	0.48
1:A:130:ASP:HA	1:A:133:THR:HB	1.96	0.47
1:B:498:ASN:ND2	1:B:501:PRO:HA	2.29	0.47
1:A:94:VAL:HG13	1:A:105:HIS:CD2	2.47	0.47
1:B:370:PHE:HB2	1:B:373:MET:HB2	1.96	0.47
1:A:216:ALA:HB1	1:A:287:PHE:HB2	1.95	0.47
1:A:611:TYR:CE2	1:A:613:PHE:HB2	2.49	0.47
1:B:315:ASN:OD1	1:B:317:TYR:N	2.47	0.47
1:B:324:GLY:O	1:B:328:ASP:N	2.43	0.47
1:A:381:THR:O	1:A:385:VAL:HG23	2.14	0.47
1:B:224:LYS:HD3	1:B:287:PHE:HE1	1.80	0.47
1:A:599:VAL:O	1:A:604:GLY:HA2	2.15	0.47
1:B:210:THR:OG1	1:B:224:LYS:NZ	2.43	0.47
1:B:316:ALA:O	1:B:318:GLY:N	2.48	0.47
1:B:197:VAL:HG13	1:B:404:TRP:NE1	2.30	0.46
1:A:51:GLN:HG3	1:A:139:GLU:C	2.36	0.46
1:A:414:MET:HA	1:A:418:SER:HB2	1.97	0.46
1:B:198:SER:CB	1:B:399:THR:CG2	2.93	0.46
1:B:454:ALA:HA	1:B:483:HIS:CE1	2.50	0.46
1:A:178:VAL:HB	1:A:453:LEU:HD21	1.97	0.46
1:B:398:PRO:O	1:B:401:ARG:N	2.48	0.46
1:A:557:ARG:O	1:A:561:LEU:HG	2.16	0.46
1:A:574:VAL:HG21	1:A:631:TYR:HA	1.98	0.46
1:A:34:LYS:HG3	1:A:35:ARG:N	2.31	0.46
1:A:569:LEU:HD12	1:A:573:ALA:HB3	1.97	0.46
1:B:270:TYR:HE1	1:B:280:ARG:HG3	1.81	0.46
1:B:72:VAL:O	1:B:76:GLN:N	2.44	0.46
1:B:124:ALA:N	1:B:125:GLY:CA	2.75	0.46
1:A:346:ILE:HD12	1:A:455:TYR:HB3	1.98	0.46
1:A:468:ASN:HB2	1:A:474:TYR:CD2	2.50	0.46
1:B:139:GLU:OE1	1:B:578:ARG:NH2	2.47	0.46
1:B:587:ILE:O	1:B:590:THR:HG22	2.16	0.46
1:B:223:THR:OG1	1:B:226:ASP:OD2	2.34	0.46
1:A:230:ASP:O	1:A:234:CYS:HB3	2.15	0.46
1:B:338:GLU:H	1:B:342:LYS:HZ2	1.64	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:200:GLU:O	1:B:204:LEU:HB2	2.16	0.45
1:B:320:SER:HB2	1:B:614:LEU:O	2.16	0.45
1:B:399:THR:OG1	1:A:668:ARG:HD2	2.17	0.45
1:A:314:PHE:CE2	1:A:346:ILE:HD11	2.51	0.45
1:A:222:ARG:HG3	1:A:227:ALA:HB2	1.98	0.45
1:A:283:PHE:N	1:A:438:THR:O	2.38	0.45
1:A:346:ILE:HG22	1:A:355:ILE:HG23	1.99	0.45
1:B:372:GLN:OE1	1:B:372:GLN:HA	2.16	0.45
1:B:139:GLU:HG3	1:B:534:PHE:HZ	1.81	0.45
1:B:362:LYS:HB2	1:B:364:TYR:CE1	2.51	0.45
1:B:95:CYS:HB3	1:B:97:ILE:HD12	1.99	0.45
1:A:128:THR:C	1:A:129:SER:CB	2.74	0.45
1:B:249:THR:HA	1:B:252:ARG:CG	2.46	0.45
1:B:212:ALA:HB1	1:B:225:ARG:HG3	1.98	0.45
1:B:218:PRO:HG3	1:B:221:TRP:CH2	2.51	0.45
1:A:591:ALA:HB3	1:A:594:VAL:HG22	1.99	0.45
1:B:423:ASP:OD1	1:B:424:GLY:N	2.49	0.45
1:A:317:TYR:HA	1:A:349:TYR:CD1	2.52	0.44
1:B:244:VAL:HA	1:A:671:ILE:HG22	1.98	0.44
1:B:570:VAL:HG11	1:B:620:PHE:CE2	2.53	0.44
1:B:518:LYS:HD2	1:B:531:ASN:ND2	2.33	0.44
1:A:270:TYR:HE1	1:A:280:ARG:HG3	1.83	0.44
1:B:198:SER:HB2	1:B:399:THR:CG2	2.43	0.43
1:A:332:TRP:NE1	1:A:345:ARG:HD2	2.33	0.43
1:A:171:ARG:HG3	1:A:375:GLY:O	2.18	0.43
1:A:532:LEU:HB2	1:A:537:TRP:CE2	2.53	0.43
1:B:562:PHE:CD2	1:B:590:THR:HG21	2.53	0.43
1:B:442:GLY:O	1:B:445:LEU:N	2.47	0.43
1:B:545:ARG:NH2	1:B:610:ALA:HB1	2.32	0.43
1:B:354:ASP:HB2	1:B:512:PHE:CD1	2.53	0.43
1:A:520:VAL:HG11	1:A:575:PHE:CD1	2.54	0.43
1:B:417:GLN:CD	1:B:433:LYS:HD3	2.39	0.43
1:B:346:ILE:HG21	1:B:455:TYR:HB3	2.00	0.43
1:B:399:THR:O	1:A:668:ARG:CG	2.59	0.43
1:A:30:LEU:HA	1:A:33:LEU:HB2	2.00	0.43
1:A:430:LYS:HB3	1:A:430:LYS:HE2	1.78	0.43
1:B:95:CYS:HA	1:B:96:GLY:HA3	1.65	0.43
1:A:69:LYS:HE3	1:A:69:LYS:HB3	1.76	0.43
1:B:521:ARG:HG2	1:B:522:ARG:N	2.34	0.43
1:A:238:LEU:HD21	1:A:251:LEU:HD11	1.99	0.43
1:A:23:ALA:O	1:A:27:ARG:HG3	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:113:ASN:O	1:B:115:LYS:N	2.52	0.43
1:B:218:PRO:HB3	1:B:257:GLU:HG2	2.00	0.43
1:B:228:LEU:HD21	1:A:19:LEU:HG	2.01	0.43
1:B:534:PHE:CD2	1:B:578:ARG:HB3	2.54	0.43
1:A:211:ASN:OD1	1:A:214:ALA:N	2.51	0.43
1:B:139:GLU:HG2	1:B:582:LEU:HD13	2.00	0.42
1:A:617:ASP:C	1:A:619:VAL:H	2.22	0.42
1:B:607:ALA:HA	1:B:608:PRO:HD3	1.84	0.42
1:B:444:THR:O	1:B:448:THR:OG1	2.28	0.42
1:A:42:ARG:NH2	1:A:587:ILE:O	2.52	0.42
1:B:332:TRP:CE2	1:B:345:ARG:HD2	2.55	0.42
1:B:76:GLN:OE1	1:B:123:ALA:O	2.38	0.42
1:B:15:LEU:HD21	1:A:287:PHE:HZ	1.84	0.42
1:A:137:ARG:HG2	1:A:138:PHE:CD2	2.55	0.42
1:B:504:GLY:HA2	1:B:518:LYS:HE2	2.02	0.42
1:B:470:LEU:O	1:B:492:TRP:NE1	2.48	0.42
1:B:662:LYS:HB2	1:B:662:LYS:HE3	1.74	0.42
1:A:240:TYR:O	1:A:244:VAL:HG23	2.18	0.42
1:A:599:VAL:HG11	1:A:606:GLY:HA2	2.01	0.42
1:B:132:LEU:HD13	1:B:542:VAL:HG13	2.01	0.42
1:B:78:LEU:HD11	1:B:159:ALA:H	1.84	0.41
1:B:591:ALA:O	1:B:595:VAL:HG23	2.20	0.41
1:B:92:TYR:CD2	1:B:256:PRO:HB2	2.54	0.41
1:B:307:MET:HE3	1:B:347:ALA:HB1	2.02	0.41
1:A:263:CYS:HB2	1:A:283:PHE:CE2	2.56	0.41
1:A:33:LEU:O	1:A:37:LEU:HB2	2.21	0.41
1:A:204:LEU:HD23	1:A:204:LEU:HA	1.91	0.41
1:A:264:LYS:HB3	1:A:264:LYS:HE2	1.85	0.41
1:B:180:ARG:NH1	1:B:391:ASP:OD2	2.54	0.41
1:A:549:ARG:HB3	1:A:552:GLU:HB3	2.03	0.41
1:A:51:GLN:HG3	1:A:139:GLU:O	2.21	0.41
1:B:317:TYR:HA	1:B:349:TYR:CE1	2.55	0.41
1:B:117:SER:HA	1:B:425:THR:HA	2.03	0.41
1:B:218:PRO:HG3	1:B:221:TRP:CZ2	2.55	0.41
1:B:503:PHE:CD2	1:B:521:ARG:HB2	2.56	0.41
1:B:272:VAL:HA	1:B:275:LEU:HG	2.01	0.41
1:A:650:PHE:HA	1:A:651:PRO:HD3	1.86	0.41
1:A:411:TRP:NE1	1:A:441:VAL:O	2.42	0.41
1:A:499:GLU:O	1:A:501:PRO:CD	2.70	0.40
1:B:110:HIS:HA	1:B:111:ASN:O	2.21	0.40
1:A:95:CYS:HA	1:A:96:GLY:HA3	1.75	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:478:PHE:O	1:B:482:LYS:HB2	2.21	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:53:GLY:N	1:A:55:GLU:OE1[2_555]	1.93	0.27
1:A:53:GLY:CA	1:A:55:GLU:OE1[2_555]	2.11	0.09

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	658/705 (93%)	600 (91%)	46 (7%)	12 (2%)	11	53
1	B	662/705 (94%)	585 (88%)	68 (10%)	9 (1%)	14	58
All	All	1320/1410 (94%)	1185 (90%)	114 (9%)	21 (2%)	12	55

All (21) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	38	ALA
1	B	123	ALA
1	B	317	TYR
1	B	399	THR
1	B	550	SER
1	A	58	GLN
1	A	376	SER
1	B	114	GLY
1	B	196	ASP
1	A	113	ASN
1	A	196	ASP

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Mol	Chain	Res	Type
1	A	492	TRP
1	A	602	GLY
1	B	55	GLU
1	A	125	GLY
1	A	317	TYR
1	A	372	GLN
1	A	441	VAL
1	B	252	ARG
1	A	119	LYS
1	A	57	ALA

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	557/598 (93%)	547 (98%)	10 (2%)	66	88
1	B	560/598 (94%)	550 (98%)	10 (2%)	66	88
All	All	1117/1196 (93%)	1097 (98%)	20 (2%)	66	88

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	52	CYS
1	B	56	SER
1	B	78	LEU
1	B	115	LYS
1	B	129	SER
1	B	145	ARG
1	B	172	LYS
1	B	197	VAL
1	B	200	GLU
1	B	523	GLU
1	A	30	LEU
1	A	78	LEU
1	A	120	ARG

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Mol	Chain	Res	Type
1	A	181	LYS
1	A	221	TRP
1	A	366	ILE
1	A	485	LEU
1	A	616	ARG
1	A	656	LEU
1	A	661	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	76	GLN

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 1 is monoatomic - leaving 7 for Mogul analysis.

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	$\# Z > 2$
3	SO4	A	702	-	4,4,4	0.28	0	6,6,6	0.06	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	PO4	A	703	-	4,4,4	0.71	0	6,6,6	0.23	0
4	PO4	A	704	-	4,4,4	0.73	0	6,6,6	0.24	0
4	PO4	A	705	-	4,4,4	0.68	0	6,6,6	0.23	0
2	ATP	B	701	-	26,33,33	0.99	1 (3%)	26,52,52	1.65	1 (3%)
3	SO4	B	702	-	4,4,4	0.28	0	6,6,6	0.06	0
4	PO4	B	703	-	4,4,4	0.75	0	6,6,6	0.23	0

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
3	SO4	A	702	-	-	0/0/0/0	0/0/0/0
4	PO4	A	703	-	-	0/0/0/0	0/0/0/0
4	PO4	A	704	-	-	0/0/0/0	0/0/0/0
4	PO4	A	705	-	-	0/0/0/0	0/0/0/0
2	ATP	B	701	-	-	0/18/38/38	0/3/3/3
3	SO4	B	702	-	-	0/0/0/0	0/0/0/0
4	PO4	B	703	-	-	0/0/0/0	0/0/0/0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	701	ATP	C5-C4	3.21	1.47	1.40

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	701	ATP	N3-C2-N1	-6.59	123.69	128.87

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	701	ATP	2	0
4	B	703	PO4	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	6
1	B	2

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	50:GLU	C	51:GLN	N	2.12
1	A	57:ALA	C	58:GLN	N	1.87
1	A	497:VAL	C	498:ASN	N	1.70
1	B	122:PHE	C	123:ALA	N	1.63
1	A	129:SER	C	130:ASP	N	1.11
1	A	496:LEU	C	497:VAL	N	1.04
1	B	128:THR	C	129:SER	N	0.98
1	A	128:THR	C	129:SER	N	0.80

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	662/705 (93%)	-0.04	5 (0%) 87 80	11, 32, 56, 83	0
1	B	659/705 (93%)	-0.09	3 (0%) 91 88	15, 30, 51, 77	0
All	All	1321/1410 (93%)	-0.07	8 (0%) 90 85	11, 31, 54, 83	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	524	ASN	2.5
1	A	550	SER	2.5
1	A	600	GLN	2.4
1	A	552	GLU	2.1
1	A	548	TYR	2.1
1	A	645	ASP	2.1
1	B	552	GLU	2.0
1	B	553	THR	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(Å ²)	Q<0.9
2	ATP	B	701	31/31	0.73	0.41	3.99	48,79,158,168	0
4	PO4	A	705	5/5	0.83	0.32	1.14	39,42,49,53	0
4	PO4	A	703	5/5	0.90	0.26	0.10	38,39,40,44	0
3	SO4	B	702	5/5	0.93	0.24	-0.24	36,36,43,45	0
3	SO4	A	702	5/5	0.93	0.20	-1.38	31,32,36,36	0
5	MG	A	701	1/1	0.89	0.13	-	18,18,18,18	0
4	PO4	A	704	5/5	0.94	0.19	-	31,32,33,34	0
4	PO4	B	703	5/5	0.93	0.20	-	37,38,41,43	0

6.5 Other polymers [i](#)

There are no such residues in this entry.