



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 11:33 AM UTC

PDB ID : 6S2V / pdb_00006s2v
Title : Structure of the N-terminal catalytic region of T. thermophilus Rel
Authors : Garcia-Pino, A.
Deposited on : 2019-06-22
Resolution : 2.96 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

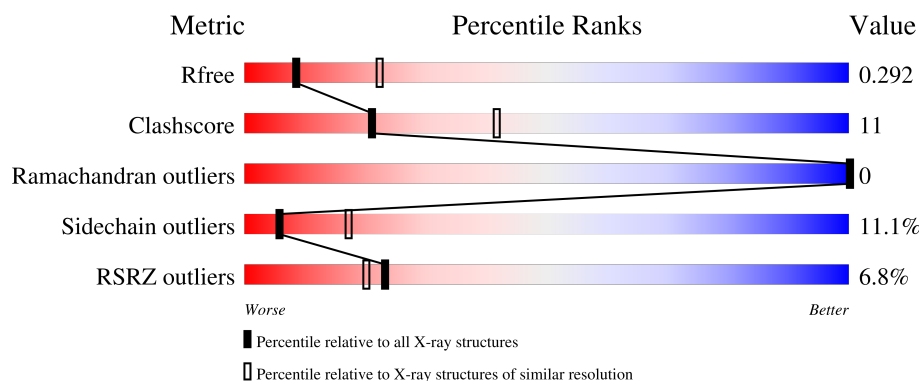
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.96 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1130 (2.98-2.94)
Clashscore	190562	1157 (2.98-2.94)
Ramachandran outliers	187476	1101 (2.98-2.94)
Sidechain outliers	187428	1101 (2.98-2.94)
RSRZ outliers	180081	1130 (2.98-2.94)

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

Mol	Chain	Length	Quality of chain
1	A	356	8% 75% 20% • •
1	B	356	7% 65% 22% • 8%
1	C	356	3% 69% 23% • 6%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 7576 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 (P)ppGpp synthetase I, SpoT/RelA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	342	Total	C	N	O	S	0	1	0
			2588	1649	465	466	8			
1	B	326	Total	C	N	O	S	0	0	0
			2386	1513	434	433	6			
1	C	336	Total	C	N	O	S	0	1	0
			2523	1605	459	451	8			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	ALA	-	expression tag	UNP F6DES6
A	1	SER	-	expression tag	UNP F6DES6
B	0	ALA	-	expression tag	UNP F6DES6
B	1	SER	-	expression tag	UNP F6DES6
C	0	ALA	-	expression tag	UNP F6DES6
C	1	SER	-	expression tag	UNP F6DES6

- Molecule 2 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Mn	0	0
			1	1		
2	B	1	Total	Mn	0	0
			1	1		
2	C	1	Total	Mn	0	0
			1	1		

- Molecule 3 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

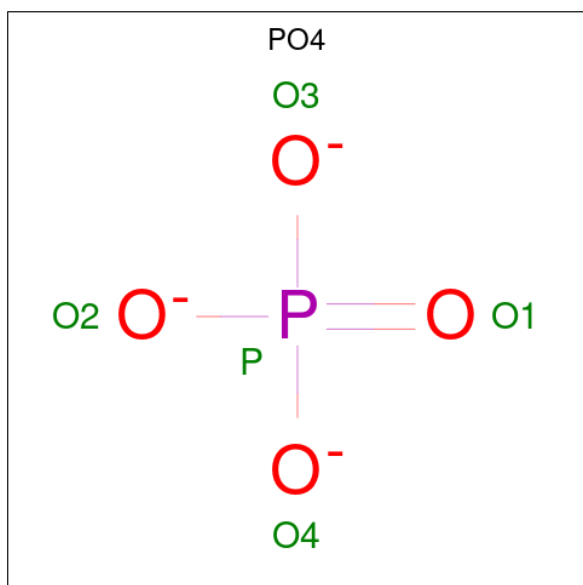
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Cl	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Cl	0	0
			1	1		
3	C	4	Total	Cl	0	0
			4	4		

- Molecule 4 is PHOSPHATE ION (CCD ID: PO4) (formula: O_4P).



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

- Molecule 5 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	C	1	Total	Na	0	0
			1	1		

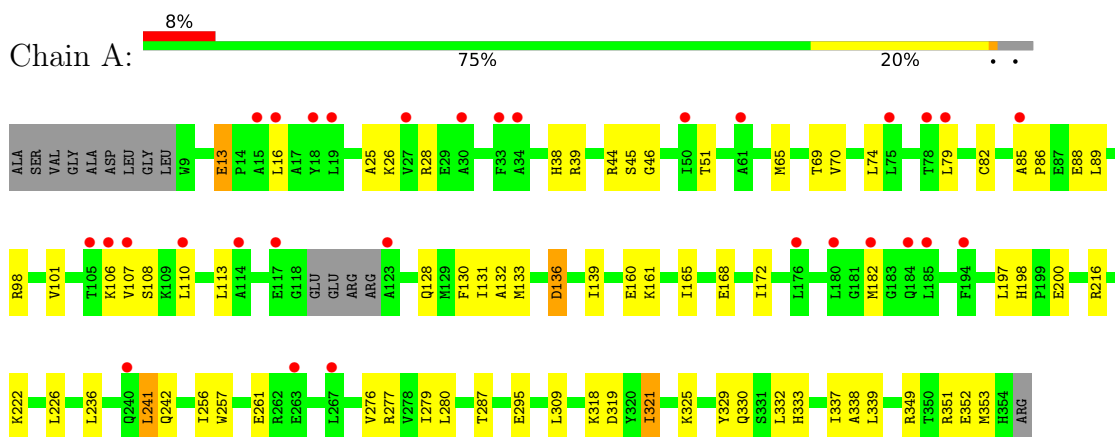
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	27	Total	O	0	1
			28	28		
6	B	11	Total	O	0	0
			11	11		
6	C	25	Total	O	0	0
			25	25		

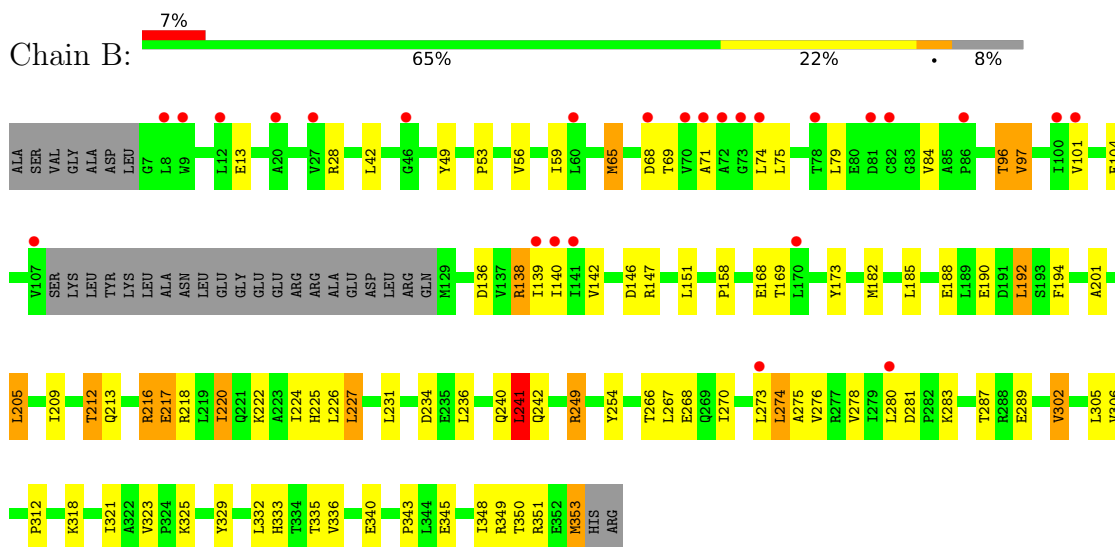
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes 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.

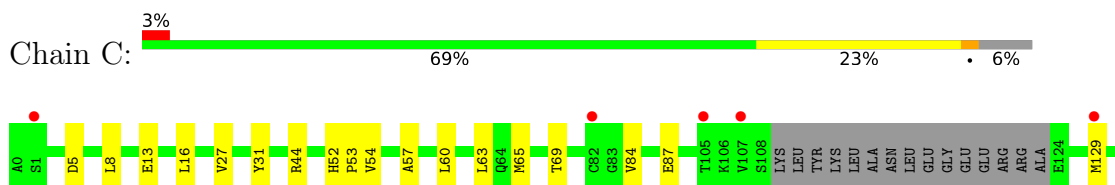
- Molecule 1: (P)ppGpp synthetase I, SpoT/RelA

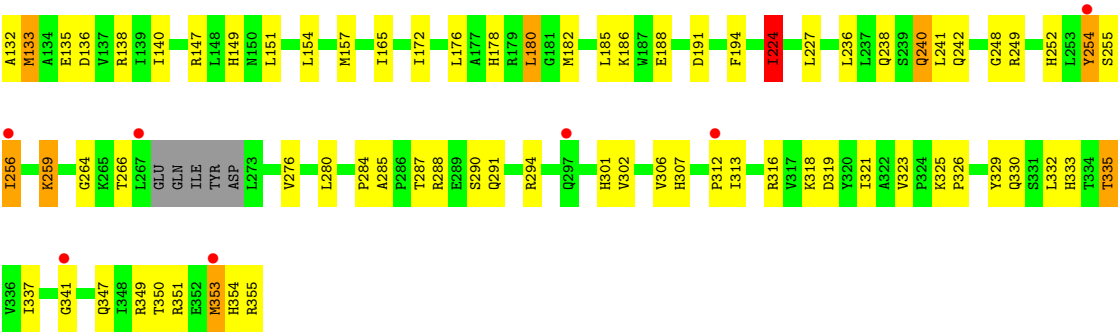


- Molecule 1: (P)ppGpp synthetase I, SpoT/RelA



- Molecule 1: (P)ppGpp synthetase I, SpoT/RelA





4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	105.74Å 105.74Å 241.46Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.77 – 2.96 40.77 – 2.96	Depositor EDS
% Data completeness (in resolution range)	72.4 (40.77-2.96) 72.4 (40.77-2.96)	Depositor EDS
R_{merge}	0.21	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.03 (at 2.95Å)	Xtriage
Refinement program	BUSTER 2.10.3	Depositor
R, R_{free}	0.221 , 0.275 0.244 , 0.292	Depositor DCC
R_{free} test set	1066 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	84.7	Xtriage
Anisotropy	0.085	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 119.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	7576	wwPDB-VP
Average B, all atoms (Å ²)	101.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.44% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: MN, CL, NA, PO4

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	A	0.86	1/2643 (0.0%)	1.48	18/3599 (0.5%)
1	B	0.91	1/2432 (0.0%)	1.45	11/3320 (0.3%)
1	C	0.90	1/2577 (0.0%)	1.48	13/3516 (0.4%)
All	All	0.89	3/7652 (0.0%)	1.47	42/10435 (0.4%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	353	MET	SD-CE	-5.34	1.66	1.79
1	B	158	PRO	CA-C	5.33	1.54	1.51
1	C	323	VAL	CA-C	5.11	1.58	1.52

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	255	SER	N-CA-C	-8.70	102.68	113.38
1	A	321	ILE	N-CA-C	-7.46	103.51	110.53
1	C	242	GLN	N-CA-C	-6.92	105.38	114.31
1	C	319	ASP	CA-CB-CG	6.61	119.21	112.60
1	A	338	ALA	N-CA-C	6.35	113.99	108.78
1	C	5	ASP	CA-CB-CG	5.98	118.58	112.60
1	B	234	ASP	CA-C-N	5.87	128.06	120.44
1	B	234	ASP	C-N-CA	5.87	128.06	120.44
1	A	46	GLY	CA-C-N	5.81	128.47	120.39
1	A	46	GLY	C-N-CA	5.81	128.47	120.39
1	C	172	ILE	N-CA-C	5.68	116.11	111.62
1	B	194	PHE	CA-C-N	5.67	127.87	120.28
1	B	194	PHE	C-N-CA	5.67	127.87	120.28
1	C	31	TYR	CA-C-N	5.54	128.02	120.54
1	C	31	TYR	C-N-CA	5.54	128.02	120.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	132	ALA	CA-C-N	5.48	127.57	120.44
1	A	132	ALA	C-N-CA	5.48	127.57	120.44
1	C	254	TYR	CA-C-N	5.48	131.74	122.26
1	C	254	TYR	C-N-CA	5.48	131.74	122.26
1	B	241	LEU	N-CA-C	5.46	117.80	108.90
1	C	63	LEU	N-CA-C	-5.45	106.68	113.38
1	A	113	LEU	CA-C-N	5.35	127.45	120.28
1	A	113	LEU	C-N-CA	5.35	127.45	120.28
1	A	161	LYS	CA-C-N	5.30	127.38	120.28
1	A	161	LYS	C-N-CA	5.30	127.38	120.28
1	A	257	TRP	CA-C-N	5.28	127.30	120.44
1	A	257	TRP	C-N-CA	5.28	127.30	120.44
1	C	224	ILE	N-CA-CB	5.26	117.69	110.54
1	C	288	ARG	CA-C-N	5.24	127.25	120.44
1	C	288	ARG	C-N-CA	5.24	127.25	120.44
1	A	287	THR	CA-C-N	5.23	127.29	120.28
1	A	287	THR	C-N-CA	5.23	127.29	120.28
1	B	222	LYS	CA-C-N	5.20	127.24	120.28
1	B	222	LYS	C-N-CA	5.20	127.24	120.28
1	A	319	ASP	CA-CB-CG	5.18	117.78	112.60
1	A	13	GLU	CB-CG-CD	5.15	121.36	112.60
1	B	65	MET	CA-C-N	5.14	128.38	120.82
1	B	65	MET	C-N-CA	5.14	128.38	120.82
1	B	217	GLU	CB-CG-CD	5.13	121.33	112.60
1	A	45	SER	N-CA-C	-5.09	102.15	110.20
1	B	146	ASP	CA-CB-CG	5.09	117.69	112.60
1	A	242	GLN	N-CA-C	-5.04	107.26	113.41

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2588	0	2495	38	0
1	B	2386	0	2259	71	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	2523	0	2437	56	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	4	0	0	1	0
4	B	5	0	0	0	0
5	C	1	0	0	0	0
6	A	28	0	0	1	0
6	B	11	0	0	0	0
6	C	25	0	0	0	0
All	All	7576	0	7191	162	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 11.

All (162) 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:C:349:ARG:NH2	1:C:354:HIS:CE1	2.28	1.02
1:B:212:THR:HG23	1:B:216:ARG:HD3	1.38	0.99
1:C:316:ARG:HB3	1:C:335:THR:OG1	1.63	0.96
1:B:138:ARG:HG2	1:B:138:ARG:HH11	1.34	0.89
1:B:287:THR:HG22	1:B:289:GLU:H	1.37	0.89
1:C:349:ARG:NE	1:C:354:HIS:ND1	2.21	0.89
1:C:313:ILE:HD11	1:C:337:ILE:HD11	1.56	0.87
1:A:321:ILE:HD12	1:A:332:LEU:HD13	1.58	0.86
1:C:337:ILE:CG2	1:C:341:GLY:HA2	2.07	0.84
1:C:349:ARG:CZ	1:C:354:HIS:ND1	2.40	0.84
1:C:349:ARG:NH2	1:C:354:HIS:HE1	1.72	0.82
1:B:69:THR:CG2	1:B:138:ARG:HD2	2.09	0.82
1:B:69:THR:HG21	1:B:138:ARG:HD2	1.63	0.81
1:B:79:LEU:HA	1:B:84:VAL:HA	1.62	0.81
1:C:349:ARG:HH21	1:C:354:HIS:CE1	2.02	0.77
1:B:231:LEU:CD2	1:B:302:VAL:HG22	2.16	0.75
1:B:138:ARG:HG2	1:B:138:ARG:NH1	2.02	0.73
1:C:87:GLU:H	1:C:87:GLU:CD	1.98	0.72
1:B:151:LEU:HD23	1:B:192:LEU:HB3	1.73	0.71
1:B:278:VAL:HG12	1:B:280:LEU:CD1	2.21	0.70
1:C:87:GLU:OE1	1:C:87:GLU:N	2.19	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:139:ILE:HG13	1:B:140:ILE:H	1.56	0.70
1:B:273:LEU:CD2	1:B:343:PRO:HB2	2.22	0.70
1:B:69:THR:HG21	1:B:138:ARG:CD	2.22	0.69
1:B:139:ILE:HG13	1:B:140:ILE:N	2.06	0.69
1:B:325:LYS:HB2	1:B:329:TYR:HB3	1.75	0.68
1:A:226:LEU:HD21	1:C:326:PRO:HA	1.74	0.68
1:A:89:LEU:HD13	1:A:101:VAL:HG21	1.76	0.68
1:C:316:ARG:HB3	1:C:335:THR:HG1	1.57	0.67
1:B:231:LEU:HD21	1:B:302:VAL:HG22	1.75	0.67
1:A:279:ILE:HG22	1:A:351:ARG:NH2	2.10	0.67
1:A:279:ILE:CG2	1:A:351:ARG:HD3	2.26	0.66
1:B:318:LYS:HB2	1:B:333:HIS:HB2	1.76	0.66
1:C:321:ILE:HD12	1:C:332:LEU:HD13	1.76	0.66
1:C:178:HIS:HB2	1:C:186:LYS:HD2	1.77	0.65
1:B:273:LEU:HD23	1:B:343:PRO:HB2	1.77	0.65
1:B:212:THR:HG23	1:B:216:ARG:CD	2.21	0.64
1:A:279:ILE:HG21	1:A:351:ARG:HD3	1.80	0.64
1:A:339:LEU:HD12	1:A:339:LEU:N	2.12	0.64
1:B:212:THR:O	1:B:216:ARG:HG2	1.99	0.63
1:A:106:LYS:O	1:A:107:VAL:HG23	1.99	0.63
1:B:242:GLN:HG3	1:B:283:LYS:HA	1.82	0.62
1:B:302:VAL:HG12	1:B:348:ILE:HD13	1.82	0.62
1:C:259:LYS:HG2	1:C:264:GLY:HA2	1.82	0.61
1:C:349:ARG:CZ	1:C:354:HIS:CE1	2.83	0.60
1:C:316:ARG:CB	1:C:335:THR:OG1	2.47	0.60
1:A:339:LEU:H	1:A:339:LEU:CD1	2.14	0.60
1:B:69:THR:CG2	1:B:138:ARG:HG3	2.32	0.60
1:B:97:VAL:HG22	1:B:97:VAL:O	2.01	0.60
1:C:84:VAL:HG12	1:C:84:VAL:O	2.02	0.59
1:A:216:ARG:HD2	6:A:508:HOH:O	2.01	0.59
1:C:57:ALA:HA	1:C:60:LEU:HD12	1.83	0.59
1:B:69:THR:CG2	1:B:138:ARG:CD	2.79	0.58
1:B:278:VAL:HG12	1:B:280:LEU:HD11	1.84	0.58
1:A:136:ASP:HA	1:A:139:ILE:HD12	1.86	0.58
1:C:316:ARG:HH11	1:C:335:THR:HG21	1.69	0.57
1:A:106:LYS:O	1:A:107:VAL:CG2	2.53	0.57
1:B:69:THR:CG2	1:B:138:ARG:CG	2.83	0.57
1:C:337:ILE:HG23	1:C:341:GLY:HA2	1.87	0.57
1:B:69:THR:HG22	1:B:138:ARG:HG3	1.87	0.57
1:A:168:GLU:HG3	1:A:172:ILE:HD12	1.87	0.57
1:C:8:LEU:HB3	1:C:54:VAL:HG11	1.86	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:339:LEU:N	1:A:339:LEU:CD1	2.67	0.56
1:C:194:PHE:CE1	1:C:256:ILE:HD13	2.41	0.56
1:A:106:LYS:C	1:A:107:VAL:HG23	2.31	0.56
1:C:325:LYS:HB2	1:C:329:TYR:HB3	1.88	0.56
1:C:240:GLN:HE21	1:C:294:ARG:HE	1.54	0.56
1:B:231:LEU:HD22	1:B:302:VAL:HG22	1.87	0.55
1:A:318:LYS:HB2	1:A:333:HIS:HB2	1.88	0.55
1:C:280:LEU:HD21	1:C:302:VAL:HG21	1.88	0.55
1:B:138:ARG:HG3	1:B:138:ARG:O	2.06	0.54
1:A:160:GLU:N	1:A:160:GLU:OE1	2.41	0.53
1:A:85:ALA:HB1	1:A:88:GLU:HB3	1.89	0.53
1:B:139:ILE:O	1:B:142:VAL:N	2.41	0.53
1:C:191:ASP:OD1	1:C:252:HIS:O	2.27	0.53
1:B:68:ASP:OD1	1:B:96:THR:HB	2.09	0.53
1:C:326:PRO:HG2	3:C:404:CL:CL	2.46	0.52
1:B:182:MET:HG2	1:B:185:LEU:HB3	1.90	0.52
1:B:249:ARG:HD3	1:B:275:ALA:HB3	1.91	0.52
1:A:82:CYS:SG	1:A:82:CYS:O	2.69	0.50
1:C:349:ARG:NH2	1:C:354:HIS:ND1	2.49	0.50
1:B:220:ILE:HG13	1:B:274:LEU:HD23	1.94	0.50
1:B:225:HIS:HD2	1:B:226:LEU:HD12	1.77	0.50
1:C:227:LEU:HD11	1:C:306:VAL:HG22	1.94	0.49
1:A:13:GLU:HA	1:A:16:LEU:HD12	1.94	0.49
1:B:56:VAL:HA	1:B:59:ILE:HD12	1.94	0.49
1:C:285:ALA:H	1:C:291:GLN:NE2	2.11	0.49
1:B:69:THR:HG22	1:B:138:ARG:CG	2.44	0.48
1:B:69:THR:HG23	1:B:138:ARG:HD2	1.94	0.47
1:B:188:GLU:O	1:B:192:LEU:HB2	2.14	0.47
1:B:213:GLN:O	1:B:217:GLU:HB2	2.14	0.47
1:C:151:LEU:HD23	1:C:154:LEU:HD22	1.97	0.47
1:A:130:PHE:HA	1:A:133:MET:HE3	1.97	0.47
1:B:79:LEU:HB2	1:B:101:VAL:HG21	1.97	0.47
1:B:182:MET:HG3	1:B:185:LEU:H	1.80	0.47
1:A:325:LYS:HB2	1:A:329:TYR:HB3	1.97	0.47
1:B:190:GLU:HB3	1:B:254:TYR:HB3	1.97	0.47
1:C:241:LEU:HD21	1:C:280:LEU:HB3	1.96	0.47
1:B:302:VAL:HG12	1:B:348:ILE:CD1	2.44	0.46
1:C:287:THR:HG23	1:C:290:SER:H	1.80	0.46
1:A:337:ILE:HG21	1:C:133:MET:HE1	1.97	0.46
1:C:330:GLN:CB	1:C:353:MET:HE2	2.45	0.46
1:B:241:LEU:HD11	1:B:280:LEU:HD23	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:38:HIS:CE1	1:A:82:CYS:SG	3.09	0.46
1:B:49:TYR:O	1:B:53:PRO:HD3	2.17	0.45
1:A:65:MET:HE2	1:A:69:THR:HG21	1.98	0.45
1:B:231:LEU:HD22	1:B:302:VAL:CG2	2.47	0.45
1:C:180:LEU:HB3	1:C:182:MET:HG3	1.98	0.44
1:C:87:GLU:CD	1:C:87:GLU:N	2.73	0.44
1:B:71:ALA:HA	1:B:74:LEU:HD12	1.99	0.44
1:B:139:ILE:CG1	1:B:140:ILE:H	2.29	0.44
1:B:321:ILE:HD12	1:B:332:LEU:HD13	2.00	0.43
1:C:333:HIS:CD2	1:C:347:GLN:HG2	2.53	0.43
1:C:318:LYS:HB2	1:C:333:HIS:HB2	1.99	0.43
1:A:226:LEU:HG	1:A:309:LEU:HD11	2.00	0.43
1:A:321:ILE:HA	1:A:330:GLN:HG2	2.00	0.43
1:A:25:ALA:HA	1:A:28:ARG:HG2	2.01	0.43
1:B:312:PRO:HA	1:B:336:VAL:HA	2.01	0.43
1:A:337:ILE:CG2	1:C:133:MET:HE1	2.48	0.43
1:B:227:LEU:HD21	1:B:306:VAL:HG22	1.99	0.43
1:A:222:LYS:HB2	1:A:339:LEU:CD2	2.48	0.43
1:C:353:MET:HE3	1:C:353:MET:HB3	1.80	0.43
1:B:287:THR:HG22	1:B:289:GLU:N	2.19	0.42
1:C:248:GLY:HA2	1:C:276:VAL:HA	2.01	0.42
1:B:169:THR:HA	1:B:173:TYR:HB2	2.01	0.42
1:C:236:LEU:HD22	1:C:301:HIS:CE1	2.54	0.42
1:B:349:ARG:HH12	1:B:351:ARG:NH2	2.18	0.42
1:C:16:LEU:HD11	1:C:27:VAL:HG11	2.00	0.42
1:B:278:VAL:HG12	1:B:280:LEU:HD12	2.00	0.42
1:B:302:VAL:O	1:B:306:VAL:HG23	2.19	0.42
1:C:330:GLN:HB2	1:C:353:MET:HE2	2.02	0.42
1:C:284:PRO:HA	1:C:291:GLN:HE22	1.84	0.42
1:B:69:THR:HG21	1:B:138:ARG:CG	2.48	0.42
1:A:85:ALA:N	1:A:86:PRO:CD	2.83	0.41
1:A:107:VAL:HG12	1:A:108:SER:N	2.35	0.41
1:B:139:ILE:CG1	1:B:140:ILE:N	2.81	0.41
1:B:201:ALA:O	1:B:205:LEU:HB2	2.20	0.41
1:A:197:LEU:HD23	1:A:198:HIS:CE1	2.55	0.41
1:A:351:ARG:HG2	1:A:351:ARG:HH21	1.85	0.41
1:B:79:LEU:HD22	1:B:84:VAL:O	2.21	0.41
1:B:136:ASP:O	1:B:139:ILE:HG12	2.20	0.41
1:B:302:VAL:CG1	1:B:348:ILE:HG12	2.51	0.41
1:B:312:PRO:HA	1:B:336:VAL:HG12	2.03	0.41
1:B:350:THR:H	1:B:353:MET:HE2	1.85	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:44:ARG:H	1:C:44:ARG:HD3	1.85	0.41
1:C:52:HIS:HB3	1:C:53:PRO:HD3	2.02	0.41
1:B:42:LEU:HD12	1:B:42:LEU:C	2.46	0.41
1:B:69:THR:HG21	1:B:138:ARG:HG3	2.00	0.41
1:B:138:ARG:NH1	1:B:138:ARG:CG	2.73	0.41
1:C:52:HIS:HA	1:C:149:HIS:CE1	2.55	0.41
1:A:106:LYS:C	1:A:107:VAL:CG2	2.94	0.41
1:B:147:ARG:HH12	1:B:168:GLU:HB3	1.85	0.41
1:C:350:THR:O	1:C:351:ARG:HB3	2.20	0.41
1:A:128:GLN:HA	1:A:131:ILE:HG22	2.03	0.41
1:A:241:LEU:HD11	1:A:280:LEU:HB3	2.02	0.41
1:B:220:ILE:O	1:B:224:ILE:HG13	2.20	0.41
1:A:70:VAL:O	1:A:74:LEU:HD22	2.21	0.41
1:C:224:ILE:HG22	1:C:276:VAL:HG21	2.03	0.41
1:C:307:HIS:CG	1:C:312:PRO:HG3	2.57	0.41
1:C:313:ILE:CD1	1:C:337:ILE:HD11	2.39	0.41
1:C:132:ALA:O	1:C:135:GLU:O	2.39	0.40
1:C:65:MET:HB3	1:C:69:THR:HB	2.01	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	339/356 (95%)	319 (94%)	20 (6%)	0	100	100
1	B	322/356 (90%)	296 (92%)	26 (8%)	0	100	100
1	C	331/356 (93%)	305 (92%)	26 (8%)	0	100	100
All	All	992/1068 (93%)	920 (93%)	72 (7%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	244/300 (81%)	223 (91%)	21 (9%)	10	25
1	B	219/300 (73%)	185 (84%)	34 (16%)	2	8
1	C	240/300 (80%)	216 (90%)	24 (10%)	7	20
All	All	703/900 (78%)	624 (89%)	79 (11%)	6	17

All (79) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	26	LYS
1	A	39	ARG
1	A	44	ARG
1	A	51	THR
1	A	79	LEU
1	A	98	ARG
1	A	110	LEU
1	A	136	ASP
1	A	165	ILE
1	A	182	MET
1	A	200	GLU
1	A	236	LEU
1	A	241	LEU
1	A	256	ILE
1	A	261[A]	GLU
1	A	261[B]	GLU
1	A	276	VAL
1	A	277	ARG
1	A	295	GLU
1	A	349	ARG
1	A	352	GLU
1	B	13	GLU
1	B	28	ARG
1	B	65	MET
1	B	75	LEU
1	B	96	THR

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Mol	Chain	Res	Type
1	B	97	VAL
1	B	104	GLU
1	B	138	ARG
1	B	192	LEU
1	B	205	LEU
1	B	209	ILE
1	B	212	THR
1	B	216	ARG
1	B	218	ARG
1	B	220	ILE
1	B	227	LEU
1	B	236	LEU
1	B	240	GLN
1	B	241	LEU
1	B	249	ARG
1	B	266	THR
1	B	267	LEU
1	B	268	GLU
1	B	270	ILE
1	B	274	LEU
1	B	276	VAL
1	B	281	ASP
1	B	302	VAL
1	B	305	LEU
1	B	323	VAL
1	B	335	THR
1	B	340	GLU
1	B	345	GLU
1	B	353	MET
1	C	13	GLU
1	C	129	MET
1	C	133	MET
1	C	136	ASP
1	C	138	ARG
1	C	140	ILE
1	C	147	ARG
1	C	157	MET
1	C	165	ILE
1	C	176	LEU
1	C	180	LEU
1	C	185	LEU
1	C	188	GLU

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Mol	Chain	Res	Type
1	C	224	ILE
1	C	238	GLN
1	C	240	GLN
1	C	249	ARG
1	C	254	TYR
1	C	256	ILE
1	C	259	LYS
1	C	266	THR
1	C	335	THR
1	C	353	MET
1	C	355	ARG

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

Mol	Chain	Res	Type
1	A	150	ASN
1	A	213	GLN
1	A	238	GLN
1	A	269	GLN
1	A	347	GLN
1	B	149	HIS
1	B	156	HIS
1	B	162	GLN
1	B	221	GLN
1	B	225	HIS
1	B	238	GLN
1	B	347	GLN
1	C	38	HIS
1	C	221	GLN
1	C	228	GLN
1	C	240	GLN
1	C	242	GLN
1	C	291	GLN
1	C	301	HIS
1	C	307	HIS

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 11 ligands modelled in this entry, 10 are monoatomic - leaving 1 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	$\# Z > 2$
4	PO4	B	402	-	4,4,4	2.56	3 (75%)	6,6,6	0.55	0

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	402	PO4	P-O1	4.23	1.60	1.50
4	B	402	PO4	P-O2	2.04	1.60	1.54
4	B	402	PO4	P-O4	2.02	1.60	1.54

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues

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	342/356 (96%)	0.63	30 (8%)	15 14	46, 98, 140, 158	1 (0%)
1	B	326/356 (91%)	0.55	26 (7%)	18 16	45, 113, 204, 224	0
1	C	336/356 (94%)	0.37	12 (3%)	46 38	45, 90, 129, 165	1 (0%)
All	All	1004/1068 (94%)	0.52	68 (6%)	23 20	45, 98, 177, 224	2 (0%)

All (68) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	61	ALA	5.4
1	B	71	ALA	4.6
1	A	105	THR	4.5
1	A	18	TYR	4.4
1	C	1	SER	4.2
1	C	107	VAL	4.0
1	A	16	LEU	4.0
1	A	34	ALA	3.8
1	C	82	CYS	3.7
1	C	129	MET	3.6
1	A	185	LEU	3.6
1	B	46	GLY	3.5
1	B	140	ILE	3.4
1	A	117	GLU	3.4
1	A	85	ALA	3.3
1	A	114	ALA	3.2
1	B	70	VAL	3.2
1	C	297	GLN	3.1
1	B	74	LEU	3.1
1	B	170	LEU	3.1
1	B	139	ILE	3.0
1	B	68	ASP	3.0
1	B	101	VAL	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	50	ILE	2.9
1	A	263	GLU	2.9
1	A	78	THR	2.9
1	B	273	LEU	2.9
1	A	27	VAL	2.8
1	B	20	ALA	2.8
1	A	176	LEU	2.8
1	B	280	LEU	2.7
1	A	75	LEU	2.7
1	C	254	TYR	2.7
1	A	107	VAL	2.6
1	B	8	LEU	2.6
1	B	12	LEU	2.6
1	C	105	THR	2.6
1	B	27	VAL	2.5
1	B	82	CYS	2.5
1	B	78	THR	2.5
1	A	79	LEU	2.4
1	B	9	TRP	2.4
1	B	141	ILE	2.4
1	A	19	LEU	2.4
1	B	73	GLY	2.3
1	A	15	ALA	2.3
1	A	30	ALA	2.3
1	B	81	ASP	2.3
1	A	110	LEU	2.3
1	A	106	LYS	2.3
1	A	194	PHE	2.3
1	B	107	VAL	2.2
1	A	180	LEU	2.2
1	B	86	PRO	2.2
1	C	341	GLY	2.2
1	C	353	MET	2.2
1	B	72	ALA	2.2
1	A	240	GLN	2.2
1	A	182	MET	2.2
1	C	256	ILE	2.2
1	A	184	GLN	2.2
1	C	267	LEU	2.1
1	B	60	LEU	2.1
1	A	33	PHE	2.1
1	C	312	PRO	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	100	ILE	2.0
1	A	123	ALA	2.0
1	A	267	LEU	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	PO4	B	402	5/5	0.78	0.11	150,151,151,152	0
3	CL	C	404	1/1	0.84	0.06	64,64,64,64	0
3	CL	A	402	1/1	0.90	0.10	95,95,95,95	0
3	CL	C	402	1/1	0.91	0.12	51,51,51,51	0
5	NA	C	406	1/1	0.91	0.09	56,56,56,56	0
3	CL	B	403	1/1	0.95	0.04	106,106,106,106	0
3	CL	C	403	1/1	0.98	0.10	67,67,67,67	0
3	CL	C	405	1/1	0.98	0.06	57,57,57,57	0
2	MN	A	401	1/1	0.99	0.02	98,98,98,98	0
2	MN	B	401	1/1	0.99	0.05	120,120,120,120	0
2	MN	C	401	1/1	1.00	0.05	73,73,73,73	0

6.5 Other polymers [i](#)

There are no such residues in this entry.