



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 02:25 PM UTC

PDB ID : 2PGS / pdb_00002pgs
Title : Crystal structure of a putative deoxyguanosinetriphosphate triphosphohydro-
lase from *Pseudomonas syringae* pv. *phaseolicola* 1448A
Authors : Rao, K.N.; Burley, S.K.; Swaminathan, S.; New York SGX Research Center
for Structural Genomics (NYSGXRC)
Deposited on : 2007-04-10
Resolution : 2.35 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

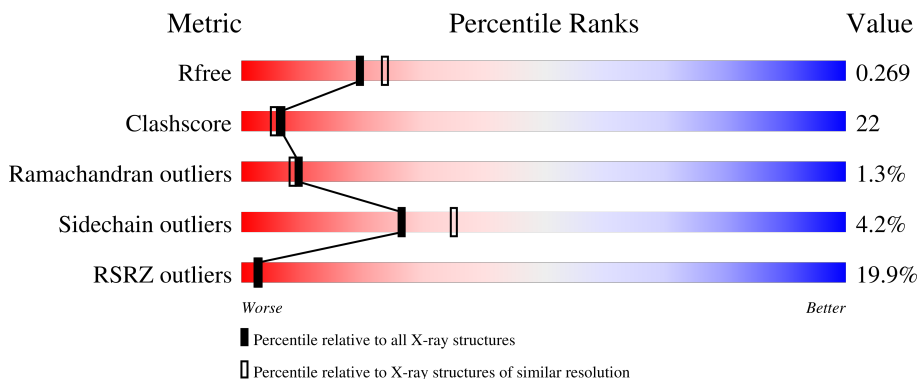
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.35 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	1596 (2.36-2.36)
Clashscore	190562	1663 (2.36-2.36)
Ramachandran outliers	187476	1646 (2.36-2.36)
Sidechain outliers	187428	1646 (2.36-2.36)
RSRZ outliers	180081	1598 (2.36-2.36)

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	451	<div style="display: flex; align-items: center;"> <div style="width: 17%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 52%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 32%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 13%; height: 10px; background-color: grey; margin-right: 5px;"></div> </div> <p style="font-size: small; margin-top: 5px;">17% 52% 32% • 13%</p>

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 3105 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 Putative deoxyguanosinetriphosphate triphosphohydrolase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	A	393	3056	1934	546	558	7	11	0	0	0

There are 21 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	-	cloning artifact	UNP Q48JX0
A	2	SER	-	cloning artifact	UNP Q48JX0
A	79	MSE	MET	modified residue	UNP Q48JX0
A	100	MSE	MET	modified residue	UNP Q48JX0
A	140	MSE	MET	modified residue	UNP Q48JX0
A	239	MSE	MET	modified residue	UNP Q48JX0
A	258	MSE	MET	modified residue	UNP Q48JX0
A	334	MSE	MET	modified residue	UNP Q48JX0
A	348	MSE	MET	modified residue	UNP Q48JX0
A	419	MSE	MET	modified residue	UNP Q48JX0
A	426	MSE	MET	modified residue	UNP Q48JX0
A	434	MSE	MET	modified residue	UNP Q48JX0
A	438	MSE	MET	modified residue	UNP Q48JX0
A	444	GLU	-	cloning artifact	UNP Q48JX0
A	445	GLY	-	cloning artifact	UNP Q48JX0
A	446	HIS	-	cloning artifact	UNP Q48JX0
A	447	HIS	-	cloning artifact	UNP Q48JX0
A	448	HIS	-	cloning artifact	UNP Q48JX0
A	449	HIS	-	cloning artifact	UNP Q48JX0
A	450	HIS	-	cloning artifact	UNP Q48JX0
A	451	HIS	-	cloning artifact	UNP Q48JX0

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	49	Total	O	0	0
			49	49		

4 Data and refinement statistics

Property	Value	Source
Space group	P 63 2 2	Depositor
Cell constants a, b, c, α , β , γ	103.73Å 103.73Å 159.12Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	49.31 – 2.35 49.31 – 2.35	Depositor EDS
% Data completeness (in resolution range)	97.7 (49.31-2.35) 97.7 (49.31-2.35)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.59 (at 2.32Å)	Xtrriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.251 , 0.273 0.251 , 0.269	Depositor DCC
R_{free} test set	619 reflections (2.78%)	wwPDB-VP
Wilson B-factor (Å ²)	39.2	Xtrriage
Anisotropy	0.087	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 46.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	3105	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.19% 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 i

5.1 Standard geometry i

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.55	1/3110 (0.0%)	0.97	13/4184 (0.3%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	390	PRO	C-N	-18.48	1.09	1.33

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	313	ALA	N-CA-C	-9.46	103.84	114.62
1	A	164	LEU	N-CA-C	8.64	121.95	111.40
1	A	233	HIS	N-CA-C	-7.30	100.88	109.93
1	A	50	GLN	N-CA-C	-7.11	100.95	110.55
1	A	255	GLY	N-CA-C	-7.09	103.87	113.37
1	A	93	CYS	N-CA-C	7.01	120.33	107.99
1	A	317	GLN	N-CA-C	6.73	119.50	110.88
1	A	294	LYS	N-CA-C	-6.52	104.59	112.54
1	A	109	HIS	N-CA-C	5.91	121.12	111.37
1	A	427	THR	N-CA-C	-5.75	102.79	110.55
1	A	32	ASP	N-CA-C	-5.60	105.08	111.07
1	A	336	GLY	N-CA-C	5.24	123.03	112.34
1	A	358	ARG	N-CA-C	-5.13	105.77	111.36

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	3056	0	2982	131	0
2	A	49	0	0	2	0
All	All	3105	0	2982	131	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 22.

All (131) 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:79:MSE:HE2	1:A:96:SER:HA	1.41	1.03
1:A:373:LEU:HD13	1:A:426:MSE:HE1	1.50	0.93
1:A:189:THR:HG22	1:A:191:ARG:H	1.35	0.90
1:A:317:GLN:O	1:A:321:LEU:HB2	1.79	0.83
1:A:273:LEU:HD23	1:A:341:CYS:SG	2.21	0.81
1:A:423:ILE:HA	1:A:426:MSE:HE2	1.65	0.78
1:A:29:PHE:HE2	1:A:176:THR:HG21	1.48	0.76
1:A:176:THR:HG22	1:A:178:ALA:H	1.49	0.76
1:A:326:LEU:HD12	1:A:327:PRO:HD2	1.70	0.73
1:A:63:ARG:NH2	1:A:113:ASN:HB2	2.02	0.73
1:A:383:GLU:HG3	1:A:394:HIS:ND1	2.05	0.72
1:A:140:MSE:HE2	1:A:413:HIS:HA	1.75	0.69
1:A:185:LYS:HA	1:A:236:VAL:HG11	1.75	0.68
1:A:185:LYS:HE3	1:A:186:TYR:CE1	2.29	0.67
1:A:80:ARG:NH1	1:A:308:ASN:HD21	1.93	0.67
1:A:176:THR:HG22	1:A:178:ALA:N	2.10	0.67
1:A:321:LEU:HD13	1:A:326:LEU:HD22	1.77	0.66
1:A:401:LEU:N	1:A:401:LEU:HD12	2.11	0.66
1:A:80:ARG:NH1	1:A:308:ASN:ND2	2.45	0.65
1:A:122:ASP:HA	1:A:125:ARG:HH12	1.63	0.64
1:A:434:MSE:O	1:A:438:MSE:HB2	1.98	0.64
1:A:273:LEU:C	1:A:273:LEU:HD12	2.23	0.64
1:A:79:MSE:HE2	1:A:96:SER:CA	2.25	0.63
1:A:436:ARG:C	1:A:438:MSE:H	2.07	0.62
1:A:48:LYS:HE2	1:A:426:MSE:O	1.99	0.62
1:A:310:ALA:HA	1:A:334:MSE:HE2	1.82	0.62
1:A:168:GLN:HG2	1:A:169:PHE:CD1	2.35	0.62
1:A:423:ILE:CA	1:A:426:MSE:HE2	2.30	0.61
1:A:129:ASN:OD1	1:A:149:LEU:HD22	2.00	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:122:ASP:HA	1:A:125:ARG:NH1	2.15	0.61
1:A:383:GLU:HG3	1:A:394:HIS:HD1	1.66	0.60
1:A:139:ALA:HB3	1:A:412:LEU:HD22	1.84	0.59
1:A:163:GLN:HG3	1:A:209:GLU:OE1	2.04	0.57
1:A:200:LYS:HD2	1:A:201:HIS:N	2.20	0.57
1:A:350:ARG:HA	1:A:354:PHE:HB3	1.87	0.57
1:A:392:PHE:O	1:A:393:LYS:HB3	2.04	0.56
1:A:332:GLU:HA	1:A:339:LYS:HE3	1.87	0.56
1:A:115:PRO:HG3	1:A:426:MSE:O	2.06	0.56
1:A:241:ALA:O	1:A:245:ILE:HG13	2.05	0.55
1:A:339:LYS:O	1:A:343:LEU:HG	2.07	0.55
1:A:144:GLU:HB3	1:A:413:HIS:CD2	2.43	0.54
1:A:401:LEU:O	1:A:403:ASN:N	2.41	0.54
1:A:3:LEU:HB2	1:A:322:LEU:O	2.07	0.54
1:A:132:ALA:C	1:A:134:ARG:H	2.15	0.53
1:A:210:LEU:HB3	1:A:211:PRO:HD3	1.89	0.53
1:A:140:MSE:HG2	1:A:144:GLU:HB2	1.90	0.52
1:A:29:PHE:CE2	1:A:176:THR:HG21	2.38	0.52
1:A:77:LEU:HD21	1:A:310:ALA:HB3	1.91	0.52
1:A:418:ARG:C	1:A:419:MSE:HE2	2.34	0.52
1:A:52:HIS:HB3	1:A:258:MSE:HE1	1.91	0.52
1:A:98:LEU:O	1:A:101:VAL:HG12	2.10	0.52
1:A:335:HIS:HD2	1:A:337:PRO:HD2	1.74	0.52
1:A:423:ILE:HG23	1:A:426:MSE:HE2	1.91	0.52
1:A:109:HIS:HE1	1:A:185:LYS:NZ	2.08	0.52
1:A:354:PHE:HD1	1:A:359:LYS:HZ3	1.54	0.52
1:A:359:LYS:O	1:A:363:GLU:HG3	2.10	0.51
1:A:273:LEU:CD1	1:A:274:VAL:HG13	2.40	0.51
1:A:423:ILE:HG23	1:A:426:MSE:CE	2.40	0.51
1:A:80:ARG:HH12	1:A:308:ASN:HD21	1.57	0.51
1:A:206:TYR:HB2	1:A:209:GLU:HG2	1.92	0.51
1:A:406:PRO:HD3	1:A:419:MSE:CE	2.41	0.51
1:A:75:ARG:O	1:A:79:MSE:HG3	2.11	0.50
1:A:318:GLN:O	1:A:322:LEU:HB2	2.11	0.50
1:A:267:GLU:O	1:A:271:LEU:HG	2.11	0.50
1:A:406:PRO:HD3	1:A:419:MSE:HE1	1.93	0.50
1:A:383:GLU:HG3	1:A:394:HIS:CE1	2.46	0.50
1:A:6:GLN:HA	1:A:90:PRO:HG3	1.94	0.50
1:A:385:PHE:O	1:A:388:ARG:HG3	2.12	0.50
1:A:189:THR:CG2	1:A:190:ALA:N	2.75	0.49
1:A:310:ALA:CA	1:A:334:MSE:HE2	2.42	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:160:VAL:HA	1:A:164:LEU:HB2	1.94	0.49
1:A:368:THR:HA	1:A:371:GLU:HG2	1.94	0.49
1:A:270:LEU:O	1:A:274:VAL:HG22	2.13	0.49
1:A:326:LEU:HD12	1:A:327:PRO:CD	2.41	0.49
1:A:406:PRO:HA	1:A:418:ARG:NH1	2.27	0.49
1:A:299:ARG:O	1:A:303:ILE:HG13	2.12	0.48
1:A:154:ASN:HD22	1:A:204:GLY:HA3	1.78	0.48
1:A:49:THR:HG22	1:A:50:GLN:N	2.29	0.48
1:A:244:ASP:HB3	1:A:346:LYS:NZ	2.29	0.48
1:A:140:MSE:HE2	1:A:413:HIS:CA	2.41	0.48
1:A:176:THR:CG2	1:A:178:ALA:H	2.24	0.47
1:A:163:GLN:HE21	1:A:169:PHE:HA	1.80	0.47
1:A:209:GLU:HG3	2:A:452:HOH:O	2.15	0.47
1:A:6:GLN:HG2	1:A:90:PRO:HG3	1.97	0.47
1:A:33:HIS:CE1	1:A:37:ILE:HD12	2.50	0.46
1:A:137:LEU:HD13	1:A:145:ARG:HG3	1.97	0.46
1:A:143:THR:HG23	1:A:144:GLU:H	1.81	0.46
1:A:325:THR:HG22	1:A:326:LEU:N	2.31	0.46
1:A:419:MSE:HE2	1:A:419:MSE:N	2.31	0.46
1:A:191:ARG:HH11	1:A:191:ARG:HG2	1.82	0.45
1:A:109:HIS:CE1	1:A:185:LYS:NZ	2.84	0.45
1:A:189:THR:HG23	2:A:457:HOH:O	2.17	0.45
1:A:247:TYR:O	1:A:251:ASP:HB2	2.17	0.45
1:A:253:GLU:O	1:A:256:LEU:HB2	2.17	0.45
1:A:335:HIS:CD2	1:A:337:PRO:HG2	2.52	0.45
1:A:392:PHE:O	1:A:393:LYS:CB	2.65	0.45
1:A:3:LEU:HB3	1:A:8:LEU:HD11	1.99	0.44
1:A:61:HIS:CD2	1:A:254:ASP:HB3	2.52	0.44
1:A:135:GLY:HA2	1:A:138:ASP:OD1	2.18	0.44
1:A:52:HIS:O	1:A:362:HIS:HD2	2.01	0.44
1:A:349:ALA:O	1:A:353:ILE:HB	2.18	0.44
1:A:215:GLN:O	1:A:219:LYS:HG2	2.18	0.43
1:A:138:ASP:C	1:A:140:MSE:H	2.25	0.43
1:A:249:LEU:HD11	1:A:306:LEU:HD12	2.00	0.43
1:A:114:PRO:HB2	1:A:115:PRO:HD2	2.01	0.42
1:A:109:HIS:HE1	1:A:185:LYS:CE	2.32	0.42
1:A:125:ARG:HD3	1:A:201:HIS:ND1	2.33	0.42
1:A:189:THR:HG22	1:A:190:ALA:N	2.34	0.42
1:A:273:LEU:HD12	1:A:274:VAL:HG13	2.02	0.42
1:A:115:PRO:HB3	1:A:428:ASP:HA	2.01	0.42
1:A:309:ALA:HB1	1:A:334:MSE:HG2	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:86:ARG:HA	1:A:86:ARG:HD2	1.86	0.42
1:A:436:ARG:C	1:A:438:MSE:N	2.71	0.42
1:A:11:ARG:HD3	1:A:92:TRP:CZ3	2.55	0.42
1:A:200:LYS:HD2	1:A:201:HIS:H	1.83	0.42
1:A:406:PRO:HA	1:A:418:ARG:HH12	1.85	0.41
1:A:163:GLN:NE2	1:A:169:PHE:HD2	2.18	0.41
1:A:401:LEU:N	1:A:401:LEU:CD1	2.81	0.41
1:A:80:ARG:HH11	1:A:308:ASN:ND2	2.19	0.41
1:A:98:LEU:O	1:A:101:VAL:CG1	2.68	0.41
1:A:108:ALA:HB1	1:A:161:LEU:HD11	2.03	0.41
1:A:5:TRP:CD1	1:A:322:LEU:HD21	2.56	0.41
1:A:128:PHE:HZ	1:A:420:ILE:HD11	1.86	0.41
1:A:163:GLN:NE2	1:A:169:PHE:HA	2.35	0.41
1:A:436:ARG:O	1:A:438:MSE:N	2.54	0.41
1:A:32:ASP:OD2	1:A:176:THR:HB	2.21	0.41
1:A:244:ASP:HB3	1:A:346:LYS:HZ1	1.86	0.40
1:A:326:LEU:CD1	1:A:327:PRO:HD2	2.47	0.40
1:A:416:PHE:O	1:A:420:ILE:HG12	2.21	0.40
1:A:41:ALA:HB3	1:A:111:ILE:HD11	2.03	0.40
1:A:376:ALA:HB1	1:A:397:ILE:HD13	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	383/451 (85%)	359 (94%)	19 (5%)	5 (1%)	9 8

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	402	GLY

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Mol	Chain	Res	Type
1	A	225	LEU
1	A	437	GLU
1	A	327	PRO
1	A	336	GLY

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	312/359 (87%)	299 (96%)	13 (4%)	26 35

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	114	PRO
1	A	143	THR
1	A	163	GLN
1	A	164	LEU
1	A	176	THR
1	A	200	LYS
1	A	228	GLN
1	A	244	ASP
1	A	262	ASP
1	A	273	LEU
1	A	355	GLN
1	A	370	LEU
1	A	417	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	61	HIS
1	A	109	HIS
1	A	146	ASN
1	A	150	ASN

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Mol	Chain	Res	Type
1	A	154	ASN
1	A	163	GLN
1	A	207	GLN
1	A	308	ASN
1	A	335	HIS
1	A	362	HIS

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	390:PRO	C	391:SER	N	1.09

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	382/451 (84%)	1.02	76 (19%) 3 3	17, 41, 65, 79	0

All (76) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	275	GLY	5.7
1	A	292	ARG	4.9
1	A	263	TYR	4.8
1	A	325	THR	4.6
1	A	262	ASP	4.5
1	A	200	LYS	4.0
1	A	392	PHE	3.9
1	A	191	ARG	3.9
1	A	437	GLU	3.7
1	A	323	ALA	3.7
1	A	2	SER	3.7
1	A	367	TYR	3.6
1	A	312	ARG	3.6
1	A	317	GLN	3.5
1	A	318	GLN	3.5
1	A	5	TRP	3.4
1	A	3	LEU	3.4
1	A	353	ILE	3.4
1	A	322	LEU	3.4
1	A	329	ASP	3.3
1	A	324	GLY	3.2
1	A	4	ASP	3.2
1	A	136	TRP	3.1
1	A	28	PRO	3.1
1	A	226	GLU	3.0
1	A	401	LEU	3.0
1	A	254	ASP	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	402	GLY	3.0
1	A	304	GLU	2.8
1	A	257	GLU	2.8
1	A	261	LEU	2.8
1	A	433	GLU	2.7
1	A	273	LEU	2.7
1	A	227	GLU	2.7
1	A	129	ASN	2.6
1	A	134	ARG	2.6
1	A	268	SER	2.6
1	A	293	ARG	2.6
1	A	327	PRO	2.6
1	A	404	SER	2.5
1	A	320	ALA	2.5
1	A	393	LYS	2.5
1	A	395	ARG	2.5
1	A	331	VAL	2.4
1	A	326	LEU	2.4
1	A	83	GLU	2.4
1	A	130	GLN	2.4
1	A	201	HIS	2.4
1	A	436	ARG	2.4
1	A	143	THR	2.3
1	A	256	LEU	2.3
1	A	89	LEU	2.3
1	A	93	CYS	2.3
1	A	260	LEU	2.3
1	A	400	LEU	2.3
1	A	53	PRO	2.3
1	A	135	GLY	2.3
1	A	350	ARG	2.3
1	A	388	ARG	2.3
1	A	396	ARG	2.3
1	A	387	GLY	2.2
1	A	88	ALA	2.2
1	A	298	LEU	2.2
1	A	141	SER	2.2
1	A	249	LEU	2.1
1	A	6	GLN	2.1
1	A	90	PRO	2.1
1	A	61	HIS	2.1
1	A	335	HIS	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	354	PHE	2.1
1	A	176	THR	2.0
1	A	189	THR	2.0
1	A	391	SER	2.0
1	A	133	GLY	2.0
1	A	328	GLY	2.0
1	A	10	ASN	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.