



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 10, 2016 – 04:27 PM EDT

PDB ID : 5LV9
Title : Crystal structure of thermophilic tryptophan halogenase (Th-Hal) enzyme from *Streptomyces violaceusniger*.
Authors : Dunstan, M.S.; Menon, B.
Deposited on : 2016-09-13
Resolution : 2.33 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	:	unknown
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20027939
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-20027939

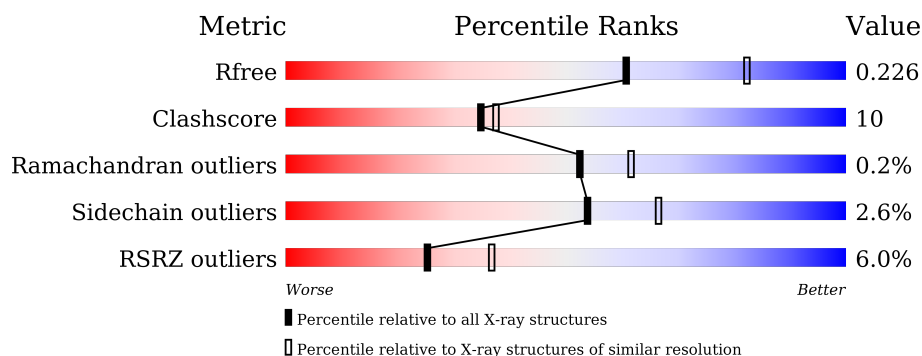
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.33 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	1406 (2.36-2.32)
Clashscore	102246	1509 (2.36-2.32)
Ramachandran outliers	100387	1490 (2.36-2.32)
Sidechain outliers	100360	1491 (2.36-2.32)
RSRZ outliers	91569	1412 (2.36-2.32)

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	513	<div> <div>7%</div> <div> <div></div> <div>77%</div> <div>20%</div> <div>..</div> </div> </div>
1	B	513	<div> <div>4%</div> <div> <div></div> <div>83%</div> <div>16%</div> <div>.</div> </div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8510 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 thermophilic tryptophan halogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	509	Total	C	N	O	S	0	0	0
			4047	2579	696	753	19			
1	B	513	Total	C	N	O	S	0	0	0
			4086	2606	701	760	19			

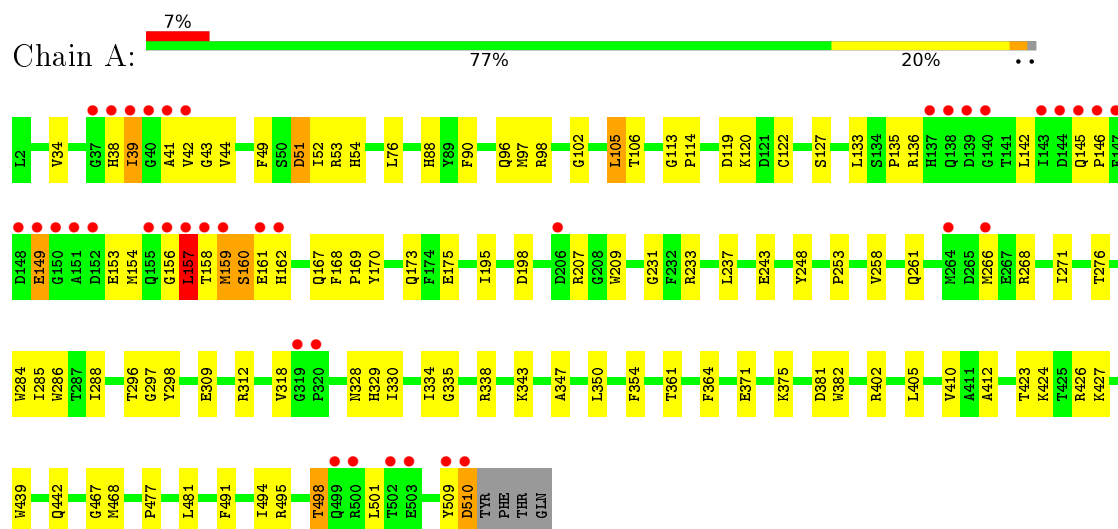
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	170	Total	O	0	0
			170	170		
2	B	207	Total	O	0	0
			207	207		

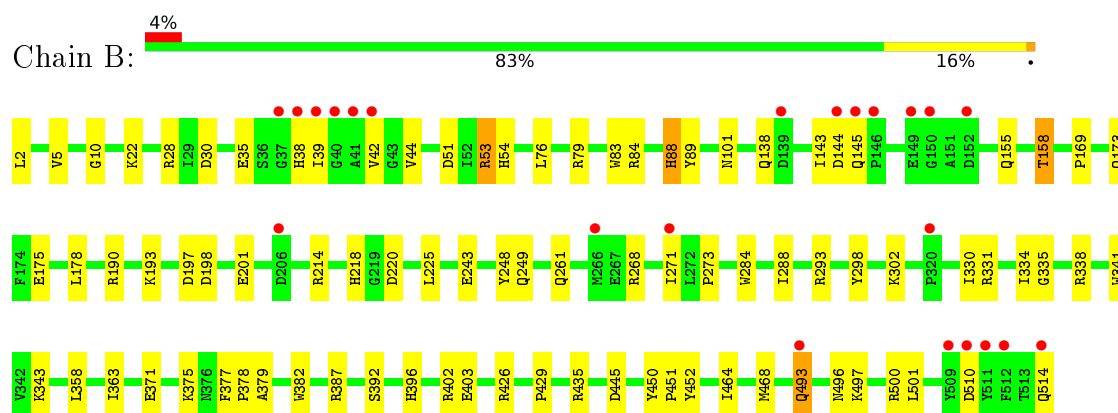
3 Residue-property plots

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: thermophilic tryptophan halogenase



● Molecule 1: thermophilic tryptophan halogenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, α , β , γ	67.14Å 67.14Å 477.37Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	56.49 – 2.33 56.49 – 2.33	Depositor EDS
% Data completeness (in resolution range)	99.7 (56.49-2.33) 99.7 (56.49-2.33)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.65 (at 2.34Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.179 , 0.224 0.179 , 0.226	Depositor DCC
R_{free} test set	2546 reflections (4.95%)	DCC
Wilson B-factor (Å ²)	34.3	Xtriage
Anisotropy	0.011	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 39.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.073 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8510	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.16% 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	0.43	0/4154	0.60	1/5639 (0.0%)
1	B	0.42	0/4195	0.61	1/5695 (0.0%)
All	All	0.43	0/8349	0.61	2/11334 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	157	LEU	CA-CB-CG	10.69	139.89	115.30
1	B	493	GLN	CA-CB-CG	5.84	126.24	113.40

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	4047	0	3915	95	0
1	B	4086	0	3948	67	0
2	A	170	0	0	15	1
2	B	207	0	0	11	1
All	All	8510	0	7863	158	1

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 10.

All (158) 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:198:ASP:OD2	2:A:601:HOH:O	1.82	0.95
1:B:53:ARG:NH1	2:B:601:HOH:O	1.99	0.95
1:A:382:TRP:O	2:A:602:HOH:O	1.89	0.89
1:A:157:LEU:HG	1:A:162:HIS:CE1	2.12	0.84
1:A:97:MET:HE1	1:A:106:THR:H	1.44	0.83
1:B:145:GLN:NE2	1:B:169:PRO:HB3	1.95	0.81
1:A:136:ARG:NH1	1:A:170:TYR:OH	2.13	0.80
1:A:145:GLN:HG2	1:A:146:PRO:HD2	1.65	0.77
1:A:157:LEU:HG	1:A:162:HIS:NE2	2.00	0.76
1:A:39:ILE:HG13	1:A:41:ALA:H	1.50	0.76
1:A:49:PHE:HB2	1:A:51:ASP:OD1	1.84	0.75
1:A:276:THR:OG1	2:A:604:HOH:O	2.05	0.73
1:B:403:GLU:OE1	2:B:603:HOH:O	2.07	0.73
1:B:343:LYS:O	1:B:387:ARG:NH1	2.21	0.72
1:A:381:ASP:OD1	2:A:605:HOH:O	2.07	0.71
1:B:371:GLU:OE2	2:B:602:HOH:O	2.07	0.71
1:B:84:ARG:HD2	1:B:88:HIS:CD2	2.25	0.71
1:A:42:VAL:HG23	1:A:43:GLY:H	1.56	0.71
1:B:83:TRP:CE3	1:B:468:MET:HE1	2.26	0.71
1:A:207:ARG:NH1	2:A:606:HOH:O	2.24	0.69
1:A:154:MET:HB2	1:A:157:LEU:HD12	1.74	0.69
1:A:309:GLU:OE2	1:A:312:ARG:NH2	2.27	0.68
1:A:97:MET:CE	1:A:106:THR:H	2.05	0.68
1:B:218:HIS:O	2:B:604:HOH:O	2.12	0.68
1:B:392:SER:O	1:B:396:HIS:HD2	1.78	0.67
1:A:481:LEU:O	1:B:28:ARG:NH1	2.27	0.67
1:B:201:GLU:OE1	1:B:214:ARG:NH1	2.29	0.66
1:A:105:LEU:HD23	1:A:122:CYS:SG	2.36	0.66
1:B:83:TRP:HE3	1:B:468:MET:HE1	1.61	0.65
1:B:268:ARG:HD3	1:B:271:ILE:HD13	1.77	0.65
1:A:296:THR:HG21	1:A:318:VAL:HG21	1.79	0.64
1:B:249:GLN:OE1	2:B:605:HOH:O	2.15	0.64
1:A:156:GLY:HA3	1:B:101:ASN:HA	1.80	0.64
1:B:248:TYR:HB2	1:B:334:ILE:HG22	1.79	0.64
1:A:410:VAL:O	1:A:424:LYS:NZ	2.31	0.64
1:B:214:ARG:HG2	1:B:220:ASP:OD1	1.98	0.63
1:A:412:ALA:HA	1:A:468:MET:HB3	1.79	0.63
1:A:495:ARG:O	1:A:498:THR:HG22	1.98	0.63
1:A:96:GLN:OE1	1:A:167:GLN:HB2	1.99	0.63
1:B:54:HIS:HD2	2:B:627:HOH:O	1.83	0.62
1:A:54:HIS:CE1	1:A:162:HIS:HD1	2.17	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:97:MET:HE1	1:A:106:THR:N	2.14	0.61
1:B:426:ARG:O	2:B:606:HOH:O	2.16	0.61
1:A:120:LYS:NZ	1:A:127:SER:OG	2.34	0.61
1:A:268:ARG:O	1:A:271:ILE:HG12	2.01	0.60
1:A:136:ARG:NH2	1:A:149:GLU:OE1	2.35	0.60
1:B:30:ASP:OD1	1:B:193:LYS:NZ	2.36	0.59
1:A:233:ARG:NH2	2:A:603:HOH:O	1.99	0.59
1:A:49:PHE:CZ	1:A:361:THR:HG22	2.38	0.59
1:A:53:ARG:NH2	1:A:153:GLU:OE1	2.36	0.58
1:A:161:GLU:OE1	1:A:161:GLU:N	2.37	0.58
1:A:412:ALA:O	1:A:424:LYS:NZ	2.36	0.57
1:A:158:THR:O	1:A:162:HIS:CE1	2.58	0.56
1:B:387:ARG:NH2	2:B:617:HOH:O	2.38	0.56
1:B:138:GLN:HG3	1:B:138:GLN:O	2.05	0.56
1:B:173:GLN:HG2	1:B:288:ILE:HG23	1.90	0.54
1:A:51:ASP:HB2	1:A:159:MET:HG3	1.89	0.54
1:B:341:TRP:CG	1:B:387:ARG:HG3	2.42	0.54
1:B:83:TRP:HB3	1:B:468:MET:CE	2.38	0.54
1:B:83:TRP:HB3	1:B:468:MET:HE3	1.88	0.54
1:A:154:MET:H	1:A:157:LEU:HD13	1.73	0.54
1:A:142:LEU:HD11	1:A:169:PRO:HG2	1.89	0.53
1:B:51:ASP:OD1	2:B:607:HOH:O	2.18	0.53
1:A:76:LEU:HD21	1:A:135:PRO:HG3	1.91	0.53
1:A:158:THR:O	1:A:162:HIS:NE2	2.42	0.53
1:B:84:ARG:CZ	1:B:88:HIS:HE2	2.22	0.52
1:B:2:LEU:HD21	1:B:382:TRP:CE2	2.45	0.51
1:A:258:VAL:HG23	1:A:329:HIS:CD2	2.45	0.51
1:A:402:ARG:HD3	2:A:633:HOH:O	2.10	0.51
1:A:423:THR:OG1	1:A:424:LYS:HE2	2.12	0.50
1:A:102:GLY:O	1:B:375:LYS:NZ	2.45	0.49
1:A:97:MET:HE3	1:A:105:LEU:N	2.27	0.49
1:A:509:TYR:HD2	1:A:510:ASP:OD1	1.95	0.49
1:A:43:GLY:HA3	1:A:286:TRP:HH2	1.77	0.49
1:A:98:ARG:HD2	1:A:494:ILE:HD11	1.94	0.49
1:A:231:GLY:HA2	1:A:350:LEU:HB2	1.95	0.49
1:A:54:HIS:CE1	1:A:162:HIS:ND1	2.81	0.48
1:A:44:VAL:HG23	2:A:609:HOH:O	2.12	0.48
1:B:158:THR:HG23	1:B:371:GLU:OE1	2.14	0.48
1:B:341:TRP:CD2	1:B:387:ARG:HG3	2.49	0.48
1:A:88:HIS:HE1	2:A:756:HOH:O	1.97	0.48
1:B:79:ARG:HD2	1:B:89:TYR:CD1	2.49	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:248:TYR:CD2	1:B:335:GLY:HA2	2.49	0.47
1:B:435:ARG:NH2	1:B:451:PRO:HD3	2.29	0.47
1:B:155:GLN:NE2	2:B:625:HOH:O	2.44	0.47
1:B:145:GLN:HE21	1:B:169:PRO:HB3	1.74	0.47
1:B:44:VAL:HG22	1:B:175:GLU:HG3	1.96	0.47
1:B:396:HIS:HB3	1:B:450:TYR:HE1	1.79	0.47
1:B:331:ARG:HD3	2:B:718:HOH:O	2.14	0.47
1:B:284:TRP:CE2	1:B:358:LEU:HD13	2.50	0.47
1:A:157:LEU:CG	1:A:162:HIS:CE1	2.93	0.47
1:A:154:MET:O	1:A:157:LEU:HB2	2.15	0.46
1:B:302:LYS:HD3	1:B:331:ARG:NH1	2.30	0.46
1:B:261:GLN:NE2	1:B:293:ARG:HD3	2.31	0.46
1:A:173:GLN:HG2	1:A:288:ILE:HG23	1.98	0.46
1:A:53:ARG:HH22	1:A:153:GLU:CD	2.19	0.45
1:B:138:GLN:HE22	1:B:514:GLN:HB3	1.81	0.45
1:A:477:PRO:HB3	1:B:379:ALA:HB2	1.98	0.45
1:A:439:TRP:HA	1:A:442:GLN:O	2.15	0.45
1:A:49:PHE:HZ	1:A:361:THR:HG22	1.80	0.45
1:B:144:ASP:O	1:B:497:LYS:HE2	2.16	0.45
1:B:5:VAL:HG22	1:B:225:LEU:HB3	1.98	0.45
1:B:396:HIS:CE1	1:B:452:TYR:CD2	3.04	0.45
1:A:237:LEU:HD22	1:A:347:ALA:HB2	1.99	0.44
1:A:209:TRP:CD1	1:A:343:LYS:HD3	2.52	0.44
1:A:350:LEU:HD22	1:A:354:PHE:HD1	1.82	0.44
1:B:88:HIS:ND1	1:B:89:TYR:N	2.65	0.44
1:A:154:MET:N	1:A:157:LEU:HD13	2.32	0.44
1:A:361:THR:O	1:A:364:PHE:HB3	2.17	0.44
1:B:22:LYS:HE3	1:B:190:ARG:O	2.17	0.44
1:A:127:SER:HB2	1:A:498:THR:OG1	2.18	0.44
1:A:266:MET:HB3	2:A:634:HOH:O	2.17	0.44
1:A:286:TRP:CE2	1:A:297:GLY:HA3	2.53	0.44
1:B:268:ARG:O	1:B:271:ILE:HG12	2.18	0.44
1:A:285:ILE:HA	1:A:297:GLY:O	2.17	0.44
1:A:268:ARG:NH2	2:A:620:HOH:O	2.50	0.43
1:A:330:ILE:HD12	1:A:330:ILE:HA	1.83	0.43
1:A:248:TYR:HB2	1:A:334:ILE:HG22	2.00	0.43
1:B:268:ARG:HH11	1:B:271:ILE:HG21	1.83	0.43
2:A:708:HOH:O	1:B:158:THR:HG21	2.18	0.43
1:B:42:VAL:HG13	1:B:330:ILE:HG21	1.99	0.43
1:B:496:ASN:O	1:B:500:ARG:HG3	2.19	0.43
1:A:160:SER:HB3	2:A:747:HOH:O	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:375:LYS:NZ	2:A:618:HOH:O	2.49	0.43
1:A:44:VAL:HG22	1:A:175:GLU:HG3	2.00	0.43
1:B:10:GLY:HA3	1:B:35:GLU:OE1	2.19	0.42
1:B:197:ASP:OD1	1:B:198:ASP:N	2.48	0.42
1:B:143:ILE:HD11	1:B:501:LEU:HD13	1.99	0.42
1:A:105:LEU:HD21	1:A:491:PHE:CE1	2.53	0.42
1:A:158:THR:O	1:A:162:HIS:CD2	2.72	0.42
1:A:162:HIS:N	1:A:162:HIS:CD2	2.88	0.42
1:A:113:GLY:HA2	1:A:114:PRO:HD3	1.90	0.42
1:B:248:TYR:CD1	1:B:402:ARG:HD2	2.55	0.42
1:B:158:THR:HG23	1:B:371:GLU:CD	2.41	0.41
1:A:142:LEU:HD21	1:A:169:PRO:HG2	2.02	0.41
1:A:136:ARG:HG2	1:A:142:LEU:HA	2.01	0.41
1:B:330:ILE:HA	1:B:330:ILE:HD12	1.89	0.41
1:A:168:PHE:HA	1:A:169:PRO:HD3	1.79	0.41
1:A:248:TYR:CD2	1:A:335:GLY:HA2	2.56	0.41
1:A:119:ASP:OD1	1:A:119:ASP:N	2.45	0.41
1:B:377:PHE:HA	1:B:378:PRO:HD3	1.88	0.41
1:A:52:ILE:HD12	1:A:52:ILE:HA	1.78	0.41
1:B:464:ILE:O	1:B:468:MET:HG2	2.21	0.41
1:A:328:ASN:HB3	2:A:741:HOH:O	2.21	0.41
1:B:178:LEU:HD23	1:B:178:LEU:HA	1.83	0.41
1:A:243:GLU:OE2	1:A:338:ARG:HB3	2.21	0.40
1:B:243:GLU:OE2	1:B:338:ARG:HB3	2.21	0.40
1:B:402:ARG:NH2	1:B:429:PRO:HG3	2.36	0.40
1:A:90:PHE:CZ	1:A:467:GLY:HA3	2.56	0.40
1:A:90:PHE:HZ	1:A:468:MET:HE3	1.86	0.40
1:A:501:LEU:HA	1:A:501:LEU:HD23	1.86	0.40
1:A:34:VAL:HA	1:A:195:ILE:O	2.22	0.40
1:A:253:PRO:HG3	1:A:426:ARG:CZ	2.50	0.40
1:B:76:LEU:HD13	1:B:273:PRO:HB2	2.03	0.40
1:A:160:SER:N	1:A:371:GLU:OE2	2.45	0.40
1:A:157:LEU:C	1:A:162:HIS:CE1	2.95	0.40
1:A:284:TRP:CZ3	1:A:286:TRP:HB3	2.57	0.40
1:A:405:LEU:HD23	1:A:405:LEU:HA	1.89	0.40

All (1) 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)
2:A:648:HOH:O	2:B:749:HOH:O[6_544]	2.11	0.09

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	507/513 (99%)	488 (96%)	18 (4%)	1 (0%)	52	61
1	B	511/513 (100%)	489 (96%)	21 (4%)	1 (0%)	52	61
All	All	1018/1026 (99%)	977 (96%)	39 (4%)	2 (0%)	52	61

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	38	HIS
1	B	510	ASP

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	424/428 (99%)	411 (97%)	13 (3%)	47	59
1	B	428/428 (100%)	419 (98%)	9 (2%)	61	75
All	All	852/856 (100%)	830 (97%)	22 (3%)	54	66

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	39	ILE
1	A	51	ASP
1	A	105	LEU
1	A	133	LEU

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Mol	Chain	Res	Type
1	A	149	GLU
1	A	157	LEU
1	A	159	MET
1	A	160	SER
1	A	261	GLN
1	A	298	TYR
1	A	427	LYS
1	A	498	THR
1	A	510	ASP
1	B	38	HIS
1	B	39	ILE
1	B	53	ARG
1	B	88	HIS
1	B	158	THR
1	B	298	TYR
1	B	363	ILE
1	B	445	ASP
1	B	493	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	145	GLN
1	B	396	HIS
1	B	514	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

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	509/513 (99%)	0.14	38 (7%)	17 26	25, 40, 74, 117	0
1	B	513/513 (100%)	0.02	23 (4%)	37 50	22, 37, 71, 106	0
All	All	1022/1026 (99%)	0.08	61 (5%)	25 37	22, 39, 73, 117	0

All (61) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	40	GLY	12.6
1	A	39	ILE	9.8
1	A	41	ALA	9.4
1	B	39	ILE	8.7
1	A	38	HIS	7.6
1	A	157	LEU	6.5
1	A	162	HIS	6.4
1	B	38	HIS	6.4
1	A	152	ASP	6.1
1	B	41	ALA	5.9
1	A	42	VAL	5.9
1	A	150	GLY	5.4
1	B	40	GLY	5.3
1	B	514	GLN	5.1
1	B	42	VAL	4.7
1	A	509	TYR	4.4
1	B	150	GLY	4.3
1	A	146	PRO	4.2
1	A	37	GLY	4.1
1	A	147	PHE	4.1
1	A	158	THR	4.0
1	A	144	ASP	4.0
1	B	512	PHE	3.9
1	B	152	ASP	3.7

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Mol	Chain	Res	Type	RSRZ
1	A	510	ASP	3.7
1	A	155	GLN	3.6
1	A	499	GLN	3.5
1	B	145	GLN	3.5
1	A	156	GLY	3.4
1	A	151	ALA	3.2
1	A	138	GLN	3.2
1	A	137	HIS	3.1
1	A	266	MET	3.1
1	B	266	MET	3.1
1	B	511	TYR	3.0
1	A	148	ASP	3.0
1	B	37	GLY	3.0
1	B	146	PRO	3.0
1	A	149	GLU	2.9
1	B	510	ASP	2.9
1	A	159	MET	2.8
1	B	320	PRO	2.8
1	A	500	ARG	2.7
1	A	503	GLU	2.7
1	B	271	ILE	2.7
1	B	144	ASP	2.6
1	B	493	GLN	2.6
1	B	206	ASP	2.5
1	A	161	GLU	2.5
1	A	145	GLN	2.4
1	A	319	GLY	2.4
1	A	140	GLY	2.3
1	A	139	ASP	2.2
1	B	139	ASP	2.2
1	A	143	ILE	2.2
1	A	320	PRO	2.2
1	B	509	TYR	2.1
1	A	206	ASP	2.1
1	A	264	MET	2.0
1	A	502	THR	2.0
1	B	149	GLU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.