



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 22, 2026 – 01:24 AM UTC

PDB ID : 1AI4 / pdb_00001ai4
Title : PENICILLIN ACYLASE COMPLEXED WITH 3,4-DIHYDROXYPHENYL ACETIC ACID
Authors : Done, S.H.
Deposited on : 1997-05-01
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

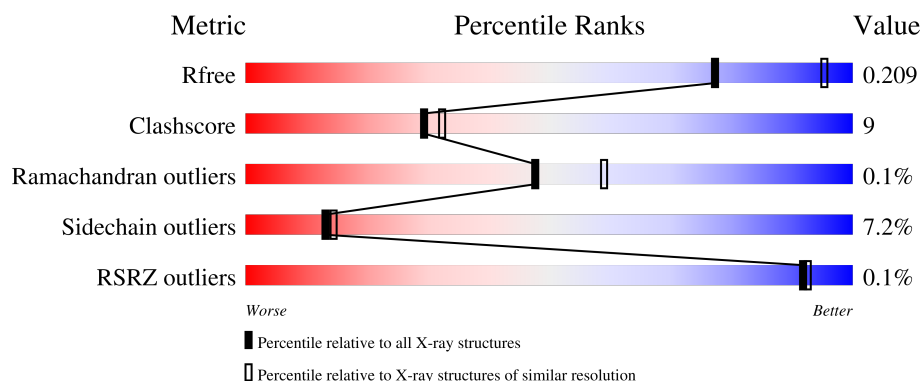
1 Overall quality at a glance

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	209	
2	B	557	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6684 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 PENICILLIN AMIDOHYDROLASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	206	Total	C	N	O	S	0	0	0
			1656	1058	278	312	8			

- Molecule 2 is a protein called PENICILLIN AMIDOHYDROLASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	557	Total	C	N	O	S	0	0	0
			4415	2805	767	833	10			

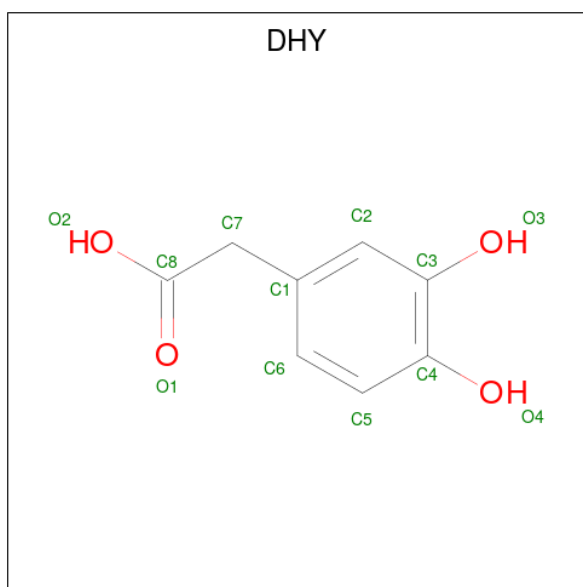
There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	165	GLN	GLU	conflict	UNP P06875

- Molecule 3 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Ca	0	0
			1	1		

- Molecule 4 is 2-(3,4-DIHYDROXYPHENYL)ACETIC ACID (CCD ID: DHY) (formula: C₈H₈O₄).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			12	8	4		

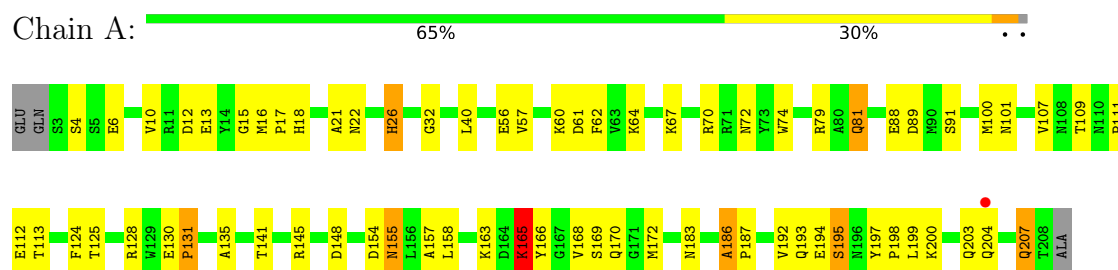
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	153	Total	O	0	0
			153	153		
5	B	447	Total	O	0	0
			447	447		

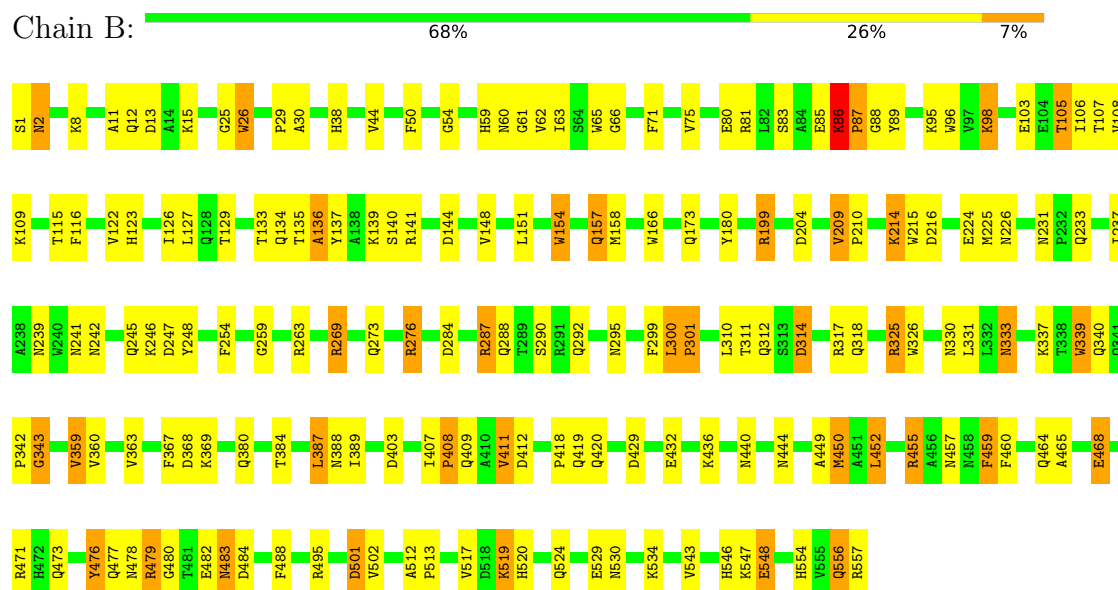
3 Residue-property plots

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: PENICILLIN AMIDOHYDROLASE



• Molecule 2: PENICILLIN AMIDOHYDROLASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	52.12Å 65.08Å 76.30Å 100.20° 111.44° 105.81°	Depositor
Resolution (Å)	19.85 – 2.35 19.85 – 2.36	Depositor EDS
% Data completeness (in resolution range)	96.5 (19.85-2.35) 95.6 (19.85-2.36)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.93 (at 2.35Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.150 , 0.236 0.138 , 0.209	Depositor DCC
R_{free} test set	2413 reflections (7.16%)	wwPDB-VP
Wilson B-factor (Å ²)	21.7	Xtriage
Anisotropy	0.279	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 91.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	6684	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.37% 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: DHY, CA

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	0.83	1/1698 (0.1%)	1.90	44/2305 (1.9%)
2	B	0.82	1/4541 (0.0%)	1.95	137/6192 (2.2%)
All	All	0.83	2/6239 (0.0%)	1.94	181/8497 (2.1%)

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
2	B	0	3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	412	ASP	C-N	-6.23	1.24	1.33
1	A	125	THR	N-CA	5.54	1.50	1.45

All (181) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	157	GLN	CG-CD-NE2	-15.25	93.52	116.40
1	A	79	ARG	CD-NE-CZ	13.88	143.83	124.40
2	B	157	GLN	CG-CD-OE1	12.49	145.77	120.80
2	B	363	VAL	CA-C-O	9.78	125.55	119.19
2	B	29	PRO	CA-C-N	8.44	132.80	120.90
2	B	29	PRO	C-N-CA	8.44	132.80	120.90
2	B	144	ASP	CA-CB-CG	8.37	120.97	112.60
2	B	464	GLN	OE1-CD-NE2	-8.37	114.23	122.60
1	A	22	ASN	CA-CB-CG	-8.29	104.31	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	330	ASN	CB-CG-ND2	8.24	128.75	116.40
2	B	157	GLN	CB-CG-CD	8.20	126.53	112.60
2	B	460	PHE	CA-CB-CG	8.09	121.89	113.80
2	B	483	ASN	OD1-CG-ND2	-8.00	114.60	122.60
2	B	60	ASN	OD1-CG-ND2	-7.79	114.81	122.60
2	B	412	ASP	O-C-N	-7.72	113.53	123.02
2	B	269	ARG	NE-CZ-NH1	7.56	129.06	121.50
2	B	409	GLN	CG-CD-NE2	-7.40	105.30	116.40
2	B	239	ASN	OD1-CG-ND2	-7.36	115.24	122.60
2	B	263	ARG	NE-CZ-NH2	-7.31	112.62	119.20
2	B	488	PHE	CA-C-N	7.29	130.71	122.83
2	B	488	PHE	C-N-CA	7.29	130.71	122.83
1	A	89	ASP	CA-CB-CG	-7.29	105.31	112.60
2	B	330	ASN	OD1-CG-ND2	-7.28	115.32	122.60
2	B	408	PRO	N-CA-CB	7.28	109.12	103.30
1	A	61	ASP	CA-CB-CG	7.27	119.87	112.60
2	B	173	GLN	CG-CD-OE1	7.25	135.30	120.80
2	B	444	ASN	CA-CB-CG	7.13	119.73	112.60
2	B	287	ARG	CD-NE-CZ	7.09	134.32	124.40
1	A	183	ASN	CA-C-N	7.03	126.73	119.56
1	A	183	ASN	C-N-CA	7.03	126.73	119.56
2	B	483	ASN	CB-CG-ND2	6.92	126.79	116.40
1	A	154	ASP	CA-C-O	-6.91	113.09	120.42
2	B	343	GLY	N-CA-C	6.71	120.76	112.64
1	A	16	MET	CA-C-N	6.69	126.92	119.90
1	A	16	MET	C-N-CA	6.69	126.92	119.90
2	B	157	GLN	CA-CB-CG	6.66	127.42	114.10
2	B	409	GLN	CB-CG-CD	6.62	123.85	112.60
2	B	339	TRP	CA-C-N	6.57	129.08	120.28
2	B	339	TRP	C-N-CA	6.57	129.08	120.28
2	B	548	GLU	N-CA-CB	6.53	119.74	110.67
2	B	331	LEU	CA-C-O	-6.53	112.77	120.60
2	B	173	GLN	CG-CD-NE2	-6.52	106.62	116.40
2	B	245	GLN	N-CA-CB	-6.48	101.71	111.56
2	B	411	VAL	N-CA-CB	6.47	122.65	111.39
1	A	26	HIS	CB-CG-CD2	-6.45	122.81	131.20
2	B	389	ILE	CA-C-O	-6.45	113.13	120.74
2	B	301	PRO	CA-C-N	6.34	128.78	120.28
2	B	301	PRO	C-N-CA	6.34	128.78	120.28
2	B	269	ARG	CA-C-O	-6.33	113.72	120.42
2	B	311	THR	CA-C-N	6.31	129.36	120.28
2	B	311	THR	C-N-CA	6.31	129.36	120.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	74	TRP	CA-C-N	6.30	126.50	119.32
1	A	74	TRP	C-N-CA	6.30	126.50	119.32
1	A	135	ALA	CA-C-O	-6.29	113.88	120.55
2	B	482	GLU	CA-C-O	-6.28	114.31	121.40
2	B	25	GLY	N-CA-C	-6.22	102.86	111.76
2	B	66	GLY	CA-C-O	-6.22	114.91	121.56
1	A	124	PHE	CA-CB-CG	6.20	120.00	113.80
2	B	333	ASN	OD1-CG-ND2	6.16	128.76	122.60
2	B	54	GLY	O-C-N	-6.13	116.70	122.96
2	B	412	ASP	CA-CB-CG	6.13	118.73	112.60
1	A	61	ASP	CA-C-O	-6.13	112.08	119.49
1	A	141	THR	O-C-N	-6.13	114.44	122.59
2	B	136	ALA	CA-C-O	-6.13	114.08	120.70
2	B	233	GLN	OE1-CD-NE2	-6.08	116.52	122.60
1	A	15	GLY	N-CA-C	-6.05	106.97	115.32
2	B	465	ALA	CA-C-O	-6.02	114.84	121.28
1	A	81	GLN	OE1-CD-NE2	-5.99	116.61	122.60
1	A	157	ALA	CA-C-O	-5.98	114.08	120.42
2	B	449	ALA	CA-C-N	5.98	129.97	121.42
2	B	449	ALA	C-N-CA	5.98	129.97	121.42
1	A	40	LEU	CA-C-O	-5.97	114.56	120.82
2	B	359	VAL	CB-CA-C	-5.96	104.33	111.97
1	A	186	ALA	CA-C-N	5.96	126.26	119.83
1	A	186	ALA	C-N-CA	5.96	126.26	119.83
2	B	30	ALA	O-C-N	5.95	130.22	122.97
2	B	1	SER	CA-C-N	-5.93	112.86	122.82
2	B	1	SER	C-N-CA	-5.93	112.86	122.82
2	B	314	ASP	CA-C-N	5.92	125.80	119.28
2	B	314	ASP	C-N-CA	5.92	125.80	119.28
2	B	548	GLU	CA-CB-CG	5.91	125.92	114.10
2	B	237	ILE	CA-C-O	-5.87	114.34	120.27
2	B	471	ARG	CD-NE-CZ	5.85	132.59	124.40
1	A	141	THR	CA-C-N	5.83	128.02	120.44
1	A	141	THR	C-N-CA	5.83	128.02	120.44
1	A	165	LYS	CA-C-O	-5.83	114.88	121.00
2	B	50	PHE	CA-CB-CG	5.81	119.61	113.80
2	B	75	VAL	CA-C-O	-5.80	114.41	120.27
2	B	534	LYS	CA-C-N	5.80	128.95	120.71
2	B	534	LYS	C-N-CA	5.80	128.95	120.71
2	B	105	THR	CA-C-O	-5.78	114.12	120.36
2	B	154	TRP	CA-C-O	5.76	126.59	119.97
1	A	111	PRO	N-CA-CB	5.68	108.96	103.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	478	ASN	OD1-CG-ND2	-5.67	116.92	122.60
2	B	295	ASN	CA-CB-CG	-5.67	106.93	112.60
1	A	192	VAL	CB-CA-C	-5.62	103.45	112.16
2	B	86	LYS	CA-C-O	5.60	127.83	120.16
2	B	88	GLY	N-CA-C	-5.60	107.37	115.27
2	B	524	GLN	CA-C-N	5.60	127.78	120.28
2	B	524	GLN	C-N-CA	5.60	127.78	120.28
2	B	547	LYS	CA-C-O	-5.58	114.56	120.92
2	B	263	ARG	N-CA-C	5.55	119.15	112.38
1	A	158	LEU	CA-C-N	5.55	127.72	120.28
1	A	158	LEU	C-N-CA	5.55	127.72	120.28
2	B	476	TYR	N-CA-CB	5.54	118.16	109.69
2	B	126	ILE	CA-C-N	5.53	127.96	120.44
2	B	126	ILE	C-N-CA	5.53	127.96	120.44
2	B	241	ASN	CA-CB-CG	5.51	118.11	112.60
2	B	292	GLN	N-CA-C	5.49	117.66	109.59
2	B	248	TYR	CA-C-O	5.49	124.64	119.76
2	B	342	PRO	CA-C-N	5.48	126.18	119.99
2	B	342	PRO	C-N-CA	5.48	126.18	119.99
1	A	81	GLN	CG-CD-NE2	5.47	124.61	116.40
2	B	388	ASN	CB-CG-ND2	-5.47	108.20	116.40
1	A	32	GLY	CA-C-O	5.45	126.29	120.40
2	B	86	LYS	CA-CB-CG	5.44	124.97	114.10
2	B	452	LEU	CA-C-O	5.42	125.98	120.24
1	A	17	PRO	N-CA-CB	5.41	108.12	103.36
2	B	199	ARG	CD-NE-CZ	-5.41	116.83	124.40
2	B	429	ASP	CA-CB-CG	5.40	118.00	112.60
2	B	455	ARG	N-CA-C	5.40	118.47	110.48
2	B	457	ASN	N-CA-CB	5.38	119.18	110.41
2	B	86	LYS	CA-C-N	5.38	125.14	119.76
2	B	86	LYS	C-N-CA	5.38	125.14	119.76
2	B	290	SER	CB-CA-C	-5.37	101.55	110.68
1	A	26	HIS	CA-CB-CG	-5.36	108.44	113.80
2	B	87	PRO	N-CA-CB	5.36	108.32	103.34
2	B	15	LYS	CA-C-O	-5.36	113.24	118.97
2	B	259	GLY	O-C-N	-5.34	119.84	123.95
2	B	209	VAL	CA-C-N	5.32	125.24	119.76
2	B	209	VAL	C-N-CA	5.32	125.24	119.76
2	B	284	ASP	CA-CB-CG	-5.31	107.29	112.60
1	A	195	SER	CA-C-O	-5.31	115.60	121.28
2	B	554	HIS	CB-CA-C	5.29	118.61	110.62
2	B	154	TRP	O-C-N	-5.28	115.78	122.27

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	455	ARG	NE-CZ-NH2	-5.26	114.46	119.20
2	B	231	ASN	CA-C-N	5.24	125.24	119.89
2	B	231	ASN	C-N-CA	5.24	125.24	119.89
2	B	318	GLN	CB-CG-CD	5.24	121.52	112.60
2	B	226	ASN	CA-CB-CG	5.24	117.84	112.60
2	B	242	ASN	OD1-CG-ND2	-5.24	117.36	122.60
2	B	367	PHE	CA-CB-CG	-5.24	108.56	113.80
2	B	62	VAL	N-CA-C	-5.23	107.72	112.12
2	B	495	ARG	CA-C-N	5.21	124.83	119.56
2	B	495	ARG	C-N-CA	5.21	124.83	119.56
2	B	263	ARG	NE-CZ-NH1	5.21	126.71	121.50
1	A	79	ARG	NE-CZ-NH2	5.21	123.89	119.20
2	B	137	TYR	CA-C-O	-5.17	114.79	120.43
2	B	479	ARG	NE-CZ-NH1	5.17	126.67	121.50
1	A	70	ARG	CA-C-O	-5.17	115.07	120.55
2	B	363	VAL	O-C-N	-5.12	116.95	121.61
2	B	440	ASN	OD1-CG-ND2	-5.12	117.47	122.60
2	B	215	TRP	CA-CB-CG	-5.12	103.87	113.60
2	B	312	GLN	CA-CB-CG	5.11	124.32	114.10
2	B	325	ARG	N-CA-CB	-5.11	101.82	110.41
2	B	276	ARG	CA-C-O	-5.11	115.47	121.44
1	A	32	GLY	CA-C-N	5.10	127.38	120.44
1	A	32	GLY	C-N-CA	5.10	127.38	120.44
1	A	100	MET	CA-C-O	5.09	125.82	120.42
2	B	71	PHE	CA-CB-CG	-5.09	108.71	113.80
2	B	455	ARG	CD-NE-CZ	5.09	131.52	124.40
2	B	418	PRO	N-CA-CB	5.08	107.83	103.31
2	B	300	LEU	CA-C-N	5.08	124.52	119.24
2	B	300	LEU	C-N-CA	5.08	124.52	119.24
2	B	301	PRO	O-C-N	-5.08	116.70	122.18
1	A	155	ASN	CB-CG-ND2	5.07	124.01	116.40
2	B	455	ARG	NH1-CZ-NH2	5.07	125.89	119.30
2	B	501	ASP	CA-CB-CG	5.06	117.66	112.60
2	B	15	LYS	O-C-N	5.06	128.60	122.48
2	B	50	PHE	O-C-N	-5.05	114.10	122.23
2	B	449	ALA	O-C-N	-5.04	117.26	122.96
2	B	269	ARG	CD-NE-CZ	5.04	131.46	124.40
2	B	546	HIS	CA-CB-CG	-5.04	108.76	113.80
1	A	131	PRO	CA-C-N	5.04	127.03	120.28
1	A	131	PRO	C-N-CA	5.04	127.03	120.28
1	A	62	PHE	CA-CB-CG	-5.04	108.77	113.80
1	A	57	VAL	CA-C-O	-5.02	114.50	120.78

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	484	ASP	O-C-N	5.02	129.00	123.33
2	B	480	GLY	CA-C-N	5.01	127.50	120.28
2	B	480	GLY	C-N-CA	5.01	127.50	120.28
2	B	141	ARG	CD-NE-CZ	5.00	131.41	124.40

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	122	VAL	Mainchain
2	B	26	TRP	Mainchain
2	B	299	PHE	Mainchain

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	1656	0	1604	34	0
2	B	4415	0	4244	86	0
3	B	1	0	0	0	0
4	B	12	0	7	0	0
5	A	153	0	0	0	0
5	B	447	0	0	1	0
All	All	6684	0	5855	103	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 9.

All (103) 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:207:GLN:HE22	2:B:204:ASP:H	1.03	1.01
2:B:59:HIS:HD2	2:B:61:GLY:H	1.23	0.85
2:B:157:GLN:OE1	2:B:166:TRP:NE1	2.12	0.82
2:B:59:HIS:CD2	2:B:61:GLY:H	1.99	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:387:LEU:HD21	2:B:450:MET:HE1	1.62	0.80
1:A:207:GLN:NE2	2:B:204:ASP:H	1.79	0.80
2:B:556:GLN:NE2	2:B:557:ARG:H	1.85	0.73
2:B:86:LYS:N	2:B:87:PRO:HD3	2.05	0.72
1:A:18:HIS:HD2	2:B:38:HIS:NE2	1.88	0.71
2:B:214:LYS:H	2:B:214:LYS:HD2	1.54	0.70
2:B:339:TRP:HH2	2:B:450:MET:HE2	1.57	0.69
2:B:512:ALA:HB1	2:B:513:PRO:HD2	1.75	0.69
2:B:83:SER:HB2	2:B:96:TRP:CH2	2.29	0.68
2:B:86:LYS:H	2:B:87:PRO:HD3	1.59	0.68
1:A:207:GLN:HE22	2:B:204:ASP:N	1.87	0.68
2:B:479:ARG:HH21	2:B:483:ASN:HD22	1.42	0.68
1:A:72:ASN:HD21	2:B:139:LYS:NZ	1.92	0.67
2:B:479:ARG:HH21	2:B:483:ASN:ND2	1.93	0.65
2:B:556:GLN:NE2	2:B:557:ARG:N	2.43	0.65
2:B:360:VAL:HG13	2:B:368:ASP:HB2	1.79	0.64
2:B:129:THR:HG22	2:B:136:ALA:CB	2.28	0.63
2:B:12:GLN:O	2:B:13:ASP:HB2	1.98	0.61
2:B:384:THR:HG22	2:B:455:ARG:NH1	2.16	0.60
1:A:67:LYS:HG2	2:B:116:PHE:CE1	2.36	0.60
2:B:26:TRP:CD2	2:B:452:LEU:HD11	2.37	0.58
2:B:86:LYS:N	2:B:87:PRO:CD	2.66	0.58
2:B:44:VAL:HG11	2:B:158:MET:HB3	1.85	0.57
2:B:317:ARG:HG3	2:B:317:ARG:HH11	1.70	0.57
2:B:519:LYS:HG2	2:B:520:HIS:CD2	2.39	0.57
2:B:123:HIS:O	2:B:140:SER:HB2	2.06	0.56
2:B:129:THR:HG22	2:B:136:ALA:HA	1.88	0.56
2:B:214:LYS:H	2:B:214:LYS:CD	2.18	0.56
2:B:89:TYR:CZ	2:B:98:LYS:HG3	2.42	0.55
2:B:384:THR:HG22	2:B:455:ARG:HH12	1.72	0.55
2:B:459:PHE:C	2:B:459:PHE:CD1	2.84	0.55
1:A:165:LYS:HD3	1:A:166:TYR:CE2	2.42	0.54
1:A:56:GLU:O	2:B:109:LYS:HB2	2.08	0.54
2:B:384:THR:HG22	2:B:455:ARG:CZ	2.39	0.53
1:A:155:ASN:ND2	2:B:254:PHE:HB3	2.24	0.53
2:B:512:ALA:HB1	2:B:513:PRO:CD	2.38	0.52
1:A:72:ASN:HD21	2:B:139:LYS:HZ1	1.58	0.52
1:A:101:ASN:HB3	1:A:128:ARG:HH12	1.75	0.52
2:B:333:ASN:ND2	2:B:340:GLN:HG2	2.25	0.52
2:B:11:ALA:O	2:B:276:ARG:NH1	2.34	0.51
1:A:101:ASN:HB3	1:A:128:ARG:NH1	2.25	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:128:ARG:HH11	1:A:128:ARG:HG2	1.75	0.51
1:A:145:ARG:HG3	2:B:459:PHE:HZ	1.74	0.51
1:A:155:ASN:HD21	2:B:254:PHE:HB3	1.74	0.51
2:B:209:VAL:HB	2:B:210:PRO:HD2	1.93	0.51
1:A:21:ALA:HB1	1:A:26:HIS:HB3	1.93	0.51
2:B:65:TRP:HA	2:B:180:TYR:O	2.12	0.50
2:B:556:GLN:HE21	2:B:557:ARG:N	2.10	0.50
1:A:128:ARG:NH1	1:A:128:ARG:HG2	2.26	0.50
1:A:81:GLN:NE2	2:B:148:VAL:H	2.11	0.49
1:A:163:LYS:HG2	1:A:168:VAL:HA	1.95	0.49
2:B:288:GLN:HG2	5:B:980:HOH:O	2.13	0.49
1:A:166:TYR:O	1:A:170:GLN:HB3	2.13	0.48
1:A:130:GLU:HB2	1:A:131:PRO:HD2	1.95	0.48
2:B:123:HIS:NE2	2:B:216:ASP:OD1	2.46	0.48
2:B:384:THR:HG22	2:B:455:ARG:NH2	2.29	0.48
2:B:384:THR:CG2	2:B:455:ARG:HH12	2.25	0.48
2:B:59:HIS:HB2	2:B:63:ILE:O	2.14	0.48
1:A:10:VAL:HG11	2:B:543:VAL:HG12	1.94	0.47
2:B:310:LEU:HD13	2:B:314:ASP:OD2	2.14	0.47
2:B:317:ARG:HG3	2:B:317:ARG:NH1	2.29	0.47
2:B:326:TRP:CZ3	2:B:343:GLY:HA3	2.50	0.47
2:B:473:GLN:HE22	2:B:477:GLN:NE2	2.13	0.47
2:B:325:ARG:HH11	2:B:325:ARG:HD2	1.59	0.47
2:B:246:LYS:O	2:B:247:ASP:HB2	2.14	0.46
1:A:197:TYR:CD1	1:A:198:PRO:HD2	2.50	0.46
1:A:199:LEU:HG	2:B:225:MET:HE1	1.98	0.46
2:B:387:LEU:HD22	2:B:476:TYR:CE2	2.50	0.46
2:B:339:TRP:CH2	2:B:450:MET:HE2	2.44	0.46
2:B:83:SER:HB2	2:B:96:TRP:CZ3	2.51	0.45
2:B:199:ARG:HH11	2:B:199:ARG:HD2	1.63	0.45
2:B:129:THR:HG22	2:B:136:ALA:CA	2.47	0.45
1:A:12:ASP:O	2:B:548:GLU:HG2	2.17	0.45
2:B:108:VAL:O	2:B:109:LYS:C	2.60	0.44
2:B:151:LEU:HD23	2:B:151:LEU:C	2.43	0.44
2:B:80:GLU:O	2:B:135:THR:HA	2.18	0.44
2:B:106:ILE:HD11	2:B:116:PHE:HE1	1.83	0.44
2:B:287:ARG:HG2	2:B:287:ARG:HH11	1.83	0.44
1:A:172:MET:HE2	1:A:204:GLN:HA	1.99	0.43
1:A:148:ASP:OD2	2:B:139:LYS:NZ	2.52	0.43
1:A:130:GLU:HB2	1:A:131:PRO:CD	2.49	0.43
1:A:186:ALA:HA	1:A:187:PRO:HD3	1.89	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:194:GLU:O	1:A:195:SER:HB3	2.18	0.41
2:B:407:ILE:HA	2:B:408:PRO:HD3	1.89	0.41
2:B:129:THR:HG22	2:B:136:ALA:HB2	2.02	0.41
1:A:88:GLU:O	1:A:91:SER:HB2	2.20	0.41
1:A:56:GLU:HG2	2:B:107:THR:HB	2.03	0.41
2:B:133:THR:O	2:B:134:GLN:HB2	2.19	0.41
2:B:483:ASN:O	2:B:501:ASP:HA	2.20	0.41
2:B:529:GLU:HG2	2:B:530:ASN:OD1	2.20	0.41
2:B:468:GLU:CD	2:B:468:GLU:H	2.28	0.41
1:A:197:TYR:HA	1:A:198:PRO:HD3	1.92	0.40
2:B:209:VAL:HB	2:B:210:PRO:CD	2.50	0.40
2:B:300:LEU:N	2:B:301:PRO:CD	2.84	0.40
2:B:479:ARG:HE	2:B:483:ASN:ND2	2.19	0.40
2:B:479:ARG:NH2	2:B:483:ASN:HD22	2.15	0.40
2:B:2:ASN:C	2:B:2:ASN:HD22	2.29	0.40
1:A:172:MET:HE2	1:A:203:GLN:O	2.21	0.40
2:B:105:THR:HG23	2:B:115:THR:HG22	2.03	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	204/209 (98%)	196 (96%)	8 (4%)	0	100	100
2	B	555/557 (100%)	537 (97%)	17 (3%)	1 (0%)	43	52
All	All	759/766 (99%)	733 (97%)	25 (3%)	1 (0%)	48	59

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	86	LYS

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	178/180 (99%)	164 (92%)	14 (8%)	11	12
2	B	460/460 (100%)	428 (93%)	32 (7%)	14	15
All	All	638/640 (100%)	592 (93%)	46 (7%)	13	14

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	SER
1	A	6	GLU
1	A	13	GLU
1	A	60	LYS
1	A	64	LYS
1	A	107	VAL
1	A	109	THR
1	A	112	GLU
1	A	113	THR
1	A	165	LYS
1	A	169	SER
1	A	193	GLN
1	A	200	LYS
1	A	207	GLN
2	B	2	ASN
2	B	8	LYS
2	B	81	ARG
2	B	85	GLU
2	B	86	LYS
2	B	95	LYS
2	B	98	LYS
2	B	103	GLU
2	B	127	LEU
2	B	154	TRP
2	B	214	LYS
2	B	224	GLU
2	B	269	ARG

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Mol	Chain	Res	Type
2	B	273	GLN
2	B	337	LYS
2	B	359	VAL
2	B	369	LYS
2	B	380	GLN
2	B	387	LEU
2	B	403	ASP
2	B	411	VAL
2	B	419	GLN
2	B	420	GLN
2	B	432	GLU
2	B	436	LYS
2	B	450	MET
2	B	459	PHE
2	B	468	GLU
2	B	502	VAL
2	B	517	VAL
2	B	519	LYS
2	B	556	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	18	HIS
1	A	72	ASN
1	A	81	GLN
1	A	120	ASN
1	A	155	ASN
1	A	207	GLN
2	B	2	ASN
2	B	59	HIS
2	B	110	ASN
2	B	241	ASN
2	B	245	GLN
2	B	304	GLN
2	B	318	GLN
2	B	348	ASN
2	B	419	GLN
2	B	440	ASN
2	B	472	HIS
2	B	477	GLN
2	B	483	ASN

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Mol	Chain	Res	Type
2	B	507	GLN
2	B	556	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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 2 ligands modelled in this entry, 1 is 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	DHY	B	559	-	12,12,12	1.01	1 (8%)	16,16,16	2.56	8 (50%)

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
4	DHY	B	559	-	-	2/4/4/4	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	559	DHY	O2-C8	-2.16	1.23	1.30

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
4	B	559	DHY	O3-C3-C2	4.05	130.38	119.45
4	B	559	DHY	O2-C8-O1	-3.96	113.15	123.33
4	B	559	DHY	O2-C8-C7	3.69	128.04	113.98
4	B	559	DHY	O4-C4-C5	3.38	128.43	119.36
4	B	559	DHY	C7-C1-C2	-3.37	115.27	120.38
4	B	559	DHY	O3-C3-C4	-3.06	110.25	118.42
4	B	559	DHY	O4-C4-C3	-2.97	110.49	118.42
4	B	559	DHY	C7-C1-C6	2.88	125.13	120.89

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	559	DHY	C1-C7-C8-O1
4	B	559	DHY	C1-C7-C8-O2

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 [i](#)

There are no chain breaks in this entry.

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	206/209 (98%)	-0.57	1 (0%) 87 90	14, 26, 59, 80	0
2	B	557/557 (100%)	-0.66	0 100 100	9, 24, 56, 95	0
All	All	763/766 (99%)	-0.64	1 (0%) 92 92	9, 25, 57, 95	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	204	GLN	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	DHY	B	559	12/12	0.91	0.12	27,34,37,37	0
3	CA	B	558	1/1	1.00	0.02	16,16,16,16	0

6.5 Other polymers ⓘ

There are no such residues in this entry.