



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2026 – 07:41 PM UTC

PDB ID : 1IVV / pdb_00001ivv
Title : Crystal structure of copper amine oxidase from *Arthrobacter globiformis*:
Early intermediate in topaquinone biogenesis
Authors : Kim, M.; Okajima, T.; Kishishita, S.; Yoshimura, M.; Kawamori, A.;
Tanizawa, K.; Yamaguchi, H.
Deposited on : 2002-03-29
Resolution : 2.10 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

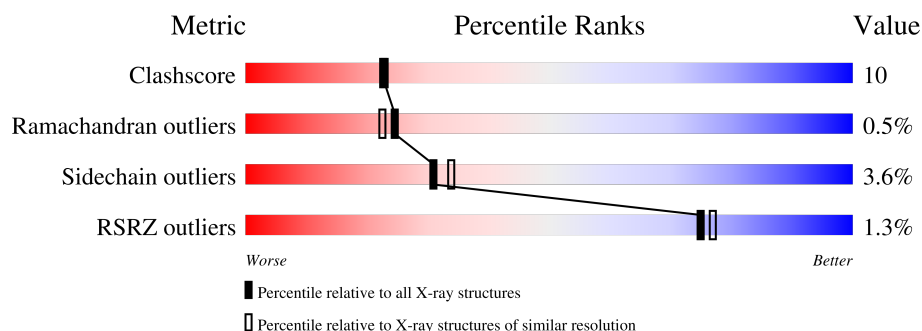
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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))
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	638	% 70% 24% • •
1	B	638	% 77% 18% • •

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10959 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 Phenylethylamine oxidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	620	Total	C	N	O	S	0	1	0
			4872	3077	857	929	9			
1	B	620	Total	C	N	O	S	0	1	0
			4872	3077	857	929	9			

- Molecule 2 is COPPER (II) ION (CCD ID: CU) (formula: Cu).

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

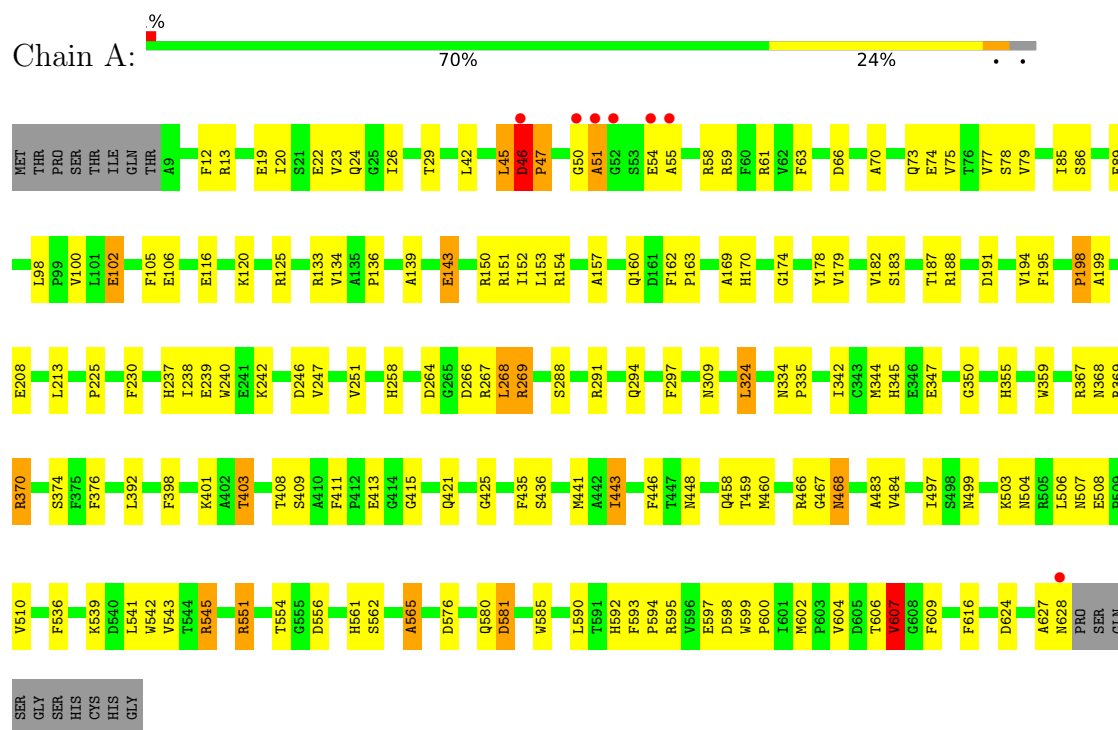
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	579	Total	O	0	0
			579	579		
3	B	634	Total	O	0	0
			634	634		

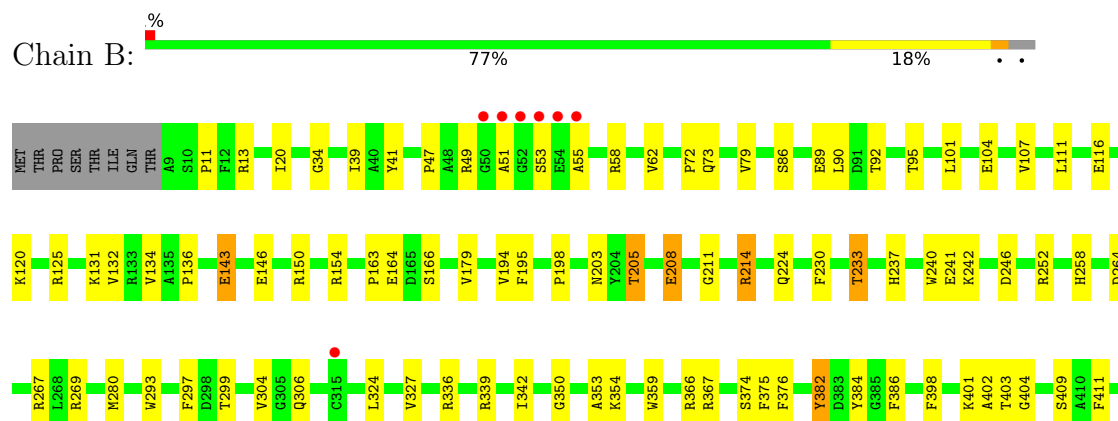
3 Residue-property plots

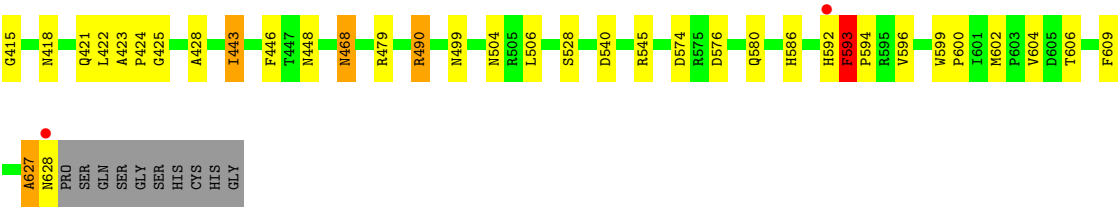
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Phenylethylamine oxidase



• Molecule 1: Phenylethylamine oxidase





4 Data and refinement statistics

Property	Value	Source
Space group	I 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	157.80Å 63.07Å 183.46Å 90.00° 111.64° 90.00°	Depositor
Resolution (Å)	7.00 – 2.10 7.00 – 2.10	Depositor EDS
% Data completeness (in resolution range)	(Not available) (7.00-2.10) 78.5 (7.00-2.10)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.05 (at 1.90Å)	Xtriage
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.190 , 0.263 0.172 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	21.4	Xtriage
Anisotropy	0.724	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.47 , 124.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	10959	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 35.95 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 5.3863e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 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: CU, DAH

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	0.61	0/4986	1.06	26/6789 (0.4%)
1	B	0.62	0/4986	1.05	23/6789 (0.3%)
All	All	0.61	0/9972	1.05	49/13578 (0.4%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	2

There are no bond length outliers.

All (49) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	565	ALA	N-CA-C	13.14	125.13	111.07
1	B	593	PHE	O-C-N	-9.79	110.06	121.32
1	B	627	ALA	N-CA-C	-8.04	103.06	113.12
1	A	446	PHE	N-CA-C	7.84	120.91	111.82
1	B	34	GLY	CA-C-N	7.72	128.42	119.47
1	B	34	GLY	C-N-CA	7.72	128.42	119.47
1	A	288	SER	CA-C-N	6.52	126.46	119.28
1	A	288	SER	C-N-CA	6.52	126.46	119.28
1	B	224	GLN	CA-C-N	6.48	126.71	119.32
1	B	224	GLN	C-N-CA	6.48	126.71	119.32
1	B	604	VAL	N-CA-C	6.48	117.79	109.30
1	A	46	ASP	CA-C-N	6.34	127.77	119.84
1	A	46	ASP	C-N-CA	6.34	127.77	119.84
1	A	604	VAL	N-CA-C	6.28	118.49	109.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	581	ASP	N-CA-C	-6.28	100.11	109.15
1	B	107	VAL	N-CA-C	6.03	116.69	110.36
1	B	267	ARG	N-CA-C	6.00	118.73	109.07
1	A	562	SER	N-CA-C	5.96	119.81	112.54
1	A	409	SER	N-CA-C	-5.92	100.99	108.45
1	A	46	ASP	CB-CA-C	5.91	117.37	108.86
1	A	324	LEU	N-CA-C	-5.91	99.77	109.40
1	A	374	SER	N-CA-C	5.78	117.57	108.96
1	B	374	SER	N-CA-C	5.69	117.43	108.96
1	B	304	VAL	N-CA-C	5.63	116.41	110.72
1	B	195	PHE	CA-C-N	-5.58	114.18	119.76
1	B	195	PHE	C-N-CA	-5.58	114.18	119.76
1	A	268	LEU	N-CA-C	-5.51	98.78	108.20
1	B	95	THR	N-CA-C	5.50	119.99	113.28
1	A	408	THR	N-CA-C	5.45	118.52	110.46
1	A	195	PHE	N-CA-C	-5.41	100.02	109.48
1	A	545	ARG	N-CA-C	-5.35	102.36	110.28
1	B	92	THR	N-CA-C	5.32	118.87	112.38
1	B	132	VAL	N-CA-C	5.28	115.38	107.51
1	A	499	ASN	CA-C-N	5.26	125.57	119.47
1	A	499	ASN	C-N-CA	5.26	125.57	119.47
1	A	199	ALA	N-CA-C	5.21	118.10	111.69
1	A	607	VAL	CB-CA-C	-5.20	101.39	110.38
1	B	359	TRP	N-CA-C	5.20	117.62	111.33
1	A	297	PHE	N-CA-C	-5.19	98.38	107.73
1	A	593	PHE	N-CA-C	-5.18	98.36	109.81
1	A	179	VAL	N-CA-C	5.17	115.34	108.11
1	A	157	ALA	N-CA-C	5.15	118.61	109.96
1	B	297	PHE	N-CA-C	-5.13	97.61	107.57
1	B	446	PHE	N-CA-C	5.13	119.26	112.89
1	A	359	TRP	N-CA-C	5.08	116.51	111.07
1	B	504	ASN	N-CA-C	-5.08	103.34	110.35
1	B	211	GLY	N-CA-C	-5.06	102.01	112.34
1	B	545	ARG	N-CA-C	-5.05	101.73	109.76
1	B	409	SER	N-CA-C	-5.00	102.15	108.45

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	41	TYR	Sidechain
1	B	593	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	4872	0	4689	125	0
1	B	4872	0	4688	85	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	579	0	0	13	0
3	B	634	0	0	16	0
All	All	10959	0	9377	198	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 10.

All (198) 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:B:208:GLU:HG3	3:B:1573:HOH:O	1.60	0.99
1:A:238:ILE:HD11	1:A:344:MET:HE2	1.46	0.96
1:A:45:LEU:HD11	1:A:61:ARG:HB2	1.49	0.93
1:A:170:HIS:HD2	1:A:198:PRO:O	1.62	0.83
1:A:458:GLN:HE22	1:B:418:ASN:HD21	1.22	0.83
1:A:247:VAL:HG21	1:A:344:MET:CE	2.09	0.82
1:A:46:ASP:HB2	3:A:1025:HOH:O	1.79	0.81
1:B:490:ARG:HH11	1:B:490:ARG:HB2	1.45	0.80
1:A:403:THR:HG21	1:B:353:ALA:HA	1.63	0.80
1:A:592[A]:HIS:CE1	3:A:1005:HOH:O	2.33	0.79
1:B:55:ALA:HB2	3:B:1588:HOH:O	1.85	0.76
1:A:590:LEU:HG	1:A:607:VAL:HG22	1.69	0.74
1:A:460:MET:HE1	1:B:528:SER:HA	1.71	0.72
1:B:73:GLN:HE21	1:B:89:GLU:HG2	1.54	0.71
1:B:382:DAH:HD2	1:B:403:THR:O	1.90	0.71
1:A:350:GLY:HA2	1:A:367:ARG:NH2	2.07	0.70
1:A:264:ASP:HB2	1:A:269:ARG:HD3	1.76	0.68
1:B:443:ILE:H	1:B:448:ASN:HD21	1.42	0.68
1:B:166:SER:H	1:B:203:ASN:HD21	1.42	0.67
1:A:484:VAL:HG12	1:A:539:LYS:HG3	1.78	0.64
1:A:45:LEU:CD1	1:A:61:ARG:HB2	2.27	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:324:LEU:HB2	1:A:342:ILE:HB	1.81	0.62
1:A:443:ILE:H	1:A:448:ASN:HD21	1.47	0.61
1:A:102:GLU:O	1:A:106:GLU:HG3	2.00	0.61
1:A:238:ILE:HD11	1:A:344:MET:CE	2.27	0.61
1:A:504:ASN:ND2	1:A:508:GLU:HG2	2.16	0.61
1:A:125:ARG:HG2	1:A:194:VAL:HG23	1.83	0.60
1:A:403:THR:HG22	1:A:602:MET:O	2.02	0.60
1:B:20:ILE:HD12	1:B:327:VAL:HG12	1.84	0.59
1:B:592[A]:HIS:HE2	1:B:594:PRO:HA	1.68	0.59
1:A:576:ASP:O	1:A:580:GLN:HG3	2.02	0.59
1:A:187:THR:O	1:A:188:ARG:HG3	2.02	0.58
1:B:490:ARG:HH11	1:B:490:ARG:CB	2.14	0.57
1:B:280:MET:HE1	1:B:384:TYR:CE2	2.40	0.56
1:B:241:GLU:O	1:B:242:LYS:HB2	2.04	0.56
1:B:233:THR:HG22	1:B:237:HIS:HB3	1.87	0.56
1:A:506:LEU:HD13	1:B:421:GLN:HG2	1.86	0.56
1:B:143:GLU:H	1:B:143:GLU:CD	2.14	0.56
1:A:19:GLU:OE2	1:A:58:ARG:HG2	2.06	0.56
1:A:458:GLN:NE2	1:B:418:ASN:HD21	2.00	0.55
1:A:63:PHE:CE1	1:A:74:GLU:HG3	2.42	0.55
1:A:78:SER:HB2	1:A:85:ILE:HD11	1.88	0.55
1:A:19:GLU:O	1:A:23:VAL:HG23	2.07	0.55
1:B:425:GLY:HA2	3:B:1424:HOH:O	2.07	0.55
1:A:595:ARG:HG2	1:A:598:ASP:OD2	2.07	0.55
1:A:46:ASP:CB	1:A:47:PRO:HD2	2.36	0.54
1:A:590:LEU:HG	1:A:607:VAL:CG2	2.37	0.54
1:A:24:GLN:HE22	1:A:334:ASN:ND2	2.06	0.54
1:A:246:ASP:HB2	1:A:258:HIS:HB2	1.89	0.54
1:A:230:PHE:HB3	1:A:240:TRP:HB2	1.89	0.54
1:A:504:ASN:HD21	1:A:508:GLU:HG2	1.71	0.54
1:B:479:ARG:HB3	1:B:574:ASP:OD1	2.08	0.54
1:A:73:GLN:HG2	1:A:89:GLU:HA	1.90	0.53
1:B:422:LEU:HD11	1:B:428:ALA:HB2	1.91	0.53
1:A:599:TRP:CD2	1:A:600:PRO:HA	2.44	0.53
1:B:339:ARG:HG3	3:B:1550:HOH:O	2.09	0.53
1:A:421:GLN:HG2	1:B:506:LEU:HD13	1.91	0.53
1:B:599:TRP:CD2	1:B:600:PRO:HA	2.43	0.53
1:A:98:LEU:HB2	1:A:151:ARG:HH21	1.73	0.53
1:A:411:PHE:CZ	1:A:415:GLY:HA2	2.43	0.53
1:A:66:ASP:HB3	1:A:70:ALA:O	2.09	0.52
1:A:368:ASN:ND2	1:A:369:ARG:H	2.07	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:19:GLU:HG2	1:A:79:VAL:HG11	1.90	0.52
1:B:101:LEU:HB2	1:B:104:GLU:HG3	1.91	0.52
1:B:13:ARG:HD2	3:B:1483:HOH:O	2.10	0.52
1:A:247:VAL:HG21	1:A:344:MET:HE3	1.87	0.52
1:B:72:PRO:HG2	1:B:90:LEU:HB2	1.91	0.52
1:B:125:ARG:HG2	1:B:194:VAL:HG13	1.91	0.52
1:B:237:HIS:HD2	1:B:246:ASP:OD1	1.93	0.52
1:A:460:MET:HE2	1:A:466:ARG:C	2.35	0.51
1:B:366:ARG:CZ	1:B:627:ALA:HB2	2.41	0.51
1:A:627:ALA:O	1:A:628:ASN:HB2	2.11	0.51
1:A:484:VAL:CG1	1:A:539:LYS:HG3	2.40	0.51
1:B:593:PHE:O	1:B:594:PRO:C	2.53	0.51
1:B:443:ILE:H	1:B:448:ASN:ND2	2.09	0.51
1:B:592[A]:HIS:NE2	1:B:594:PRO:HA	2.25	0.51
1:A:504:ASN:CG	1:A:508:GLU:HG2	2.36	0.50
1:B:205:THR:HB	3:B:1174:HOH:O	2.10	0.50
1:B:120:LYS:HE2	3:B:1508:HOH:O	2.11	0.50
1:A:150:ARG:HD3	3:A:1536:HOH:O	2.11	0.50
1:B:402:ALA:HB1	3:B:1473:HOH:O	2.12	0.50
1:A:46:ASP:HB3	1:A:47:PRO:HD2	1.94	0.50
1:A:624:ASP:CG	1:B:214:ARG:HD2	2.37	0.50
1:A:436:SER:HB2	1:A:536:PHE:CE2	2.47	0.50
1:A:134:VAL:O	1:A:136:PRO:HD3	2.12	0.49
1:A:403:THR:HG21	1:B:354:LYS:H	1.77	0.49
1:A:345:HIS:HE1	3:A:1280:HOH:O	1.95	0.49
1:A:561:HIS:CE1	1:A:565:ALA:HB2	2.48	0.49
1:B:264:ASP:HB3	1:B:269:ARG:HD2	1.94	0.49
1:A:98:LEU:HB2	1:A:151:ARG:NH2	2.28	0.49
1:A:483:ALA:HB1	1:A:543:VAL:HB	1.93	0.49
1:A:13:ARG:NH2	3:A:1399:HOH:O	2.45	0.49
1:B:324:LEU:HB2	1:B:342:ILE:HB	1.95	0.48
1:A:368:ASN:HD22	1:A:392:LEU:HD12	1.79	0.48
1:A:347:GLU:OE1	1:A:370:ARG:HD2	2.13	0.48
1:A:20:ILE:HB	1:A:335:PRO:HB3	1.96	0.48
1:A:13:ARG:HH12	1:A:58:ARG:HE	1.62	0.48
1:B:134:VAL:O	1:B:136:PRO:HD3	2.13	0.47
1:B:468:ASN:HD22	1:B:468:ASN:H	1.62	0.47
1:A:12:PHE:CZ	1:A:47:PRO:HG2	2.49	0.47
1:A:63:PHE:HE1	1:A:74:GLU:HG3	1.79	0.47
1:A:13:ARG:NH1	1:A:58:ARG:NE	2.61	0.47
1:A:66:ASP:CG	1:A:70:ALA:HB3	2.38	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:75:VAL:HG13	1:A:86:SER:O	2.15	0.47
1:A:139:ALA:O	1:A:154:ARG:NH1	2.48	0.47
1:B:443:ILE:N	1:B:448:ASN:HD21	2.11	0.47
1:A:504:ASN:OD1	1:A:508:GLU:HG2	2.14	0.47
1:B:540:ASP:O	1:B:586:HIS:HD2	1.97	0.47
1:B:111:LEU:HD12	1:B:179:VAL:CG1	2.45	0.47
1:A:355:HIS:HB2	1:B:404:GLY:HA2	1.97	0.46
1:A:242:LYS:NZ	3:A:1438:HOH:O	2.48	0.46
1:B:592[B]:HIS:CE1	1:B:602:MET:HE1	2.50	0.46
1:B:146:GLU:O	1:B:150:ARG:HD3	2.16	0.46
1:B:375:PHE:CE1	1:B:386:PHE:HB2	2.50	0.46
1:A:266:ASP:OD2	1:A:267:ARG:HD3	2.16	0.46
1:A:309:ASN:HD22	1:B:354:LYS:HE3	1.80	0.45
1:A:508:GLU:HB2	3:A:1370:HOH:O	2.16	0.45
1:A:510:VAL:HB	1:A:616:PHE:HA	1.98	0.45
1:A:178:TYR:HD2	1:A:188:ARG:HD2	1.80	0.45
1:B:339:ARG:HD2	3:B:1512:HOH:O	2.15	0.45
1:B:39:ILE:HD12	1:B:62:VAL:HG11	1.97	0.45
1:B:411:PHE:CZ	1:B:415:GLY:HA2	2.51	0.45
1:A:398:PHE:HB3	1:A:609:PHE:CE2	2.51	0.45
1:A:162:PHE:HB2	1:A:163:PRO:CD	2.46	0.45
1:A:403:THR:CG2	1:B:353:ALA:HA	2.40	0.45
1:B:86:SER:HA	3:B:1569:HOH:O	2.16	0.45
1:A:13:ARG:HH12	1:A:58:ARG:NE	2.14	0.45
1:B:280:MET:HE1	1:B:384:TYR:CD2	2.52	0.44
1:B:252:ARG:O	1:B:299:THR:HG23	2.17	0.44
1:B:230:PHE:HB3	1:B:240:TRP:HB2	1.99	0.44
1:A:160:GLN:NE2	1:A:163:PRO:O	2.51	0.44
1:A:421:GLN:NE2	1:A:425:GLY:H	2.15	0.44
1:A:551:ARG:HG2	1:A:551:ARG:HH11	1.83	0.44
1:B:586:HIS:HE1	3:B:1010:HOH:O	2.00	0.44
1:A:435:PHE:CD1	1:A:435:PHE:N	2.86	0.44
1:A:225:PRO:HA	3:A:1493:HOH:O	2.18	0.44
1:A:369:ARG:HG2	1:A:392:LEU:HD11	2.00	0.44
1:A:182:VAL:HG21	3:A:1487:HOH:O	2.18	0.43
1:A:291:ARG:HG2	1:A:294:GLN:NE2	2.33	0.43
1:A:29:THR:HG21	3:A:1464:HOH:O	2.17	0.43
1:A:628:ASN:HB2	3:A:1425:HOH:O	2.17	0.43
1:A:61:ARG:HH22	1:A:556:ASP:CG	2.26	0.43
1:B:208:GLU:CG	3:B:1573:HOH:O	2.41	0.43
1:A:264:ASP:HB2	1:A:269:ARG:CD	2.47	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:401:LYS:HG2	1:A:606:THR:HG22	2.00	0.43
1:A:503:LYS:HD3	1:A:507:ASN:ND2	2.34	0.43
1:B:576:ASP:O	1:B:580:GLN:HG3	2.18	0.43
1:A:542:TRP:HB2	1:A:585:TRP:HB2	2.01	0.43
1:A:152:ILE:HG22	1:A:153:LEU:N	2.34	0.43
1:B:154:ARG:HD2	1:B:293:TRP:CD2	2.53	0.43
1:B:411:PHE:HB3	3:B:1424:HOH:O	2.18	0.43
1:A:143:GLU:CD	1:A:143:GLU:H	2.27	0.43
1:B:398:PHE:HB3	1:B:609:PHE:CE2	2.54	0.43
1:B:214:ARG:HG2	3:B:1317:HOH:O	2.20	0.42
1:A:50:GLY:O	1:A:51:ALA:HB2	2.19	0.42
1:A:268:LEU:C	1:A:268:LEU:HD13	2.44	0.42
1:B:596:VAL:HG23	3:B:1203:HOH:O	2.19	0.42
1:A:22:GLU:O	1:A:26:ILE:HG13	2.19	0.42
1:B:246:ASP:HB2	1:B:258:HIS:HB2	2.00	0.42
1:B:401:LYS:HG2	1:B:606:THR:HG22	2.01	0.42
1:B:366:ARG:NH1	1:B:627:ALA:HB2	2.34	0.42
1:A:441:MET:HG3	1:A:497:ILE:HG21	2.02	0.42
1:A:458:GLN:HE22	1:B:418:ASN:ND2	2.03	0.42
1:B:423:ALA:HB1	1:B:424:PRO:CD	2.50	0.42
1:A:162:PHE:HB2	1:A:163:PRO:HD2	2.02	0.42
1:A:174:GLY:O	1:A:191:ASP:HA	2.19	0.42
1:A:413:GLU:HA	3:A:1381:HOH:O	2.19	0.42
1:A:120:LYS:HB2	1:A:120:LYS:HE3	1.75	0.42
1:A:169:ALA:O	1:A:198:PRO:HB2	2.19	0.42
1:A:182:VAL:HG23	1:A:183:SER:N	2.35	0.42
1:A:194:VAL:HG22	3:A:1029:HOH:O	2.19	0.41
1:B:154:ARG:HD2	1:B:293:TRP:CE3	2.54	0.41
1:B:350:GLY:HA2	1:B:367:ARG:NH2	2.35	0.41
1:A:468:ASN:HD22	1:A:468:ASN:H	1.68	0.41
1:A:545:ARG:HD2	1:A:581:ASP:O	2.20	0.41
1:A:153:LEU:HD23	1:A:153:LEU:HA	1.88	0.41
1:B:11:PRO:HB2	1:B:47:PRO:HG3	2.02	0.41
1:B:73:GLN:NE2	1:B:89:GLU:HG2	2.27	0.41
1:B:382:DAH:OE2	1:B:602:MET:SD	2.79	0.41
1:A:170:HIS:CD2	1:A:198:PRO:O	2.54	0.41
1:B:131:LYS:HE3	3:B:1556:HOH:O	2.20	0.41
1:B:443:ILE:O	1:B:499:ASN:HB2	2.20	0.41
1:A:59:ARG:NE	1:A:85:ILE:HD13	2.36	0.41
1:A:133:ARG:NE	1:A:133:ARG:HA	2.36	0.41
1:A:42:LEU:C	1:A:42:LEU:HD23	2.46	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:51:ALA:O	1:B:53:SER:N	2.54	0.41
1:A:73:GLN:HG2	1:A:89:GLU:HG2	2.02	0.41
1:A:247:VAL:HG21	1:A:344:MET:HE1	2.01	0.41
1:A:100:VAL:HG11	1:A:105:PHE:CE1	2.55	0.40
1:B:164:GLU:CD	1:B:164:GLU:H	2.29	0.40
1:A:460:MET:HE3	1:A:467:GLY:HA3	2.02	0.40
1:A:182:VAL:CG2	1:A:183:SER:N	2.84	0.40
1:A:237:HIS:HD2	1:A:246:ASP:OD1	2.05	0.40
1:B:58:ARG:O	1:B:79:VAL:HG12	2.22	0.40
1:B:163:PRO:HB2	1:B:164:GLU:OE1	2.22	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	618/638 (97%)	582 (94%)	32 (5%)	4 (1%)	21	18
1	B	618/638 (97%)	584 (94%)	32 (5%)	2 (0%)	36	36
All	All	1236/1276 (97%)	1166 (94%)	64 (5%)	6 (0%)	24	22

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	51	ALA
1	A	47	PRO
1	A	55	ALA
1	B	593	PHE
1	B	443	ILE
1	A	443	ILE

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	514/529 (97%)	491 (96%)	23 (4%)	24	25
1	B	514/529 (97%)	500 (97%)	14 (3%)	39	45
All	All	1028/1058 (97%)	991 (96%)	37 (4%)	31	34

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	45	LEU
1	A	46	ASP
1	A	54	GLU
1	A	77	VAL
1	A	102	GLU
1	A	116	GLU
1	A	143	GLU
1	A	198	PRO
1	A	208	GLU
1	A	213	LEU
1	A	239	GLU
1	A	251	VAL
1	A	269	ARG
1	A	370	ARG
1	A	376	PHE
1	A	403	THR
1	A	459	THR
1	A	468	ASN
1	A	541	LEU
1	A	551	ARG
1	A	554	THR
1	A	597	GLU
1	A	607	VAL
1	B	49	ARG
1	B	116	GLU
1	B	143	GLU
1	B	198	PRO

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Mol	Chain	Res	Type
1	B	205	THR
1	B	208	GLU
1	B	214	ARG
1	B	233	THR
1	B	306	GLN
1	B	336	ARG
1	B	376	PHE
1	B	468	ASN
1	B	490	ARG
1	B	628	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (31) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	170	HIS
1	A	237	HIS
1	A	294	GLN
1	A	306	GLN
1	A	309	ASN
1	A	334	ASN
1	A	345	HIS
1	A	368	ASN
1	A	421	GLN
1	A	448	ASN
1	A	468	ASN
1	A	507	ASN
1	A	515	HIS
1	A	573	GLN
1	B	73	GLN
1	B	203	ASN
1	B	224	GLN
1	B	237	HIS
1	B	273	ASN
1	B	306	GLN
1	B	309	ASN
1	B	334	ASN
1	B	340	ASN
1	B	345	HIS
1	B	418	ASN
1	B	421	GLN
1	B	448	ASN
1	B	468	ASN

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Mol	Chain	Res	Type
1	B	515	HIS
1	B	519	GLN
1	B	586	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues are modelled in this entry.

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
1	DAH	B	382	1	12,13,14	0.95	1 (8%)	12,17,19	1.30	2 (16%)
1	DAH	A	382	1,2	12,13,14	1.04	0	12,17,19	1.12	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	DAH	B	382	1	-	4/5/6/8	0/1/1/1
1	DAH	A	382	1,2	-	5/5/6/8	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	382	DAH	CD1-CG	2.06	1.43	1.38

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	B	382	DAH	CB-CG-CD2	2.65	124.97	120.43
1	B	382	DAH	CB-CG-CD1	-2.03	117.12	120.90

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	382	DAH	O-C-CA-CB
1	A	382	DAH	C-CA-CB-CG
1	B	382	DAH	C-CA-CB-CG
1	A	382	DAH	N-CA-CB-CG
1	A	382	DAH	CA-CB-CG-CD2
1	A	382	DAH	CA-CB-CG-CD1
1	B	382	DAH	CA-CB-CG-CD2
1	B	382	DAH	N-CA-CB-CG
1	B	382	DAH	CA-CB-CG-CD1

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	B	382	DAH	2	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

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

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 [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	619/638 (97%)	-0.61	7 (1%) 78 80	8, 22, 42, 82	1 (0%)
1	B	619/638 (97%)	-0.74	9 (1%) 72 74	4, 19, 37, 79	1 (0%)
All	All	1238/1276 (97%)	-0.68	16 (1%) 75 77	4, 21, 40, 82	2 (0%)

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	50	GLY	7.1
1	B	51	ALA	5.8
1	A	52	GLY	5.2
1	A	51	ALA	4.0
1	A	50	GLY	3.7
1	A	628	ASN	3.6
1	B	592[A]	HIS	3.4
1	B	628	ASN	3.2
1	A	46	ASP	3.0
1	A	55	ALA	3.0
1	A	54	GLU	2.9
1	B	52	GLY	2.7
1	B	54	GLU	2.4
1	B	55	ALA	2.4
1	B	53	SER	2.3
1	B	315	CYS	2.1

6.2 Non-standard residues in protein, DNA, RNA chains [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($\text{\AA}^2$)	Q<0.9
1	DAH	B	382	13/14	0.80	0.14	9,15,19,24	1
1	DAH	A	382	13/14	0.95	0.07	15,20,25,25	1

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($\text{\AA}^2$)	Q<0.9
2	CU	B	1002	1/1	0.89	0.06	25,25,25,25	0
2	CU	A	1001	1/1	0.98	0.06	26,26,26,26	0

6.5 Other polymers [i](#)

There are no such residues in this entry.