



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 11:22 AM UTC

PDB ID : 2CWT / pdb_00002cwt
Title : Catalytic base deletion in copper amine oxidase from arthrobacter globiformis
Authors : Chiu, Y.C.; Okajima, T.; Murakawa, T.; Uchida, M.; Taki, M.; Hirota, S.; Kim, M.; Yamaguchi, H.; Kawano, Y.; Kamiya, N.; Kuroda, S.; Hayashi, H.; Yamamoto, Y.; Tanizawa, K.
Deposited on : 2005-06-26
Resolution : 1.82 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)
Xtrriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

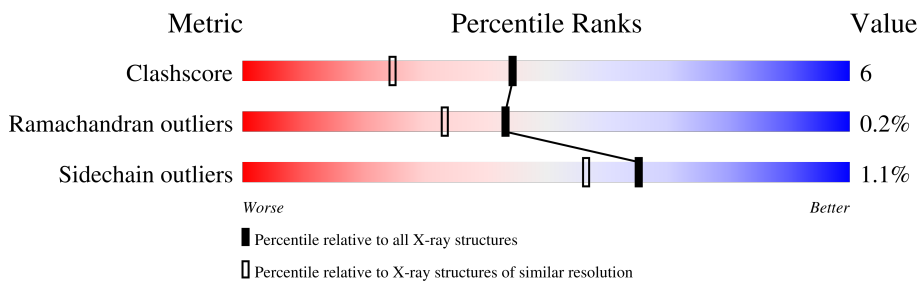
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.82 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	1148 (1.82-1.82)
Ramachandran outliers	187476	1140 (1.82-1.82)
Sidechain outliers	187428	1140 (1.82-1.82)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	638	
1	B	638	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 10543 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
			Total	C	N	O	S			
1	A	620	4865	3073	855	928	9	0	0	0
1	B	620	4865	3073	855	928	9	0	0	0

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	298	ALA	ASP	engineered mutation	UNP P46881
A	382	TPQ	TYR	modified residue	UNP P46881
B	298	ALA	ASP	engineered mutation	UNP P46881
B	382	TPQ	TYR	modified residue	UNP P46881

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

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

- Molecule 3 is water.

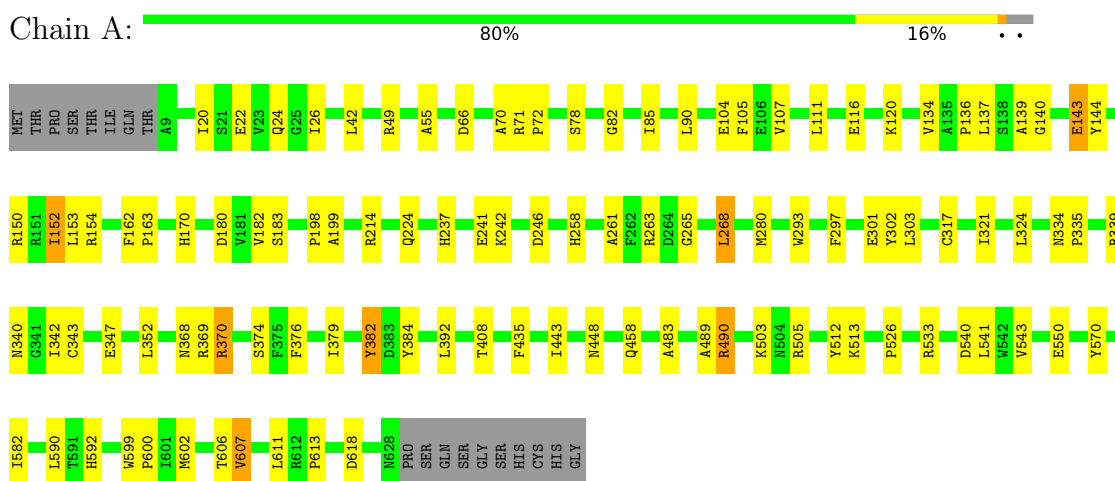
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	395	Total 395	O 395	0	0
3	B	416	Total 416	O 416	0	0

3 Residue-property plots [i](#)

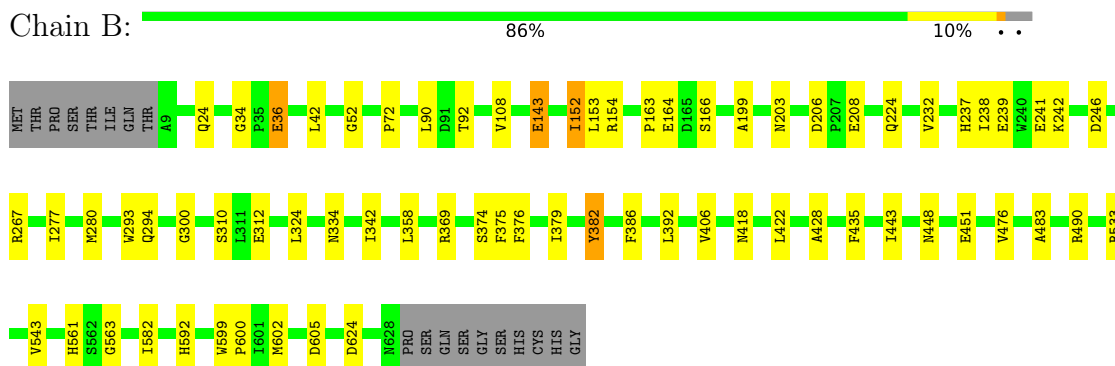
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded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

Note EDS was not executed.

- Molecule 1: Phenylethylamine oxidase



- Molecule 1: Phenylethylamine oxidase



4 Data and refinement statistics

Xtrriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	I 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	157.67Å 63.84Å 184.45Å 90.00° 112.40° 90.00°	Depositor
Resolution (Å)	30.02 – 1.82	Depositor
% Data completeness (in resolution range)	(Not available) (30.02-1.82)	Depositor
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS	Depositor
R, R_{free}	0.191 , 0.204	Depositor
Estimated twinning fraction	No twinning to report.	Xtrriage
Total number of atoms	10543	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: CU, TPQ

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	0.33	0/4972	0.90	12/6770 (0.2%)
1	B	0.34	0/4972	0.92	18/6770 (0.3%)
All	All	0.34	0/9944	0.91	30/13540 (0.2%)

There are no bond length outliers.

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	602	MET	N-CA-C	7.80	119.60	109.93
1	B	294	GLN	N-CA-C	6.85	119.01	109.15
1	A	602	MET	N-CA-C	6.71	118.52	110.07
1	A	224	GLN	CA-C-N	6.56	126.80	119.32
1	A	224	GLN	C-N-CA	6.56	126.80	119.32
1	A	199	ALA	N-CA-C	6.27	119.41	111.69
1	B	224	GLN	CA-C-N	6.24	126.44	119.32
1	B	224	GLN	C-N-CA	6.24	126.44	119.32
1	B	310	SER	N-CA-C	-6.09	99.28	108.96
1	B	277	ILE	N-CA-C	-6.05	97.68	106.88
1	A	592	HIS	N-CA-C	6.04	118.57	108.73
1	A	374	SER	N-CA-C	6.03	117.94	108.96
1	B	34	GLY	CA-C-N	5.88	126.29	119.47
1	B	34	GLY	C-N-CA	5.88	126.29	119.47
1	B	108	VAL	N-CA-C	5.88	116.34	110.23
1	A	152	ILE	N-CA-C	5.77	117.67	108.89
1	A	321	ILE	N-CA-C	5.76	116.60	108.36
1	B	152	ILE	N-CA-C	5.75	117.63	108.89
1	A	582	ILE	N-CA-C	5.73	117.09	108.96
1	B	300	GLY	N-CA-C	5.72	120.09	112.77
1	B	199	ALA	N-CA-C	5.65	118.64	111.69
1	B	92	THR	N-CA-C	5.45	117.22	111.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	582	ILE	N-CA-C	5.42	117.44	108.99
1	B	374	SER	N-CA-C	5.38	116.97	108.96
1	A	408	THR	N-CA-C	5.29	117.92	110.14
1	A	606	THR	N-CA-C	5.24	117.64	109.52
1	B	592	HIS	N-CA-C	5.23	117.04	108.52
1	B	605	ASP	N-CA-C	-5.22	101.37	109.72
1	B	563	GLY	N-CA-C	-5.15	104.63	112.51
1	A	540	ASP	N-CA-C	5.05	117.68	111.82

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	4865	0	4689	83	0
1	B	4865	0	4687	35	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	395	0	0	2	0
3	B	416	0	0	4	0
All	All	10543	0	9376	115	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 6.

All (115) 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:458:GLN:HE22	1:B:418:ASN:HD21	1.26	0.82
1:A:590:LEU:HG	1:A:607:VAL:CG2	2.09	0.81
1:A:139:ALA:HA	1:A:154:ARG:NH1	1.99	0.77
1:A:590:LEU:HG	1:A:607:VAL:HG22	1.65	0.76
1:A:104:GLU:O	1:A:107:VAL:HG12	1.85	0.76
1:B:166:SER:H	1:B:203:ASN:HD21	1.34	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:150:ARG:HA	1:A:150:ARG:HE	1.57	0.70
1:A:150:ARG:HD3	1:A:180:ASP:OD2	1.93	0.69
1:A:154:ARG:HH11	1:A:154:ARG:HG3	1.59	0.68
1:A:72:PRO:HG2	1:A:90:LEU:HB2	1.76	0.67
1:A:347:GLU:OE1	1:A:370:ARG:HD3	1.96	0.66
1:B:143:GLU:H	1:B:143:GLU:CD	2.05	0.65
1:A:143:GLU:HG3	1:A:144:TYR:H	1.63	0.63
1:A:489:ALA:H	1:A:490:ARG:HH21	1.47	0.63
1:A:379:ILE:O	1:A:382:TPQ:H6	1.99	0.62
1:A:503:LYS:HB2	1:A:503:LYS:HZ2	1.62	0.62
1:A:49:ARG:NE	1:A:49:ARG:HA	2.15	0.62
1:A:180:ASP:OD1	1:A:182:VAL:HG22	2.00	0.62
1:B:533:ARG:HD2	1:B:561:HIS:O	2.00	0.61
1:A:24:GLN:HE22	1:A:334:ASN:ND2	1.97	0.61
1:B:599:TRP:CD2	1:B:600:PRO:HA	2.35	0.61
1:A:505:ARG:HD3	1:A:618:ASP:HB3	1.83	0.61
1:A:590:LEU:HG	1:A:607:VAL:HG21	1.80	0.61
1:A:590:LEU:CG	1:A:607:VAL:HG22	2.31	0.60
1:A:599:TRP:CD2	1:A:600:PRO:HA	2.36	0.60
1:A:66:ASP:OD1	1:A:70:ALA:HB3	2.00	0.60
1:B:379:ILE:O	1:B:382:TPQ:H6	2.01	0.60
1:B:206:ASP:OD1	1:B:208:GLU:HG2	2.01	0.59
1:A:324:LEU:HB2	1:A:342:ILE:HB	1.84	0.59
1:A:261:ALA:HB1	1:A:268:LEU:HD22	1.85	0.59
1:A:66:ASP:CG	1:A:70:ALA:HB3	2.28	0.58
1:B:443:ILE:H	1:B:448:ASN:HD21	1.52	0.58
1:A:150:ARG:HA	1:A:150:ARG:NE	2.18	0.57
1:A:170:HIS:HD2	1:A:198:PRO:O	1.87	0.57
1:A:526:PRO:HG3	3:B:1307:HOH:O	2.05	0.56
1:B:422:LEU:HD11	1:B:428:ALA:HB2	1.88	0.56
1:A:105:PHE:HB3	1:A:136:PRO:HG3	1.86	0.56
1:A:489:ALA:H	1:A:490:ARG:NH2	2.06	0.54
1:A:503:LYS:HB2	1:A:503:LYS:NZ	2.23	0.54
1:B:241:GLU:O	1:B:242:LYS:HB2	2.08	0.53
1:A:237:HIS:HD2	1:A:246:ASP:OD1	1.91	0.53
1:A:317:CYS:SG	1:A:343:CYS:SG	3.07	0.53
1:A:443:ILE:H	1:A:448:ASN:HD21	1.57	0.52
1:A:483:ALA:HB1	1:A:543:VAL:HB	1.91	0.52
1:A:49:ARG:HA	1:A:49:ARG:HE	1.73	0.52
1:A:368:ASN:ND2	1:A:369:ARG:H	2.07	0.52
1:B:163:PRO:HB2	1:B:164:GLU:OE2	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:317:CYS:HG	1:A:343:CYS:CB	2.22	0.52
1:A:49:ARG:HE	1:A:49:ARG:CA	2.24	0.51
1:B:451:GLU:HG3	3:B:1112:HOH:O	2.08	0.51
1:B:533:ARG:CD	1:B:561:HIS:O	2.57	0.51
1:A:370:ARG:O	1:A:370:ARG:HG2	2.12	0.50
1:B:237:HIS:HD2	1:B:246:ASP:OD1	1.95	0.50
1:A:297:PHE:HB2	1:A:301:GLU:HG3	1.93	0.50
1:A:20:ILE:HB	1:A:335:PRO:HG3	1.93	0.49
1:A:241:GLU:O	1:A:242:LYS:HB2	2.12	0.49
1:B:451:GLU:HG2	1:B:476:VAL:HG22	1.93	0.49
1:A:152:ILE:HG22	1:A:153:LEU:N	2.28	0.49
1:A:182:VAL:HG23	1:A:183:SER:N	2.28	0.49
1:A:368:ASN:HD22	1:A:369:ARG:H	1.61	0.49
1:B:324:LEU:HB2	1:B:342:ILE:HB	1.95	0.49
1:B:72:PRO:HG2	1:B:90:LEU:HB2	1.96	0.48
1:A:317:CYS:SG	1:A:343:CYS:CB	3.02	0.48
1:B:42:LEU:C	1:B:42:LEU:HD23	2.38	0.48
1:A:42:LEU:C	1:A:42:LEU:HD23	2.39	0.48
1:A:49:ARG:NE	1:A:49:ARG:CA	2.77	0.48
1:B:561:HIS:HD2	3:B:1083:HOH:O	1.96	0.47
1:B:239:GLU:HG3	3:B:1040:HOH:O	2.15	0.47
1:A:317:CYS:SG	1:A:343:CYS:HB2	2.55	0.47
1:B:232:VAL:HG22	1:B:238:ILE:CD1	2.45	0.47
1:A:78:SER:HB2	1:A:85:ILE:HD11	1.97	0.47
1:A:261:ALA:CB	1:A:268:LEU:HD22	2.45	0.46
1:A:152:ILE:HD13	1:A:180:ASP:HA	1.98	0.46
1:B:369:ARG:HG2	1:B:392:LEU:HD11	1.98	0.46
1:A:139:ALA:HA	1:A:154:ARG:HH11	1.75	0.46
1:A:339:ARG:NE	1:A:340:ASN:ND2	2.64	0.46
1:A:246:ASP:HB2	1:A:258:HIS:HB2	1.97	0.46
1:A:105:PHE:HD1	1:A:136:PRO:HB2	1.81	0.45
1:A:214:ARG:HE	1:B:624:ASP:CG	2.25	0.45
1:A:263:ARG:HD2	1:A:265:GLY:O	2.17	0.45
1:B:143:GLU:CD	1:B:143:GLU:N	2.72	0.45
1:B:24:GLN:HE22	1:B:334:ASN:ND2	2.14	0.45
1:A:22:GLU:O	1:A:26:ILE:HG13	2.16	0.45
1:A:107:VAL:HG22	1:A:111:LEU:HG	1.98	0.44
1:A:280:MET:HE1	1:A:384:TYR:CE2	2.52	0.44
1:B:406:VAL:HG21	1:B:422:LEU:HD11	1.98	0.44
1:B:36:GLU:H	1:B:36:GLU:CD	2.23	0.44
1:A:134:VAL:O	1:A:136:PRO:HD3	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:152:ILE:CG2	1:A:153:LEU:N	2.81	0.43
1:B:490:ARG:HH21	1:B:490:ARG:HG3	1.83	0.43
1:A:71:ARG:HB2	3:A:1341:HOH:O	2.16	0.43
1:A:154:ARG:NH1	1:A:154:ARG:HG3	2.31	0.43
1:A:26:ILE:HD11	1:A:82:GLY:HA2	2.00	0.43
1:A:458:GLN:HE22	1:B:418:ASN:ND2	2.05	0.43
1:A:280:MET:HG3	1:A:435:PHE:CE2	2.53	0.43
1:B:152:ILE:HG22	1:B:153:LEU:N	2.33	0.42
1:B:375:PHE:CE1	1:B:386:PHE:HB2	2.55	0.42
1:A:137:LEU:HD12	1:A:293:TRP:HH2	1.85	0.42
1:A:339:ARG:HD2	1:A:340:ASN:CG	2.45	0.42
1:A:513:LYS:O	1:A:611:LEU:HA	2.20	0.42
1:A:302:TYR:O	1:A:303:LEU:HB2	2.20	0.42
1:B:280:MET:HG3	1:B:435:PHE:CE2	2.54	0.41
1:A:162:PHE:HB2	1:A:163:PRO:HD2	2.02	0.41
1:A:369:ARG:HG2	1:A:392:LEU:HD11	2.02	0.41
1:A:512:TYR:CZ	1:A:613:PRO:HG3	2.55	0.41
1:A:550:GLU:HG2	1:A:570:TYR:CE1	2.56	0.41
1:A:143:GLU:HG3	1:A:144:TYR:N	2.33	0.41
1:A:533:ARG:HD3	3:A:1149:HOH:O	2.20	0.41
1:B:154:ARG:HD2	1:B:293:TRP:CE3	2.55	0.41
1:A:140:GLY:HA3	1:A:293:TRP:O	2.21	0.40
1:A:352:LEU:HB2	1:A:368:ASN:HB2	2.03	0.40
1:A:116:GLU:O	1:A:120:LYS:HG3	2.21	0.40
1:B:483:ALA:HB1	1:B:543:VAL:HB	2.02	0.40
1:B:267:ARG:HH11	1:B:267:ARG:HG2	1.86	0.40
1:A:317:CYS:HB3	1:A:343:CYS:SG	2.61	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	617/638 (97%)	590 (96%)	25 (4%)	2 (0%)	36 26
1	B	617/638 (97%)	593 (96%)	23 (4%)	1 (0%)	43 33
All	All	1234/1276 (97%)	1183 (96%)	48 (4%)	3 (0%)	43 33

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	55	ALA
1	A	143	GLU
1	B	52	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	512/528 (97%)	506 (99%)	6 (1%)	63 54
1	B	512/528 (97%)	507 (99%)	5 (1%)	68 60
All	All	1024/1056 (97%)	1013 (99%)	11 (1%)	65 56

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	268	LEU
1	A	370	ARG
1	A	376	PHE
1	A	490	ARG
1	A	541	LEU
1	A	607	VAL
1	B	36	GLU
1	B	143	GLU
1	B	312	GLU
1	B	358	LEU
1	B	376	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27)

such sidechains are listed below:

Mol	Chain	Res	Type
1	A	24	GLN
1	A	170	HIS
1	A	237	HIS
1	A	334	ASN
1	A	340	ASN
1	A	368	ASN
1	A	418	ASN
1	A	421	GLN
1	A	468	ASN
1	A	507	ASN
1	A	519	GLN
1	B	110	GLN
1	B	203	ASN
1	B	224	GLN
1	B	237	HIS
1	B	306	GLN
1	B	334	ASN
1	B	340	ASN
1	B	418	ASN
1	B	421	GLN
1	B	448	ASN
1	B	458	GLN
1	B	468	ASN
1	B	507	ASN
1	B	515	HIS
1	B	519	GLN
1	B	561	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](#)

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	TPQ	B	382	1	13,14,15	2.39	5 (38%)	13,19,21	1.16	1 (7%)
1	TPQ	A	382	1	13,14,15	2.44	5 (38%)	13,19,21	1.11	1 (7%)

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	TPQ	B	382	1	-	3/5/22/24	0/1/1/1
1	TPQ	A	382	1	-	3/5/22/24	0/1/1/1

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	382	TPQ	O2-C2	4.25	1.35	1.24
1	B	382	TPQ	O2-C2	4.25	1.35	1.24
1	A	382	TPQ	C3-C4	3.69	1.42	1.36
1	A	382	TPQ	O5-C5	3.58	1.34	1.24
1	B	382	TPQ	C3-C4	3.55	1.42	1.36
1	B	382	TPQ	O5-C5	3.53	1.33	1.24
1	A	382	TPQ	C1-C2	-3.27	1.44	1.49
1	B	382	TPQ	C6-C1	3.13	1.42	1.34
1	A	382	TPQ	C6-C1	3.03	1.42	1.34
1	B	382	TPQ	C1-C2	-2.84	1.45	1.49

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	382	TPQ	O2-C2-C3	-2.68	115.74	121.83
1	A	382	TPQ	O2-C2-C3	-2.56	116.00	121.83

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	382	TPQ	N-CA-CB-C1

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Mol	Chain	Res	Type	Atoms
1	A	382	TPQ	O-C-CA-CB
1	B	382	TPQ	N-CA-CB-C1
1	B	382	TPQ	O-C-CA-CB
1	A	382	TPQ	C-CA-CB-C1
1	B	382	TPQ	C-CA-CB-C1

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	B	382	TPQ	1	0
1	A	382	TPQ	1	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 [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](#)

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates [i](#)

EDS was not executed - this section is therefore empty.

6.4 Ligands [i](#)

EDS was not executed - this section is therefore empty.

6.5 Other polymers [i](#)

EDS was not executed - this section is therefore empty.