



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 15, 2026 – 12:50 AM UTC

PDB ID : 1MHY / pdb_00001mhy
Title : METHANE MONOOXYGENASE HYDROXYLASE
Authors : Elango, N.; Radhakrishnan, R.; Froland, W.A.; Waller, B.J.; Earhart, C.A.;
Lipscomb, J.D.; Ohlendorf, D.H.
Deposited on : 1996-10-21
Resolution : 2.00 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
Xtriage (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

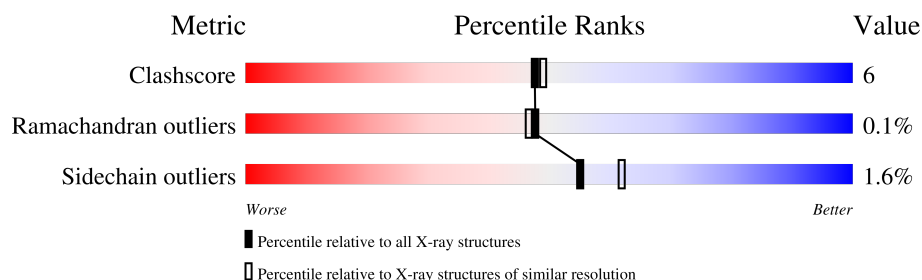
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)

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	B	395	
2	D	521	
3	G	169	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 12729 atoms, of which 3397 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 METHANE MONOOXYGENASE HYDROXYLASE.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	B	383	Total	C	H	N	O	S	0	0	0
			3808	1989	700	537	577	5			

There are 27 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	255	TYR	MET	conflict	UNP P27354
B	256	ASP	ILE	conflict	UNP P27354
B	348	VAL	-	insertion	UNP P27354
B	349	ALA	SER	conflict	UNP P27354
B	350	GLY	ARG	conflict	UNP P27354
B	352	THR	ASP	conflict	UNP P27354
B	353	ASP	ARG	conflict	UNP P27354
B	?	-	ARG	deletion	UNP P27354
B	?	-	ARG	deletion	UNP P27354
B	?	-	ARG	deletion	UNP P27354
B	?	-	LEU	deletion	UNP P27354
B	?	-	ARG	deletion	UNP P27354
B	?	-	GLY	deletion	UNP P27354
B	356	GLY	ALA	conflict	UNP P27354
B	357	VAL	ALA	conflict	UNP P27354
B	359	GLU	SER	conflict	UNP P27354
B	361	LEU	ILE	conflict	UNP P27354
B	362	GLN	GLY	conflict	UNP P27354
B	364	VAL	-	insertion	UNP P27354
B	365	PHE	-	insertion	UNP P27354
B	366	GLY	-	insertion	UNP P27354
B	367	ASP	-	insertion	UNP P27354
B	368	TRP	-	insertion	UNP P27354
B	369	LYS	SER	conflict	UNP P27354
B	371	ASP	-	insertion	UNP P27354
B	372	TYR	THR	conflict	UNP P27354
B	373	ALA	PRO	insertion	UNP P27354

- Molecule 2 is a protein called METHANE MONOOXYGENASE HYDROXYLASE.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	D	510	Total	C	H	N	O	S	0	0	0
			5058	2649	920	720	757	12			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	?	-	ASP	deletion	UNP P27353
D	?	-	ALA	deletion	UNP P27353
D	?	-	LEU	deletion	UNP P27353
D	?	-	LYS	deletion	UNP P27353
D	?	-	VAL	deletion	UNP P27353
D	37	TRP	ARG	conflict	UNP P27353
D	195	GLY	ARG	conflict	UNP P27353
D	209	GLU	ASP	conflict	UNP P27353
D	210	ALA	THR	conflict	UNP P27353
D	225	SER	ILE	conflict	UNP P27353
D	226	ALA	GLY	conflict	UNP P27353
D	329	PRO	-	insertion	UNP P27353
D	331	SER	VAL	conflict	UNP P27353
D	357	GLY	ALA	conflict	UNP P27353

- Molecule 3 is a protein called METHANE MONOOXYGENASE HYDROXYLASE.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
3	G	167	Total	C	H	N	O	S	0	0	0
			1662	868	311	230	252	1			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
G	88	ALA	ARG	conflict	UNP P27355
G	109	GLU	ASP	conflict	UNP P27355
G	110	ALA	GLY	conflict	UNP P27355
G	160	ARG	PRO	conflict	UNP P27355

- Molecule 4 is FE (III) ION (CCD ID: FE) (formula: Fe).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	2	Total	Fe	0	0
			2	2		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	283	Total 849	H 566	O 283	0	0
5	D	331	Total 993	H 662	O 331	0	0
5	G	119	Total 357	H 238	O 119	0	0

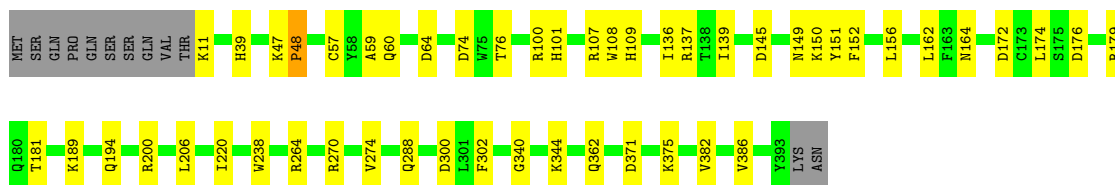
3 Residue-property plots

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. 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.

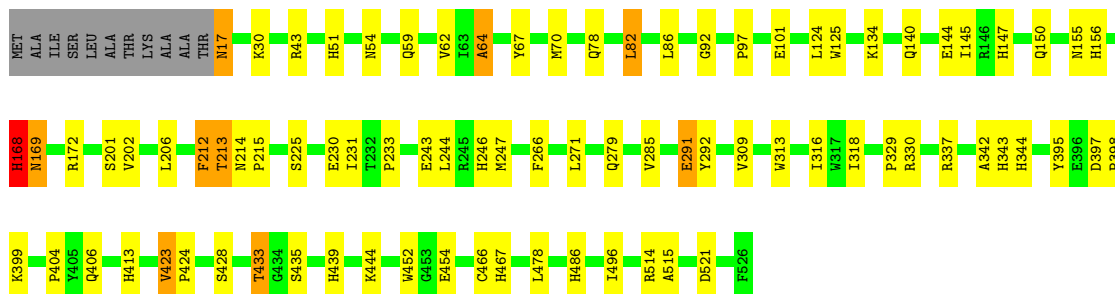
• Molecule 1: METHANE MONOOXYGENASE HYDROXYLASE

Chain B: 



• Molecule 2: METHANE MONOOXYGENASE HYDROXYLASE

Chain D: 



• Molecule 3: METHANE MONOOXYGENASE HYDROXYLASE

Chain G: 



4 Data and refinement statistics

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

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	264.49Å 71.19Å 139.44Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	5.00 – 2.00	Depositor
% Data completeness (in resolution range)	96.8 (5.00-2.00)	Depositor
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.137 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	12729	wwPDB-VP
Average B, all atoms (Å ²)	14.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

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

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	B	0.96	0/3201	1.05	11/4356 (0.3%)
2	D	0.90	0/4267	1.08	23/5797 (0.4%)
3	G	0.85	0/1377	1.05	7/1863 (0.4%)
All	All	0.91	0/8845	1.07	41/12016 (0.3%)

There are no bond length outliers.

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	174	LEU	N-CA-C	9.35	124.24	113.02
1	B	108	TRP	N-CA-C	-7.71	98.52	110.17
2	D	243	GLU	N-CA-C	7.65	120.51	111.71
1	B	145	ASP	N-CA-C	7.35	119.37	111.36
1	B	152	PHE	N-CA-C	-7.05	103.67	111.36
2	D	285	VAL	N-CA-C	7.00	117.79	110.72
2	D	309	VAL	N-CA-C	6.96	117.08	110.74
2	D	168	HIS	N-CA-C	-6.70	104.05	111.82
2	D	478	LEU	N-CA-C	6.50	118.91	111.11
3	G	84	THR	N-CA-C	-6.48	105.34	113.18
3	G	119	TYR	N-CA-C	6.43	121.92	113.30
2	D	213	THR	N-CA-C	6.30	117.81	111.07
1	B	172	ASP	N-CA-C	6.21	120.93	113.23
2	D	433	THR	N-CA-C	6.14	120.91	113.17
3	G	82	ALA	N-CA-C	-6.11	101.47	110.52
3	G	40	TYR	N-CA-C	6.11	122.62	114.12
1	B	109	HIS	N-CA-C	5.94	117.76	111.28
2	D	145	ILE	N-CA-C	-5.90	104.76	110.72
2	D	496	ILE	N-CA-C	-5.84	105.04	110.53
3	G	7	ILE	N-CA-C	5.81	115.93	110.30
2	D	423	VAL	CA-C-N	5.75	125.66	119.85

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	423	VAL	C-N-CA	5.75	125.66	119.85
1	B	100	ARG	N-CA-C	5.75	118.32	111.71
2	D	266	PHE	N-CA-C	5.69	121.09	114.04
3	G	126	VAL	N-CA-C	5.55	116.19	110.36
3	G	8	HIS	N-CA-C	5.54	120.10	113.23
1	B	340	GLY	N-CA-C	-5.51	108.61	114.67
1	B	59	ALA	N-CA-C	-5.50	105.44	111.82
2	D	82	LEU	N-CA-C	5.45	117.02	111.14
2	D	247	MET	N-CA-C	-5.35	105.53	111.36
2	D	521	ASP	N-CA-C	-5.32	97.99	109.01
1	B	302	PHE	N-CA-C	5.32	118.23	111.69
2	D	329	PRO	N-CA-C	5.31	119.21	111.03
2	D	64	ALA	N-CA-C	5.26	117.70	111.33
2	D	212	PHE	N-CA-C	5.21	119.31	111.96
2	D	291	GLU	N-CA-C	5.12	116.95	111.36
2	D	316	ILE	N-CA-C	5.11	115.33	110.53
2	D	125	TRP	N-CA-C	-5.08	105.74	111.28
2	D	225	SER	N-CA-CB	5.06	117.56	110.12
1	B	107	ARG	N-CA-C	5.05	116.53	108.41
2	D	342	ALA	N-CA-C	5.00	116.73	111.28

There are no chirality outliers.

There are no planarity outliers.

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	B	3108	700	2942	37	1
2	D	4138	920	3934	56	0
3	G	1351	311	1387	14	0
4	D	2	0	0	0	0
5	B	283	566	0	4	1
5	D	331	662	0	8	0
5	G	119	238	0	0	0
All	All	9332	3397	8263	95	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 6.

All (95) 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:47:LYS:HB2	1:B:48:PRO:HD3	1.48	0.96
2:D:78:GLN:HE22	2:D:150:GLN:HE21	1.13	0.90
1:B:194:GLN:HE22	2:D:64:ALA:H	1.26	0.82
2:D:155:ASN:HD22	2:D:168:HIS:HD2	1.31	0.79
2:D:202:VAL:HG11	2:D:271:LEU:HA	1.65	0.78
2:D:337:ARG:HG2	2:D:433:THR:HG21	1.67	0.77
2:D:292:TYR:OH	2:D:344:HIS:HD2	1.71	0.74
1:B:264:ARG:HE	1:B:288:GLN:HE22	1.32	0.73
3:G:24:THR:H	3:G:28:GLN:HE22	1.39	0.70
1:B:39:HIS:HE1	2:D:172:ARG:H	1.36	0.69
1:B:149:ASN:HD21	1:B:200:ARG:HH21	1.42	0.68
2:D:439:HIS:HE1	2:D:454:GLU:OE2	1.78	0.67
1:B:149:ASN:ND2	1:B:200:ARG:HH21	1.95	0.65
1:B:74:ASP:H	2:D:467:HIS:HD2	1.43	0.64
1:B:64:ASP:OD2	3:G:8:HIS:HD2	1.82	0.62
1:B:149:ASN:HD22	1:B:200:ARG:HE	1.48	0.62
2:D:78:GLN:NE2	2:D:150:GLN:HE21	1.93	0.61
2:D:330:ARG:HD3	2:D:452:TRP:CZ3	2.36	0.61
2:D:156:HIS:HD2	5:D:1012:HOH:O	1.82	0.60
1:B:47:LYS:CB	1:B:48:PRO:HD3	2.29	0.60
1:B:101:HIS:HE1	1:B:181:THR:OG1	1.83	0.60
2:D:54:ASN:HB2	5:D:1133:HOH:O	2.03	0.58
2:D:140:GLN:O	2:D:144:GLU:HG2	2.05	0.56
1:B:151:TYR:OH	1:B:344:LYS:HE3	2.05	0.56
1:B:57:CYS:HB3	5:B:576:HOH:O	2.05	0.56
3:G:24:THR:H	3:G:28:GLN:NE2	2.02	0.56
2:D:213:THR:HG21	5:D:807:HOH:O	2.07	0.55
2:D:514:ARG:HH11	2:D:514:ARG:HG3	1.71	0.54
2:D:330:ARG:HD3	2:D:452:TRP:CH2	2.42	0.54
1:B:149:ASN:HD21	1:B:200:ARG:NH2	2.06	0.53
2:D:444:LYS:HD3	3:G:48:TYR:CD2	2.44	0.52
2:D:413:HIS:HD2	2:D:428:SER:OG	1.93	0.52
1:B:39:HIS:CE1	2:D:172:ARG:H	2.23	0.51
2:D:212:PHE:O	2:D:215:PRO:HD2	2.10	0.51
1:B:149:ASN:ND2	1:B:200:ARG:HE	2.09	0.51
3:G:114:GLU:HA	3:G:117:GLN:HE21	1.75	0.51
1:B:194:GLN:NE2	2:D:64:ALA:H	2.04	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:467:HIS:HE1	3:G:51:ASP:O	1.95	0.50
1:B:371:ASP:O	1:B:375:LYS:HD2	2.13	0.49
1:B:164:ASN:HB3	1:B:238:TRP:CE2	2.47	0.49
2:D:147:HIS:CE1	5:D:803:HOH:O	2.66	0.49
2:D:214:ASN:HB3	2:D:215:PRO:HD3	1.94	0.49
2:D:92:GLY:HA2	2:D:231:ILE:HD11	1.94	0.49
2:D:82:LEU:HD23	2:D:86:LEU:HD12	1.95	0.48
1:B:60:GLN:HE22	1:B:179:ARG:HH12	1.61	0.48
2:D:406:GLN:NE2	5:D:1042:HOH:O	2.47	0.48
2:D:246:HIS:CD2	2:D:246:HIS:N	2.78	0.47
2:D:144:GLU:HA	2:D:144:GLU:OE1	2.15	0.47
2:D:486:HIS:HD2	5:D:982:HOH:O	1.97	0.47
1:B:382:VAL:O	1:B:386:VAL:HG23	2.15	0.46
2:D:155:ASN:HD22	2:D:168:HIS:CD2	2.22	0.46
1:B:189:LYS:HD3	1:B:189:LYS:HA	1.72	0.46
1:B:270:ARG:HD3	1:B:270:ARG:C	2.41	0.46
1:B:164:ASN:HD21	2:D:134:LYS:HD3	1.82	0.45
2:D:439:HIS:CE1	2:D:454:GLU:OE2	2.64	0.45
3:G:69:LYS:O	3:G:73:LEU:HB2	2.17	0.45
1:B:76:THR:HA	2:D:466:CYS:HB2	1.99	0.45
2:D:230:GLU:C	2:D:233:PRO:HD2	2.41	0.45
2:D:343:HIS:H	2:D:343:HIS:CD2	2.35	0.45
1:B:11:LYS:HG2	5:B:641:HOH:O	2.16	0.45
2:D:313:TRP:CZ2	2:D:318:ILE:HD11	2.53	0.44
1:B:136:ILE:O	1:B:139:ILE:HG22	2.17	0.44
1:B:156:LEU:C	1:B:156:LEU:HD12	2.42	0.44
1:B:74:ASP:H	2:D:467:HIS:CD2	2.31	0.43
1:B:150:LYS:HB3	1:B:220:ILE:HD13	2.00	0.43
3:G:5:GLU:HB3	3:G:6:PRO:CD	2.47	0.43
3:G:154:GLU:H	3:G:154:GLU:CD	2.25	0.43
2:D:17:ASN:ND2	2:D:17:ASN:N	2.67	0.43
1:B:150:LYS:HB3	1:B:220:ILE:CD1	2.48	0.43
2:D:124:LEU:HD21	2:D:201:SER:HB2	1.99	0.43
2:D:397:ASP:HA	2:D:398:PRO:HD3	1.96	0.43
2:D:51:HIS:HB2	5:D:1020:HOH:O	2.19	0.43
2:D:97:PRO:O	2:D:101:GLU:HG2	2.19	0.43
2:D:30:LYS:HB2	2:D:30:LYS:HE3	1.77	0.42
2:D:423:VAL:HA	2:D:424:PRO:HD3	1.83	0.42
1:B:39:HIS:CE1	2:D:172:ARG:HB2	2.54	0.42
3:G:74:PRO:HG2	3:G:77:ASP:CG	2.45	0.42
2:D:292:TYR:OH	2:D:344:HIS:CD2	2.60	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:395:TYR:CE1	2:D:404:PRO:HG2	2.55	0.42
1:B:238:TRP:CD1	1:B:238:TRP:C	2.97	0.42
2:D:67:TYR:HA	2:D:70:MET:HE2	2.01	0.42
1:B:101:HIS:HD2	1:B:300:ASP:OD2	2.02	0.41
5:B:441:HOH:O	3:G:8:HIS:CE1	2.73	0.41
2:D:291:GLU:OE1	2:D:343:HIS:HE1	2.02	0.41
2:D:330:ARG:HD3	2:D:452:TRP:HZ3	1.84	0.41
1:B:362:GLN:NE2	1:B:386:VAL:HG21	2.35	0.41
2:D:59:GLN:HG3	5:D:1017:HOH:O	2.20	0.41
1:B:264:ARG:NE	1:B:288:GLN:HE22	2.08	0.41
3:G:163:LYS:HB2	3:G:163:LYS:HE3	1.64	0.41
2:D:399:LYS:HA	2:D:399:LYS:HD3	1.81	0.41
1:B:162:LEU:HD23	1:B:162:LEU:HA	1.89	0.41
2:D:406:GLN:NE2	2:D:515:ALA:HB1	2.36	0.41
5:B:441:HOH:O	3:G:8:HIS:HE1	2.04	0.41
2:D:169:ASN:C	2:D:169:ASN:HD22	2.29	0.41
3:G:62:GLU:HG2	3:G:122:PRO:HD3	2.03	0.41

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)
1:B:137:ARG:HH21	5:B:520:HOH:H2[4_575]	1.17	0.43

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	B	381/395 (96%)	368 (97%)	12 (3%)	1 (0%)	36	35
2	D	508/521 (98%)	490 (96%)	18 (4%)	0	100	100
3	G	165/169 (98%)	161 (98%)	4 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	1054/1085 (97%)	1019 (97%)	34 (3%)	1 (0%)	48 46

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	48	PRO

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	B	315/327 (96%)	312 (99%)	3 (1%)	68 75
2	D	423/430 (98%)	414 (98%)	9 (2%)	47 52
3	G	144/145 (99%)	142 (99%)	2 (1%)	59 66
All	All	882/902 (98%)	868 (98%)	14 (2%)	55 62

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	176	ASP
1	B	206	LEU
1	B	274	VAL
2	D	17	ASN
2	D	43	ARG
2	D	62	VAL
2	D	168	HIS
2	D	169	ASN
2	D	206	LEU
2	D	244	LEU
2	D	279	GLN
2	D	435	SER
3	G	27	ASP
3	G	73	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (33) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	39	HIS
1	B	60	GLN
1	B	101	HIS
1	B	149	ASN
1	B	164	ASN
1	B	192	ASN
1	B	194	GLN
1	B	288	GLN
1	B	384	GLN
2	D	78	GLN
2	D	156	HIS
2	D	168	HIS
2	D	169	ASN
2	D	203	ASN
2	D	214	ASN
2	D	273	ASN
2	D	278	GLN
2	D	279	GLN
2	D	343	HIS
2	D	344	HIS
2	D	369	GLN
2	D	388	ASN
2	D	406	GLN
2	D	413	HIS
2	D	422	GLN
2	D	439	HIS
2	D	456	GLN
2	D	467	HIS
2	D	486	HIS
3	G	8	HIS
3	G	28	GLN
3	G	55	GLN
3	G	117	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.