



Full wwPDB X-ray Structure Validation Report ⓘ

Dec 13, 2016 – 12:11 PM EST

PDB ID : 5M10
Title : Crystal structure of cyclohexanone monooxygenase from *Thermocrispum municipale* in the oxidised state with a bound nicotinamide.
Authors : Gomez-Castellanos, J.R.; Mattevi, A.
Deposited on : 2016-10-06
Resolution : 1.22 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20028442
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20028442

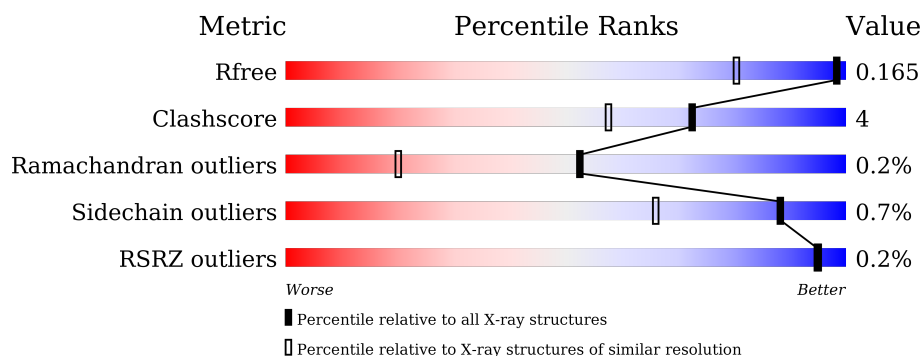
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.22 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	1197 (1.26-1.18)
Clashscore	102246	1295 (1.26-1.18)
Ramachandran outliers	100387	1239 (1.26-1.18)
Sidechain outliers	100360	1237 (1.26-1.18)
RSRZ outliers	91569	1201 (1.26-1.18)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	541	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	TRS	A	608	-	-	-	X
6	GOL	A	609	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
6	GOL	A	610	-	-	-	X

2 Entry composition [i](#)

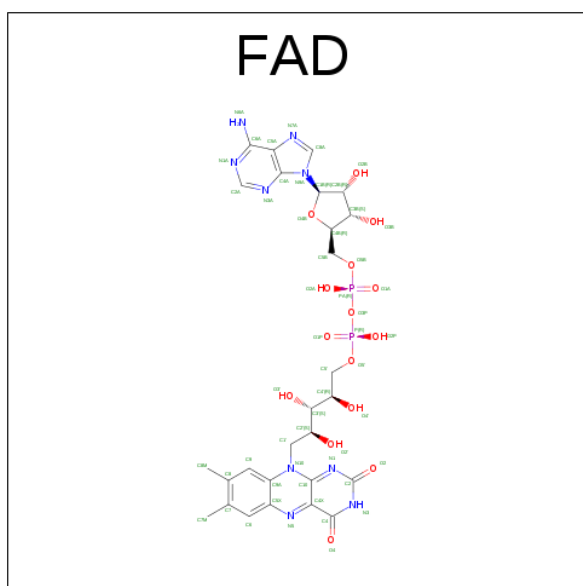
There are 7 unique types of molecules in this entry. The entry contains 5034 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Cyclohexanone Monooxygenase from *Thermocrispum municipale*.

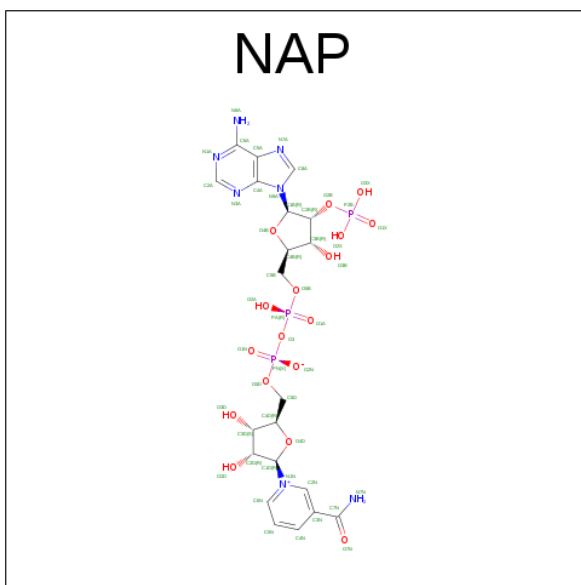
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	529	4315	2724	741	835	15	0	24	0

- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).



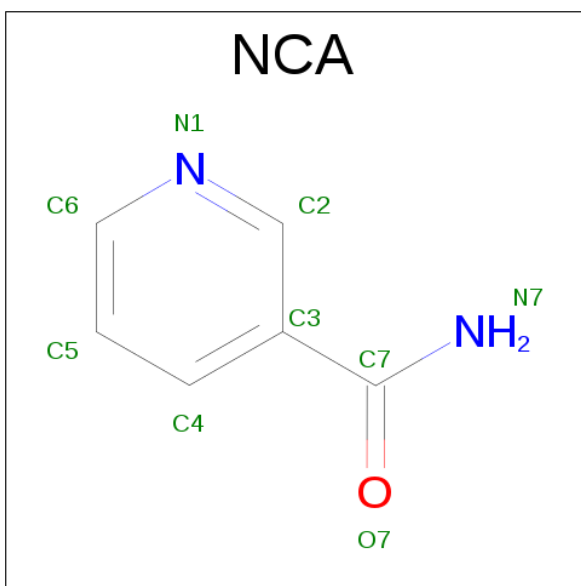
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
2	A	1	53	27	9	15	2	0	0

- Molecule 3 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NAP) (formula: $C_{21}H_{28}N_7O_{17}P_3$).



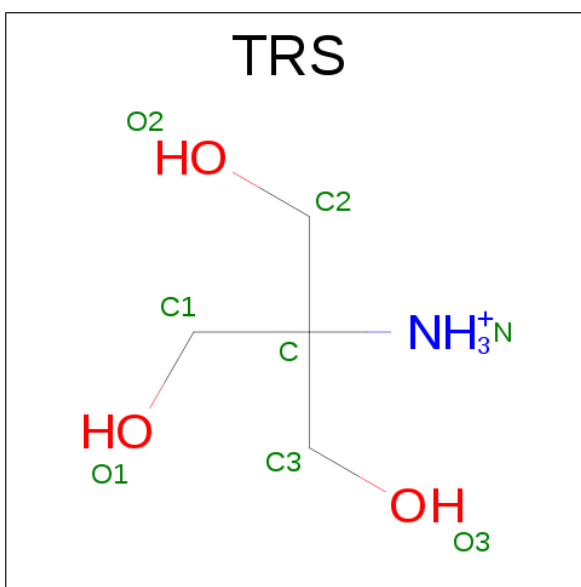
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	
			48	21	7	17	3	

- Molecule 4 is NICOTINAMIDE (three-letter code: NCA) (formula: $C_6H_6N_2O$).



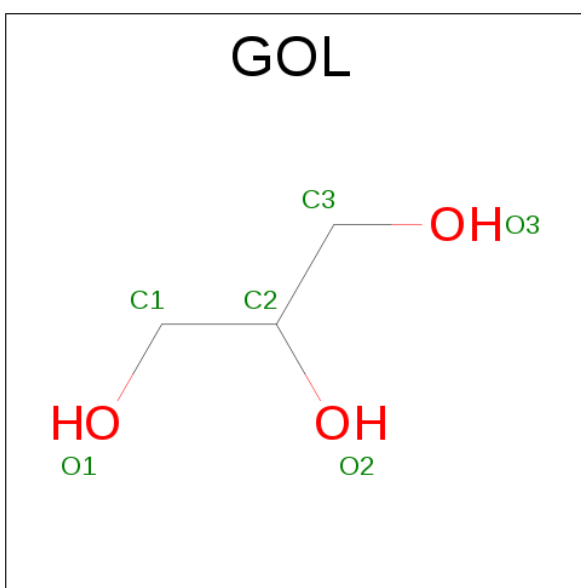
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O		
			9	6	2	1		

- Molecule 5 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (three-letter code: TRS) (formula: $C_4H_{12}NO_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			8	4	1	3		
5	A	1	Total	C	N	O	0	0
			8	4	1	3		

- Molecule 6 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			6	3	3		
6	A	1	Total	C	O	0	0
			6	3	3		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			6	3	3		
6	A	1	Total	C	O	0	0
			6	3	3		
6	A	1	Total	C	O	0	0
			6	3	3		

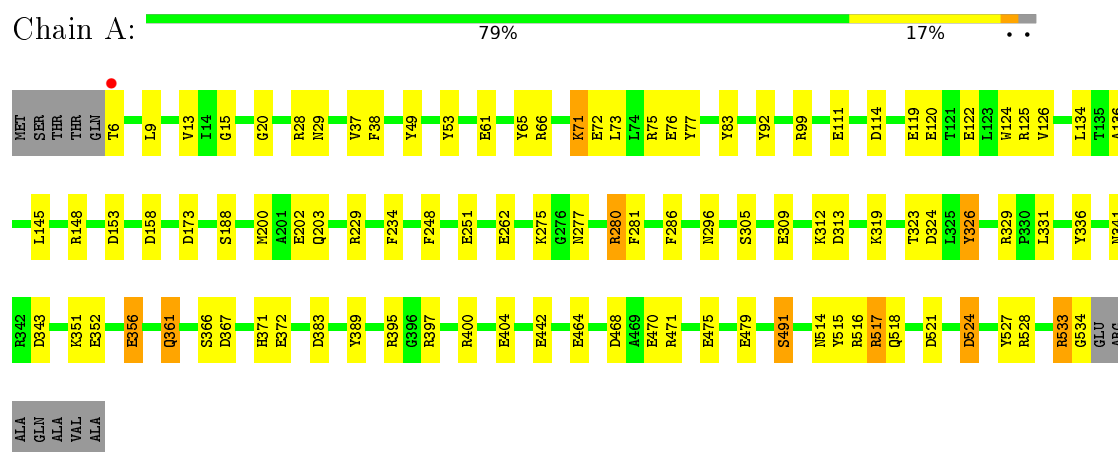
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	563	Total	O	0	0
			563	563		

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: Cyclohexanone Monooxygenase from *Thermocrispum municipale*



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	68.84Å 113.63Å 155.48Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.90 – 1.22 38.90 – 1.22	Depositor EDS
% Data completeness (in resolution range)	93.6 (38.90-1.22) 93.6 (38.90-1.22)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.62 (at 1.22Å)	Xtriage
Refinement program	REFMAC 5.8.0155	Depositor
R, R_{free}	0.131 , 0.165 0.130 , 0.165	Depositor DCC
R_{free} test set	5115 reflections (3.13%)	DCC
Wilson B-factor (Å ²)	9.6	Xtriage
Anisotropy	1.013	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 44.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.015 for 1/2*h-1/2*k,-3/2*h-1/2*k,-l 0.025 for 1/2*h+1/2*k,3/2*h-1/2*k,-l	Xtriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	5034	wwPDB-VP
Average B, all atoms (Å ²)	16.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.13% 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: GOL, NAP, TRS, FAD, NCA

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	1.68	60/4447 (1.3%)	1.72	73/6037 (1.2%)

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	A	0	1

All (60) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	76	GLU	CD-OE2	-17.32	1.06	1.25
1	A	71	LYS	CE-NZ	14.26	1.84	1.49
1	A	404	GLU	CD-OE2	12.87	1.39	1.25
1	A	352	GLU	CD-OE1	11.69	1.38	1.25
1	A	534	GLY	N-CA	11.29	1.62	1.46
1	A	356	GLU	CD-OE2	-9.88	1.14	1.25
1	A	125[A]	ARG	CZ-NH1	9.41	1.45	1.33
1	A	125[B]	ARG	CZ-NH1	9.41	1.45	1.33
1	A	251	GLU	CG-CD	9.30	1.66	1.51
1	A	119	GLU	CD-OE2	8.01	1.34	1.25
1	A	475	GLU	CD-OE2	7.74	1.34	1.25
1	A	203	GLN	CD-NE2	7.59	1.51	1.32
1	A	479	GLU	CD-OE2	-7.51	1.17	1.25
1	A	491[A]	SER	CA-CB	7.34	1.64	1.52
1	A	491[B]	SER	CA-CB	7.34	1.64	1.52
1	A	83	TYR	CE1-CZ	-7.17	1.29	1.38
1	A	248	PHE	CG-CD1	-6.86	1.28	1.38
1	A	313	ASP	CG-OD2	-6.82	1.09	1.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	122	GLU	CD-OE1	-6.72	1.18	1.25
1	A	372	GLU	CD-OE2	-6.64	1.18	1.25
1	A	72	GLU	CD-OE2	6.63	1.32	1.25
1	A	528	ARG	CD-NE	-6.50	1.35	1.46
1	A	521	ASP	CB-CG	6.18	1.64	1.51
1	A	125[A]	ARG	CD-NE	6.14	1.56	1.46
1	A	125[B]	ARG	CD-NE	6.14	1.56	1.46
1	A	72	GLU	CB-CG	-6.11	1.40	1.52
1	A	72	GLU	CD-OE1	-6.10	1.19	1.25
1	A	389	TYR	CG-CD1	-6.04	1.31	1.39
1	A	38	PHE	CG-CD2	-6.01	1.29	1.38
1	A	125[A]	ARG	CZ-NH2	5.99	1.40	1.33
1	A	125[B]	ARG	CZ-NH2	5.99	1.40	1.33
1	A	49	TYR	CE2-CZ	-5.99	1.30	1.38
1	A	83	TYR	CG-CD2	-5.91	1.31	1.39
1	A	395	ARG	CZ-NH1	5.80	1.40	1.33
1	A	148	ARG	CZ-NH1	5.78	1.40	1.33
1	A	202[A]	GLU	CD-OE2	5.76	1.31	1.25
1	A	202[B]	GLU	CD-OE2	5.76	1.31	1.25
1	A	404	GLU	CD-OE1	5.61	1.31	1.25
1	A	111	GLU	CD-OE2	-5.59	1.19	1.25
1	A	336	TYR	CG-CD1	5.58	1.46	1.39
1	A	319	LYS	CD-CE	-5.57	1.37	1.51
1	A	92	TYR	CE1-CZ	-5.52	1.31	1.38
1	A	251	GLU	CD-OE1	-5.52	1.19	1.25
1	A	262	GLU	CD-OE2	-5.51	1.19	1.25
1	A	124	TRP	CG-CD1	5.50	1.44	1.36
1	A	491[A]	SER	CB-OG	5.43	1.49	1.42
1	A	491[B]	SER	CB-OG	5.43	1.49	1.42
1	A	281	PHE	CG-CD2	-5.42	1.30	1.38
1	A	77	TYR	CE1-CZ	5.32	1.45	1.38
1	A	120[A]	GLU	CD-OE1	-5.29	1.19	1.25
1	A	120[B]	GLU	CD-OE1	-5.29	1.19	1.25
1	A	442	GLU	CD-OE1	5.27	1.31	1.25
1	A	203	GLN	CG-CD	-5.24	1.39	1.51
1	A	286	PHE	CG-CD1	5.21	1.46	1.38
1	A	65	TYR	CE2-CZ	-5.17	1.31	1.38
1	A	145	LEU	CB-CG	-5.13	1.37	1.52
1	A	126	VAL	CB-CG1	-5.12	1.42	1.52
1	A	352	GLU	CD-OE2	-5.11	1.20	1.25
1	A	119	GLU	CD-OE1	-5.11	1.20	1.25
1	A	336	TYR	CE2-CZ	5.10	1.45	1.38

All (73) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	125[A]	ARG	NE-CZ-NH2	-33.18	103.71	120.30
1	A	125[B]	ARG	NE-CZ-NH2	-33.18	103.71	120.30
1	A	524[A]	ASP	CB-CG-OD1	23.37	139.33	118.30
1	A	524[B]	ASP	CB-CG-OD1	23.37	139.33	118.30
1	A	125[A]	ARG	NE-CZ-NH1	22.18	131.39	120.30
1	A	125[B]	ARG	NE-CZ-NH1	22.18	131.39	120.30
1	A	148	ARG	NE-CZ-NH1	13.33	126.97	120.30
1	A	528	ARG	NE-CZ-NH2	-13.28	113.66	120.30
1	A	395	ARG	NE-CZ-NH2	-10.94	114.83	120.30
1	A	61	GLU	OE1-CD-OE2	-9.73	111.62	123.30
1	A	479	GLU	OE1-CD-OE2	9.48	134.68	123.30
1	A	521	ASP	CB-CG-OD2	-9.08	110.13	118.30
1	A	524[A]	ASP	CB-CG-OD2	-8.84	110.35	118.30
1	A	524[B]	ASP	CB-CG-OD2	-8.84	110.35	118.30
1	A	153	ASP	CB-CG-OD1	8.79	126.21	118.30
1	A	200	MET	CG-SD-CE	-8.75	86.20	100.20
1	A	309	GLU	OE1-CD-OE2	8.23	133.18	123.30
1	A	145	LEU	CB-CG-CD1	8.19	124.92	111.00
1	A	313	ASP	CB-CG-OD1	8.12	125.61	118.30
1	A	471	ARG	NE-CZ-NH2	-7.85	116.38	120.30
1	A	280[A]	ARG	NE-CZ-NH1	7.55	124.08	120.30
1	A	280[B]	ARG	NE-CZ-NH1	7.55	124.08	120.30
1	A	280[A]	ARG	NE-CZ-NH2	7.44	124.02	120.30
1	A	280[B]	ARG	NE-CZ-NH2	7.44	124.02	120.30
1	A	275[A]	LYS	CD-CE-NZ	-7.40	94.69	111.70
1	A	275[B]	LYS	CD-CE-NZ	-7.40	94.69	111.70
1	A	400	ARG	NE-CZ-NH1	7.39	123.99	120.30
1	A	367	ASP	CB-CG-OD1	6.97	124.57	118.30
1	A	280[A]	ARG	NH1-CZ-NH2	-6.93	111.77	119.40
1	A	280[B]	ARG	NH1-CZ-NH2	-6.93	111.77	119.40
1	A	524[A]	ASP	OD1-CG-OD2	-6.84	110.30	123.30
1	A	524[B]	ASP	OD1-CG-OD2	-6.84	110.30	123.30
1	A	336	TYR	CD1-CE1-CZ	6.75	125.88	119.80
1	A	99	ARG	NE-CZ-NH2	-6.71	116.95	120.30
1	A	517	ARG	NE-CZ-NH1	-6.68	116.96	120.30
1	A	336	TYR	CB-CG-CD2	6.65	124.99	121.00
1	A	533	ARG	CB-CG-CD	6.61	128.80	111.60
1	A	372	GLU	OE1-CD-OE2	6.61	131.23	123.30
1	A	343	ASP	CB-CG-OD2	-6.54	112.41	118.30
1	A	521	ASP	OD1-CG-OD2	6.46	135.57	123.30
1	A	229	ARG	NE-CZ-NH2	-6.45	117.07	120.30
1	A	468	ASP	CB-CG-OD1	6.44	124.10	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	173	ASP	CB-CG-OD1	6.43	124.09	118.30
1	A	326	TYR	CD1-CE1-CZ	6.26	125.43	119.80
1	A	533	ARG	NE-CZ-NH1	6.19	123.39	120.30
1	A	158	ASP	CB-CG-OD1	6.13	123.82	118.30
1	A	124	TRP	CG-CD1-NE1	-6.04	104.06	110.10
1	A	125[A]	ARG	CD-NE-CZ	5.96	131.95	123.60
1	A	125[B]	ARG	CD-NE-CZ	5.96	131.95	123.60
1	A	324	ASP	CB-CG-OD1	5.78	123.50	118.30
1	A	528	ARG	CD-NE-CZ	5.76	131.66	123.60
1	A	124	TRP	CD1-NE1-CE2	5.75	114.17	109.00
1	A	352	GLU	CG-CD-OE1	5.73	129.76	118.30
1	A	352	GLU	CG-CD-OE2	-5.65	107.00	118.30
1	A	71	LYS	CG-CD-CE	5.58	128.63	111.90
1	A	383	ASP	CB-CG-OD1	5.47	123.22	118.30
1	A	53	TYR	CG-CD2-CE2	5.39	125.61	121.30
1	A	471	ARG	CG-CD-NE	-5.33	100.61	111.80
1	A	528	ARG	NH1-CZ-NH2	5.32	125.25	119.40
1	A	158	ASP	CB-CG-OD2	-5.30	113.53	118.30
1	A	527	TYR	CD1-CE1-CZ	5.27	124.54	119.80
1	A	75	ARG	NE-CZ-NH1	5.22	122.91	120.30
1	A	527	TYR	CB-CG-CD1	-5.20	117.88	121.00
1	A	323	THR	OG1-CB-CG2	-5.17	98.11	110.00
1	A	470	GLU	OE1-CD-OE2	5.17	129.50	123.30
1	A	356	GLU	CG-CD-OE1	5.16	128.62	118.30
1	A	397	ARG	NE-CZ-NH2	5.14	122.87	120.30
1	A	114	ASP	CB-CG-OD2	-5.13	113.68	118.30
1	A	516	ARG	CD-NE-CZ	5.11	130.75	123.60
1	A	515	TYR	CB-CG-CD2	-5.11	117.94	121.00
1	A	99	ARG	NE-CZ-NH1	5.09	122.84	120.30
1	A	134	LEU	CB-CG-CD1	5.09	119.65	111.00
1	A	527	TYR	CB-CG-CD2	5.05	124.03	121.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	66	ARG	Sidechain

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	4315	0	4175	35	0
2	A	53	0	30	1	0
3	A	48	0	25	1	0
4	A	9	0	6	2	0
5	A	16	0	23	0	0
6	A	30	0	40	3	0
7	A	563	0	0	13	2
All	All	5034	0	4299	37	2

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 4.

All (37) 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:71:LYS:NZ	1:A:71:LYS:CE	1.84	1.38
1:A:464[B]:GLU:OE2	7:A:701:HOH:O	1.54	1.21
1:A:524[A]:ASP:OD2	7:A:702:HOH:O	1.63	1.14
1:A:491[A]:SER:O	7:A:703:HOH:O	1.85	0.93
1:A:188[B]:SER:OG	1:A:331:LEU:HD22	1.70	0.90
1:A:28:ARG:HH21	1:A:29:ASN:HD21	1.22	0.88
1:A:280[A]:ARG:NH1	6:A:610:GOL:O2	2.11	0.81
1:A:329:ARG:HH11	4:A:603:NCA:HN72	1.27	0.81
1:A:305[B]:SER:OG	7:A:704:HOH:O	1.99	0.80
1:A:296:ASN:HD21	1:A:326:TYR:H	1.31	0.79
1:A:361:GLN:H	1:A:361:GLN:HE21	1.30	0.77
1:A:464[A]:GLU:OE2	7:A:705:HOH:O	2.04	0.74
1:A:312:LYS:H	1:A:341:ASN:HD21	1.34	0.74
1:A:188[B]:SER:OG	1:A:331:LEU:CD2	2.39	0.70
1:A:514[A]:ASN:OD1	1:A:517:ARG:NH2	2.28	0.66
1:A:188[B]:SER:HG	1:A:331:LEU:HD22	1.62	0.64
1:A:518[B]:GLN:NE2	7:A:708:HOH:O	2.30	0.61
1:A:464[B]:GLU:CD	7:A:701:HOH:O	2.14	0.60
1:A:356:GLU:HG3	1:A:366:SER:HA	1.85	0.56
1:A:371:HIS:HD2	7:A:926:HOH:O	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:601:FAD:HM73	3:A:602:NAP:C5N	2.37	0.54
1:A:533:ARG:NH1	7:A:701:HOH:O	2.37	0.54
1:A:312:LYS:H	1:A:341:ASN:ND2	2.06	0.51
1:A:361:GLN:N	1:A:361:GLN:HE21	2.04	0.50
1:A:329:ARG:NH1	4:A:603:NCA:HN72	2.03	0.48
1:A:28:ARG:HH21	1:A:29:ASN:ND2	2.02	0.47
1:A:73:LEU:HD23	7:A:1109:HOH:O	2.16	0.46
1:A:464[A]:GLU:HG2	7:A:705:HOH:O	2.15	0.46
1:A:9:LEU:O	1:A:136:ALA:HA	2.16	0.46
1:A:13:VAL:HB	1:A:37:VAL:HG22	2.00	0.43
6:A:610:GOL:H12	7:A:856:HOH:O	2.18	0.43
1:A:280[A]:ARG:HH11	6:A:610:GOL:C2	2.28	0.43
1:A:351:LYS:HE3	7:A:1185:HOH:O	2.18	0.42
1:A:13:VAL:O	1:A:37:VAL:HA	2.20	0.42
1:A:15:GLY:O	1:A:20:GLY:HA3	2.20	0.41
1:A:296:ASN:HD21	1:A:326:TYR:N	2.07	0.41

All (2) 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)
7:A:1243:HOH:O	7:A:1243:HOH:O[4_545]	2.03	0.17
7:A:718:HOH:O	7:A:1181:HOH:O[8_445]	2.07	0.13

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	552/541 (102%)	541 (98%)	10 (2%)	1 (0%)	52 20

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	277	ASN

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	460/444 (104%)	457 (99%)	3 (1%)	88 66

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	THR
1	A	234	PHE
1	A	361	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	29	ASN
1	A	109	ASN
1	A	203	GLN
1	A	296	ASN
1	A	341	ASN
1	A	361	GLN
1	A	371	HIS
1	A	405	HIS
1	A	482	ASN

5.3.3 RNA [i](#)

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

10 ligands 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$
2	FAD	A	601	-	52,58,58	2.09	16 (30%)	52,89,89	3.32	13 (25%)
3	NAP	A	602	-	45,52,52	1.36	6 (13%)	55,80,80	1.20	2 (3%)
4	NCA	A	603	-	9,9,9	1.52	1 (11%)	11,11,11	3.03	5 (45%)
5	TRS	A	604	-	7,7,7	2.45	3 (42%)	9,9,9	2.65	6 (66%)
6	GOL	A	605	-	5,5,5	0.38	0	5,5,5	0.80	0
6	GOL	A	606	-	5,5,5	1.04	1 (20%)	5,5,5	0.59	0
6	GOL	A	607	-	5,5,5	1.42	1 (20%)	5,5,5	1.09	0
5	TRS	A	608	-	7,7,7	1.92	3 (42%)	9,9,9	7.02	5 (55%)
6	GOL	A	609	-	5,5,5	1.99	1 (20%)	5,5,5	1.24	1 (20%)
6	GOL	A	610	-	5,5,5	1.21	0	5,5,5	1.94	1 (20%)

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
2	FAD	A	601	-	-	0/30/50/50	0/6/6/6
3	NAP	A	602	-	-	0/27/67/67	0/5/5/5
4	NCA	A	603	-	-	0/4/4/4	0/1/1/1
5	TRS	A	604	-	-	0/9/9/9	0/0/0/0
6	GOL	A	605	-	-	0/4/4/4	0/0/0/0
6	GOL	A	606	-	-	0/4/4/4	0/0/0/0
6	GOL	A	607	-	-	0/4/4/4	0/0/0/0
5	TRS	A	608	-	-	0/9/9/9	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	GOL	A	609	-	-	0/4/4/4	0/0/0/0
6	GOL	A	610	-	-	0/4/4/4	0/0/0/0

All (32) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	602	NAP	C6N-C5N	-3.99	1.29	1.38
2	A	601	FAD	C5A-C4A	-3.41	1.32	1.40
4	A	603	NCA	O7-C7	-3.15	1.17	1.24
2	A	601	FAD	C4-N3	-3.15	1.27	1.33
2	A	601	FAD	C1'-N10	-2.91	1.45	1.48
3	A	602	NAP	O7N-C7N	-2.88	1.18	1.24
2	A	601	FAD	PA-O1A	-2.77	1.41	1.51
2	A	601	FAD	C10-N1	-2.73	1.31	1.35
2	A	601	FAD	C2A-N1A	-2.56	1.29	1.33
2	A	601	FAD	C4X-C10	-2.52	1.36	1.40
2	A	601	FAD	C2B-C1B	-2.24	1.50	1.53
5	A	604	TRS	O1-C1	-2.18	1.35	1.42
6	A	606	GOL	O2-C2	2.02	1.49	1.43
2	A	601	FAD	C2-N1	2.17	1.42	1.38
3	A	602	NAP	C8A-N7A	2.17	1.38	1.34
5	A	608	TRS	C1-C	2.19	1.56	1.53
2	A	601	FAD	C6A-C5A	2.22	1.54	1.42
5	A	608	TRS	O1-C1	2.29	1.49	1.42
3	A	602	NAP	C3D-C4D	2.32	1.59	1.53
5	A	604	TRS	O3-C3	2.38	1.49	1.42
6	A	607	GOL	O1-C1	2.67	1.53	1.42
3	A	602	NAP	C5A-C4A	2.73	1.46	1.40
3	A	602	NAP	O4D-C1D	2.77	1.45	1.41
2	A	601	FAD	C4X-N5	2.92	1.37	1.33
5	A	608	TRS	C-N	3.29	1.55	1.50
2	A	601	FAD	O3'-C3'	3.30	1.50	1.43
6	A	609	GOL	O1-C1	3.90	1.59	1.42
2	A	601	FAD	C4-C4X	4.28	1.50	1.41
2	A	601	FAD	C2A-N3A	4.50	1.40	1.32
5	A	604	TRS	C-N	4.99	1.57	1.50
2	A	601	FAD	C5X-N5	5.11	1.43	1.35
2	A	601	FAD	C10-N10	5.71	1.45	1.39

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	601	FAD	C4-C4X-C10	-11.79	112.39	119.94
5	A	608	TRS	C1-C-N	-10.83	89.49	107.88
5	A	608	TRS	C2-C-N	-10.75	89.62	107.88
5	A	608	TRS	O1-C1-C	-8.25	91.33	110.92
4	A	603	NCA	C3-C7-N7	-7.31	109.56	117.82
2	A	601	FAD	C4X-C10-N10	-6.11	116.08	120.52
2	A	601	FAD	N3-C2-N1	-5.18	118.97	127.69
2	A	601	FAD	C4B-O4B-C1B	-4.79	104.56	109.64
2	A	601	FAD	C4X-C4-N3	-4.74	117.32	123.52
5	A	604	TRS	C2-C-N	-4.73	99.85	107.88
3	A	602	NAP	C3N-C7N-N7N	-4.35	112.91	117.82
3	A	602	NAP	N3A-C2A-N1A	-4.13	125.62	128.87
4	A	603	NCA	C4-C3-C7	-3.33	112.26	121.11
2	A	601	FAD	N3A-C2A-N1A	-2.54	126.88	128.87
5	A	604	TRS	O1-C1-C	-2.51	104.97	110.92
2	A	601	FAD	O4B-C1B-N9A	-2.39	103.59	108.11
5	A	604	TRS	C1-C-N	-2.35	103.89	107.88
2	A	601	FAD	C4X-N5-C5X	-2.03	114.34	116.72
4	A	603	NCA	C4-C3-C2	2.01	119.95	117.66
6	A	609	GOL	C3-C2-C1	2.23	120.44	111.06
5	A	608	TRS	O2-C2-C	2.31	116.40	110.92
5	A	604	TRS	C2-C-C1	2.38	115.65	110.65
2	A	601	FAD	C2A-N1A-C6A	2.38	123.01	118.77
2	A	601	FAD	C1'-N10-C9A	2.50	121.73	118.83
4	A	603	NCA	C2-C3-C7	2.53	127.64	121.34
5	A	604	TRS	O2-C2-C	2.71	117.35	110.92
5	A	604	TRS	C3-C-C1	3.07	117.12	110.65
6	A	610	GOL	C3-C2-C1	3.62	126.32	111.06
4	A	603	NCA	O7-C7-C3	4.65	124.75	119.60
2	A	601	FAD	C5X-C9A-N10	5.04	121.36	117.58
2	A	601	FAD	N6A-C6A-N1A	5.33	127.46	118.52
5	A	608	TRS	C2-C-C1	11.61	135.11	110.65
2	A	601	FAD	C4-N3-C2	13.92	126.78	115.16

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	601	FAD	1	0
3	A	602	NAP	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	603	NCA	2	0
6	A	610	GOL	3	0

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	529/541 (97%)	-0.31	1 (0%) 95 95	9, 13, 27, 43	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	6	THR	3.3

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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
6	GOL	A	609	6/6	0.92	0.17	12.07	30,39,53,65	0
5	TRS	A	608	8/8	0.87	0.18	6.09	29,36,41,55	0
6	GOL	A	610	6/6	0.95	0.22	4.82	27,33,38,41	0
4	NCA	A	603	9/9	0.96	0.08	1.03	17,20,23,25	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	FAD	A	601	53/53	0.99	0.07	-0.29	8,9,11,15	0
6	GOL	A	607	6/6	0.98	0.05	-0.74	13,17,20,22	0
3	NAP	A	602	48/48	0.99	0.05	-0.77	9,11,15,18	0
6	GOL	A	605	6/6	0.99	0.05	-1.22	12,14,18,21	0
6	GOL	A	606	6/6	0.99	0.05	-1.61	14,16,16,19	0
5	TRS	A	604	8/8	0.96	0.09	-	17,20,25,29	0

6.5 Other polymers [i](#)

There are no such residues in this entry.