



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 09:55 AM GMT

PDB ID : 1X7D
Title : Crystal Structure Analysis of Ornithine Cyclodeaminase Complexed with NAD and ornithine to 1.6 Angstroms
Authors : Alam, S.; Goodman, J.L.; Wang, S.; Ruzicka, F.J.; Frey, P.A.; Wedekind, J.E.
Deposited on : 2004-08-13
Resolution : 1.60 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

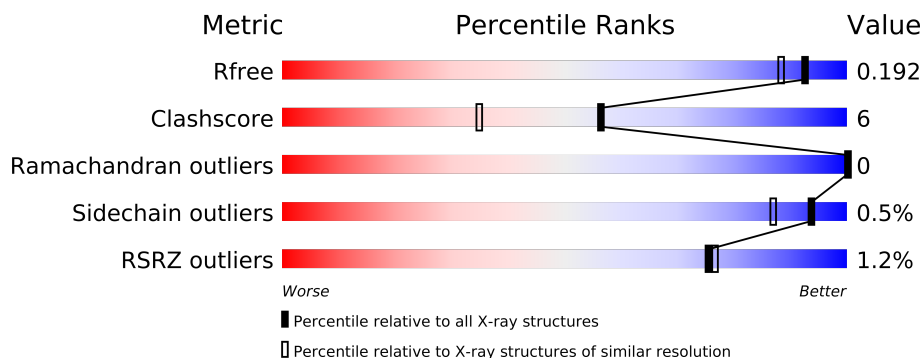
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 1.60 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}	66092	1872 (1.60-1.60)
Clashscore	79885	2199 (1.60-1.60)
Ramachandran outliers	78287	2126 (1.60-1.60)
Sidechain outliers	78261	2125 (1.60-1.60)
RSRZ outliers	66119	1872 (1.60-1.60)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	350	
1	B	350	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
4	ORN	A	1101[A]	-	X
4	ORN	A	1101[B]	-	X
4	ORN	A	1102[A]	-	X
4	ORN	A	1102[B]	-	X
5	MES	B	3158	-	X
6	MPD	A	1401	-	X

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Mol	Type	Chain	Res	Geometry	Electron density
6	MPD	A	1402	-	X
6	MPD	A	1403	-	X
6	MPD	A	1404	-	X

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 6249 atoms, of which 0 are hydrogen and 0 are deuterium.

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 ornithine cyclodeaminase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	340	Total	C	N	O	S	Se	0	0	0
			2620	1659	454	498	1	8			
1	B	339	Total	C	N	O	S	Se	0	0	0
			2615	1656	453	497	1	8			

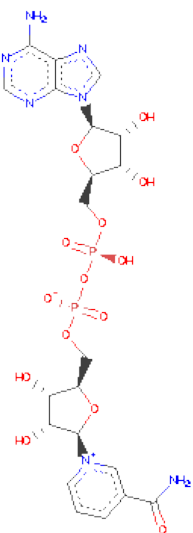
There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	10	MSE	MET	MODIFIED RESIDUE	UNP Q88H32
A	58	MSE	MET	MODIFIED RESIDUE	UNP Q88H32
A	86	MSE	MET	MODIFIED RESIDUE	UNP Q88H32
A	119	MSE	MET	MODIFIED RESIDUE	UNP Q88H32
A	132	MSE	MET	MODIFIED RESIDUE	UNP Q88H32
A	215	MSE	MET	MODIFIED RESIDUE	UNP Q88H32
A	220	MSE	MET	MODIFIED RESIDUE	UNP Q88H32
A	316	MSE	MET	MODIFIED RESIDUE	UNP Q88H32
B	10	MSE	MET	MODIFIED RESIDUE	UNP Q88H32
B	58	MSE	MET	MODIFIED RESIDUE	UNP Q88H32
B	86	MSE	MET	MODIFIED RESIDUE	UNP Q88H32
B	119	MSE	MET	MODIFIED RESIDUE	UNP Q88H32
B	132	MSE	MET	MODIFIED RESIDUE	UNP Q88H32
B	215	MSE	MET	MODIFIED RESIDUE	UNP Q88H32
B	220	MSE	MET	MODIFIED RESIDUE	UNP Q88H32
B	316	MSE	MET	MODIFIED RESIDUE	UNP Q88H32

- Molecule 2 is SODIUM ION (three-letter code: NA) (formula: Na).

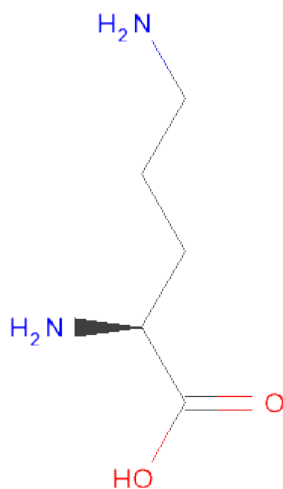
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Na	0	0
			1	1		
2	A	1	Total	Na	0	0
			1	1		

- Molecule 3 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: $C_{21}H_{27}N_7O_{14}P_2$).



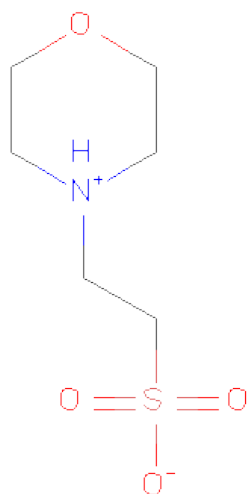
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	B	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
3	A	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

- Molecule 4 is L-ORNITHINE (three-letter code: ORN) (formula: $C_5H_{12}N_2O_2$).



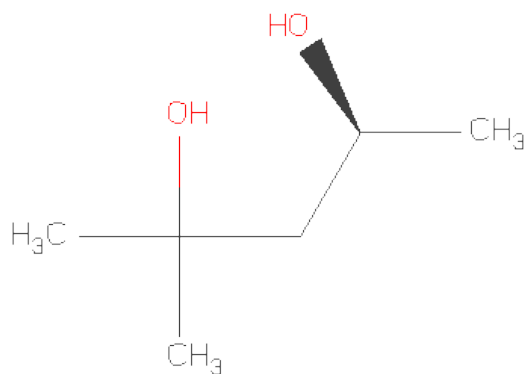
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	2	Total	C	N	O	0	2
			36	20	8	8		

- Molecule 5 is 2-(N-MORPHOLINO)-ETHANESULFONICACID (three-letter code: MES) (formula: $C_6H_{13}NO_4S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	B	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

- Molecule 6 is (4S)-2-METHYL-2,4-PENTANEDIOL (three-letter code: MPD) (formula: $C_6H_{14}O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			8	6	2		
6	A	1	Total	C	O	0	0
			8	6	2		
6	A	1	Total	C	O	0	0
			8	6	2		
6	A	1	Total	C	O	0	0
			8	6	2		
6	B	1	Total	C	O	0	0
			8	6	2		
6	B	1	Total	C	O	0	0
			8	6	2		

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	443	Total	O	0	0
			443	443		
7	B	385	Total	O	0	0
			385	385		

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.

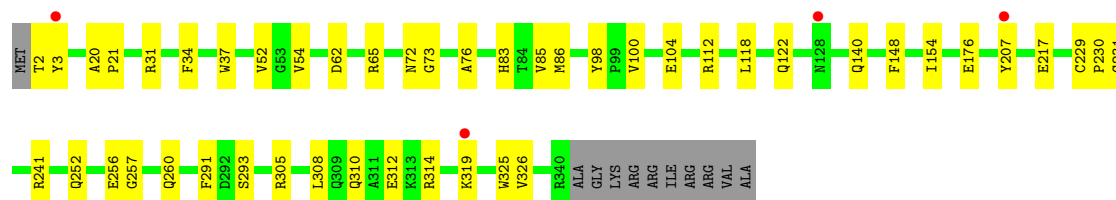
- Molecule 1: ornithine cyclodeaminase

Chain A: 



- Molecule 1: ornithine cyclodeaminase

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	69.88Å 78.62Å 119.94Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	24.95 – 1.60 24.95 – 1.60	Depositor EDS
% Data completeness (in resolution range)	98.6 (24.95-1.60) 98.8 (24.95-1.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.97 (at 1.60Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.170 , 0.190 0.170 , 0.192	Depositor DCC
R_{free} test set	4400 reflections (5.08%)	DCC
Wilson B-factor (Å ²)	13.2	Xtriage
Anisotropy	0.131	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 43.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 86671 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6249	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 4.44% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ORN, NA, MPD, MES, NAD

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	0.29	0/2664	0.63	1/3609 (0.0%)
1	B	0.29	0/2659	0.62	0/3602
All	All	0.29	0/5323	0.63	1/7211 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	281	GLY	N-CA-C	-5.04	100.49	113.10

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2620	0	2608	31	0
1	B	2615	0	2603	39	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	44	0	26	1	0
3	B	44	0	26	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	36	0	44	12	0
5	B	12	0	13	2	0
6	A	32	0	56	0	0
6	B	16	0	28	1	0
7	A	443	0	0	7	0
7	B	385	0	0	7	0
All	All	6249	0	5404	70	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 6.

All (70) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:52:VAL:HG22	7:A:1754:HOH:O	1.83	0.78
1:A:53:GLY:HA2	1:A:205:LYS:HE3	1.64	0.78
1:A:51:GLU:HG2	1:A:52:VAL:HG13	1.67	0.76
1:B:20:ALA:HB3	1:B:21:PRO:HD3	1.69	0.74
4:A:1102[B]:ORN:NE	1:B:73:GLY:HA3	2.03	0.73
1:A:183:ARG:HD2	7:A:1825:HOH:O	1.91	0.70
1:B:310:GLN:NE2	1:B:314:ARG:HH12	1.90	0.69
4:A:1102[A]:ORN:HE2	1:B:85:VAL:HG22	1.58	0.69
1:A:73:GLY:HA3	4:A:1101[A]:ORN:NE	2.06	0.68
1:A:308:LEU:O	1:A:312:GLU:HG3	1.93	0.68
1:B:2:THR:N	1:B:98:TYR:HH	1.93	0.67
1:B:3:TYR:CE1	1:B:319:LYS:HG3	2.30	0.67
1:B:308:LEU:O	1:B:312:GLU:HG3	1.98	0.63
1:A:54:VAL:CG1	4:A:1101[A]:ORN:HE2	2.12	0.62
1:A:20:ALA:HB3	1:A:21:PRO:HD3	1.82	0.62
1:A:31:ARG:HD2	7:A:1900:HOH:O	1.99	0.62
1:A:2:THR:N	1:A:98:TYR:HH	1.99	0.60
1:B:72:ASN:ND2	1:B:86:MSE:HB2	2.16	0.60
1:A:73:GLY:HA3	4:A:1101[B]:ORN:HE1	1.67	0.60
4:A:1102[B]:ORN:HE2	1:B:54:VAL:CG1	2.16	0.58
1:B:72:ASN:HD21	1:B:86:MSE:HB2	1.70	0.57
1:A:325:TRP:CE3	1:A:326:VAL:HG22	2.40	0.56
1:A:73:GLY:HA3	4:A:1101[B]:ORN:NE	2.21	0.55
1:B:100:VAL:HG21	6:B:1406:MPD:HM3	1.88	0.55
1:A:53:GLY:CA	1:A:205:LYS:HE3	2.35	0.55
1:B:229:CYS:HA	1:B:256:GLU:HB3	1.88	0.55
1:B:217:GLU:HA	7:B:3440:HOH:O	2.09	0.52
1:B:207:TYR:CZ	1:B:231:GLY:HA2	2.45	0.51
1:B:229:CYS:HB2	1:B:230:PRO:HD2	1.92	0.51

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:176:GLU:OE2	5:B:3158:MES:H32	2.11	0.51
1:B:241:ARG:HD2	7:B:3529:HOH:O	2.10	0.50
1:A:229:CYS:HA	1:A:256:GLU:HB3	1.93	0.50
1:A:73:GLY:HA3	4:A:1101[A]:ORN:HE1	1.74	0.50
1:A:83:HIS:HD2	7:B:3234:HOH:O	1.95	0.50
4:A:1102[A]:ORN:HE1	1:B:73:GLY:HA3	1.77	0.49
4:A:1102[A]:ORN:NE	1:B:85:VAL:HG22	2.26	0.49
1:B:118:LEU:O	1:B:122:GLN:HG3	2.13	0.48
1:B:65:ARG:HD2	1:B:305:ARG:HD3	1.95	0.48
1:B:325:TRP:CG	1:B:326:VAL:N	2.79	0.47
1:A:207:TYR:CZ	1:A:231:GLY:HA2	2.50	0.47
1:B:148:PHE:HB3	1:B:154:ILE:HG13	1.97	0.47
1:B:252:GLN:HG2	1:B:256:GLU:OE1	2.15	0.47
1:B:310:GLN:HE21	1:B:314:ARG:HH12	1.62	0.46
7:A:1583:HOH:O	1:B:83:HIS:HD2	1.98	0.46
1:B:34:PHE:O	1:B:37:TRP:HB2	2.15	0.46
1:A:78:THR:HG23	7:A:1760:HOH:O	2.16	0.45
1:B:176:GLU:OE2	5:B:3158:MES:H72	2.16	0.45
1:B:140:GLN:HG2	3:B:1301:NAD:PN	2.56	0.45
1:B:52:VAL:HG22	7:B:3310:HOH:O	2.17	0.45
1:A:54:VAL:HG13	4:A:1101[A]:ORN:HE2	1.81	0.45
1:A:73:GLY:CA	4:A:1101[B]:ORN:NE	2.79	0.45
1:B:257:GLY:O	1:B:260:GLN:HG2	2.17	0.45
1:B:65:ARG:NE	1:B:305:ARG:HH11	2.15	0.44
1:A:51:GLU:HG3	7:A:1932:HOH:O	2.17	0.44
1:A:72:ASN:OD1	1:A:86:MSE:HB2	2.17	0.44
1:B:31:ARG:NH2	1:B:122:GLN:OE1	2.52	0.43
1:B:310:GLN:HE21	1:B:314:ARG:NH1	2.17	0.43
1:B:217:GLU:CA	7:B:3440:HOH:O	2.66	0.43
1:B:76:ALA:HB3	7:B:3368:HOH:O	2.18	0.43
1:B:291:PHE:CE2	1:B:293:SER:HB2	2.54	0.43
1:B:62:ASP:HB2	7:B:3511:HOH:O	2.18	0.42
1:A:240:LEU:CB	1:A:262:LEU:HD11	2.49	0.42
1:A:31:ARG:HD3	7:A:1847:HOH:O	2.20	0.41
1:A:118:LEU:O	1:A:122:GLN:HG3	2.19	0.41
1:A:140:GLN:HG2	3:A:1302:NAD:PN	2.60	0.41
1:B:310:GLN:NE2	1:B:314:ARG:NH1	2.66	0.41
1:A:31:ARG:O	1:A:35:LYS:HG3	2.21	0.41
1:A:291:PHE:CE2	1:A:293:SER:HB2	2.56	0.40
1:A:58:MSE:O	1:A:68:PHE:HA	2.21	0.40
1:A:240:LEU:HB2	1:A:262:LEU:HD11	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	338/350 (97%)	326 (96%)	12 (4%)	0	100	100
1	B	337/350 (96%)	323 (96%)	14 (4%)	0	100	100
All	All	675/700 (96%)	649 (96%)	26 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	275/275 (100%)	274 (100%)	1 (0%)	95	89
1	B	275/275 (100%)	273 (99%)	2 (1%)	91	81
All	All	550/550 (100%)	547 (100%)	3 (0%)	94	87

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

Mol	Chain	Res	Type
1	A	112	ARG
1	B	104	GLU
1	B	112	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	83	HIS
1	A	144	GLN
1	A	223	ASN

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Mol	Chain	Res	Type
1	B	72	ASN
1	B	83	HIS
1	B	137	ASN
1	B	144	GLN
1	B	223	ASN
1	B	309	GLN
1	B	310	GLN

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 15 ligands modelled in this entry, 2 are monoatomic - leaving 13 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	ORN	A	1101[A]	-	8,8,8	0.71	0	9,9,9	0.60	0
4	ORN	A	1101[B]	-	8,8,8	0.74	0	9,9,9	0.59	0
4	ORN	A	1102[A]	-	8,8,8	0.68	0	9,9,9	0.59	0
4	ORN	A	1102[B]	-	8,8,8	0.69	0	9,9,9	0.61	0
3	NAD	A	1302	-	48,48,48	2.35	11 (22%)	73,73,73	2.16	10 (13%)
6	MPD	A	1401	-	7,7,7	0.73	0	10,10,10	0.76	0
6	MPD	A	1402	-	7,7,7	0.78	0	10,10,10	0.71	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	MPD	A	1403	-	7,7,7	0.97	0	10,10,10	0.63	0
6	MPD	A	1404	-	7,7,7	0.69	0	10,10,10	0.60	0
3	NAD	B	1301	-	48,48,48	2.35	11 (22%)	73,73,73	2.17	10 (13%)
6	MPD	B	1405	-	7,7,7	0.70	0	10,10,10	0.61	0
6	MPD	B	1406	-	7,7,7	0.70	0	10,10,10	0.66	0
5	MES	B	3158	-	12,12,12	2.49	6 (50%)	16,16,16	2.90	8 (50%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	ORN	A	1101[A]	-	-	0/8/8/8	0/0/0/0
4	ORN	A	1101[B]	-	-	0/8/8/8	0/0/0/0
4	ORN	A	1102[A]	-	-	0/8/8/8	0/0/0/0
4	ORN	A	1102[B]	-	-	0/8/8/8	0/0/0/0
3	NAD	A	1302	-	-	0/30/62/62	0/3/5/5
6	MPD	A	1401	-	-	0/5/5/5	0/0/0/0
6	MPD	A	1402	-	-	0/5/5/5	0/0/0/0
6	MPD	A	1403	-	-	0/5/5/5	0/0/0/0
6	MPD	A	1404	-	-	0/5/5/5	0/0/0/0
3	NAD	B	1301	-	-	0/30/62/62	0/3/5/5
6	MPD	B	1405	-	-	0/5/5/5	0/0/0/0
6	MPD	B	1406	-	-	0/5/5/5	0/0/0/0
5	MES	B	3158	-	-	0/6/14/14	0/1/1/1

All (28) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1301	NAD	C2N-N1N	8.03	1.45	1.35
3	A	1302	NAD	C2N-N1N	7.96	1.45	1.35
3	A	1302	NAD	C4N-C3N	7.92	1.53	1.39
3	B	1301	NAD	C4N-C3N	7.84	1.52	1.39
3	B	1301	NAD	C5N-C4N	5.33	1.51	1.39
5	B	3158	MES	O2S-S	5.23	1.57	1.45
3	A	1302	NAD	C5N-C4N	5.22	1.51	1.39
3	B	1301	NAD	C2A-N3A	4.38	1.40	1.32
3	A	1302	NAD	C2A-N3A	4.26	1.40	1.32
3	B	1301	NAD	C6N-N1N	4.08	1.47	1.35
3	A	1302	NAD	C6N-N1N	4.01	1.46	1.35
5	B	3158	MES	O3S-S	-3.53	1.37	1.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	3158	MES	C8-S	3.33	1.84	1.78
3	A	1302	NAD	C4A-N3A	3.29	1.40	1.35
3	B	1301	NAD	C4A-N3A	3.24	1.40	1.35
3	A	1302	NAD	C2N-C3N	-3.09	1.34	1.38
3	B	1301	NAD	C2N-C3N	-2.92	1.34	1.38
3	A	1302	NAD	C2D-C1D	2.71	1.57	1.53
3	A	1302	NAD	O4D-C1D	2.52	1.45	1.41
5	B	3158	MES	C3-C2	2.48	1.56	1.51
3	B	1301	NAD	C2D-C1D	2.46	1.56	1.53
3	B	1301	NAD	O4D-C1D	2.31	1.44	1.41
5	B	3158	MES	C3-N4	2.28	1.55	1.49
3	A	1302	NAD	C8A-N9A	2.23	1.39	1.36
3	B	1301	NAD	O4D-C4D	2.22	1.50	1.45
3	A	1302	NAD	O4D-C4D	2.18	1.50	1.45
5	B	3158	MES	O1S-S	2.07	1.50	1.45
3	B	1301	NAD	C8A-N9A	2.01	1.39	1.36

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1301	NAD	C5N-C4N-C3N	-9.20	108.37	120.32
3	A	1302	NAD	C5N-C4N-C3N	-8.99	108.64	120.32
3	B	1301	NAD	N3A-C2A-N1A	-8.27	121.79	128.71
3	A	1302	NAD	N3A-C2A-N1A	-7.97	122.04	128.71
5	B	3158	MES	O3S-S-O1S	7.42	131.57	112.48
3	B	1301	NAD	C6N-N1N-C2N	-6.53	114.67	122.04
3	A	1302	NAD	C6N-N1N-C2N	-6.43	114.77	122.04
3	A	1302	NAD	C8A-N9A-C4A	-5.15	102.97	106.90
3	B	1301	NAD	C8A-N9A-C4A	-4.92	103.14	106.90
3	A	1302	NAD	O4D-C1D-N1N	4.85	112.92	107.95
3	B	1301	NAD	O4D-C1D-N1N	4.76	112.82	107.95
5	B	3158	MES	O3S-S-C8	-4.14	95.31	105.99
3	B	1301	NAD	C2A-N1A-C6A	3.90	125.81	118.77
3	A	1302	NAD	C2A-N1A-C6A	3.83	125.69	118.77
5	B	3158	MES	C6-C5-N4	3.66	114.66	109.96
3	B	1301	NAD	C8A-N9A-C1B	3.56	133.39	126.38
3	A	1302	NAD	C8A-N9A-C1B	3.52	133.32	126.38
5	B	3158	MES	C2-C3-N4	-3.17	105.90	109.96
5	B	3158	MES	O1S-S-C8	-3.13	96.70	106.36
3	A	1302	NAD	O4D-C1D-C2D	-2.76	102.54	106.77
5	B	3158	MES	O2S-S-O1S	2.68	117.61	112.44
5	B	3158	MES	O3S-S-O2S	-2.65	105.68	112.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1302	NAD	C4B-O4B-C1B	-2.62	106.90	109.75
3	B	1301	NAD	C6N-C5N-C4N	-2.60	115.31	119.44
3	B	1301	NAD	O4D-C1D-C2D	-2.46	103.00	106.77
3	A	1302	NAD	C6N-C5N-C4N	-2.43	115.58	119.44
3	B	1301	NAD	C4B-O4B-C1B	-2.23	107.32	109.75
5	B	3158	MES	C7-N4-C3	-2.07	106.24	111.66

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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

6.1 Protein, DNA and RNA chains

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	340/350 (97%)	-0.32	4 (1%) 75 76	7, 12, 23, 38	0
1	B	339/350 (96%)	-0.25	4 (1%) 75 76	6, 13, 23, 33	0
All	All	679/700 (97%)	-0.29	8 (1%) 75 76	6, 12, 23, 38	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	51	GLU	4.2
1	A	341	ALA	3.9
1	A	207	TYR	3.3
1	A	325	TRP	2.5
1	B	207	TYR	2.4
1	B	3	TYR	2.3
1	B	319	LYS	2.3
1	B	128	ASN	2.0

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

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

6.3 Carbohydrates

There are no carbohydrates in this entry.

6.4 Ligands

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
6	MPD	A	1402	8/8	0.31	19.73	45,46,46,47	0
5	MES	B	3158	12/12	0.29	11.95	36,39,49,49	0
6	MPD	A	1401	8/8	0.18	11.08	44,45,45,46	0
4	ORN	A	1102[A]	9/9	0.19	10.06	10,13,19,28	9
4	ORN	A	1102[B]	9/9	0.19	9.55	10,13,19,28	9
6	MPD	A	1403	8/8	0.33	7.57	51,51,52,52	0
4	ORN	A	1101[A]	9/9	0.15	6.82	8,9,15,24	9
4	ORN	A	1101[B]	9/9	0.15	6.39	8,9,15,24	9
6	MPD	A	1404	8/8	0.18	2.12	27,29,30,30	0
6	MPD	B	1406	8/8	0.17	1.56	41,42,43,43	0
6	MPD	B	1405	8/8	0.15	0.28	29,32,33,34	0
2	NA	A	1501	1/1	0.08	0.23	12,12,12,12	0
3	NAD	B	1301	44/44	0.08	0.21	7,10,14,16	0
2	NA	B	1502	1/1	0.08	-0.07	13,13,13,13	0
3	NAD	A	1302	44/44	0.07	-0.26	7,9,13,14	0

6.5 Other polymers ⓘ

There are no such residues in this entry.