



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 12, 2024 – 02:14 AM EST

PDB ID : 1X87
Title : 2.4A X-ray structure of Urocanase protein complexed with NAD
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Deposited on : 2004-08-17
Resolution : 2.40 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	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	3.0
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.39

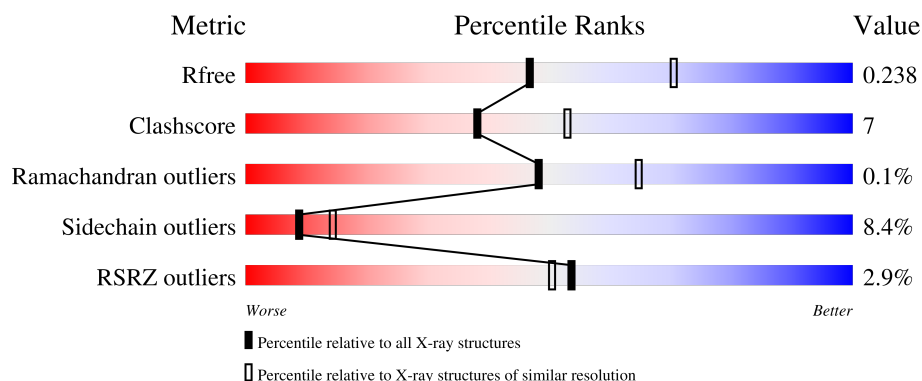
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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}	164625	4642 (2.40-2.40)
Clashscore	180529	5218 (2.40-2.40)
Ramachandran outliers	177936	5158 (2.40-2.40)
Sidechain outliers	177891	5159 (2.40-2.40)
RSRZ outliers	164620	4642 (2.40-2.40)

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	551	<div> <div> <div>0%</div> <div>71%</div> <div>17%</div> <div>• 10%</div> </div> </div>
1	B	551	<div> <div>4%</div> <div>72%</div> <div>14%</div> <div>• • 11%</div> </div>

2 Entry composition [i](#)

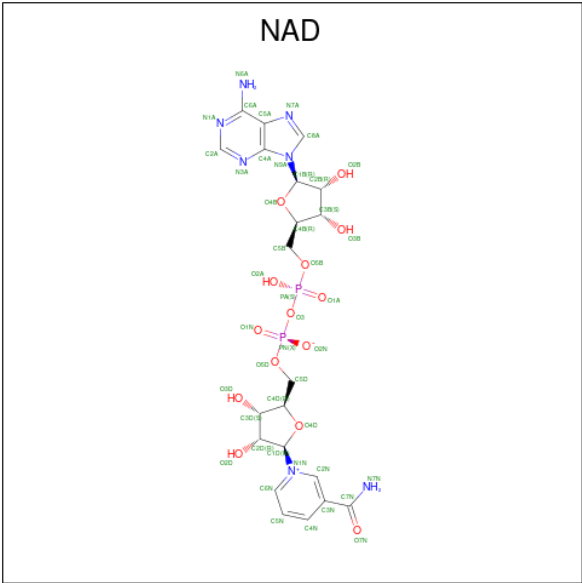
There are 3 unique types of molecules in this entry. The entry contains 7883 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Urocanase protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	A	495	3811	2412	681	699	5	14	0	2	0
1	B	493	3766	2371	671	705	5	14	0	0	0

- Molecule 2 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: C₂₁H₂₇N₇O₁₄P₂).



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

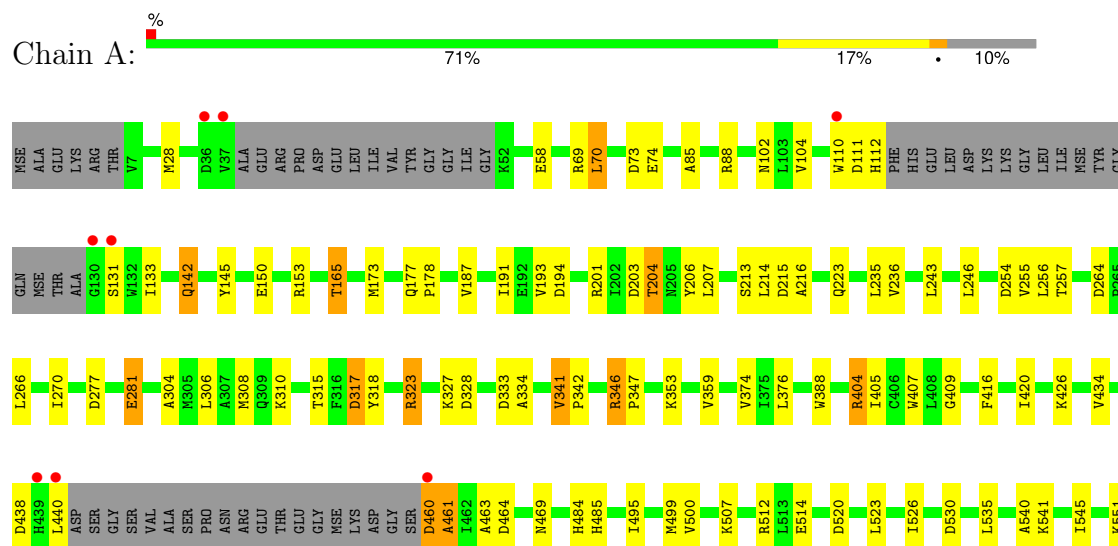
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	156	Total 156	O 156	0	0
3	B	62	Total 62	O 62	0	0

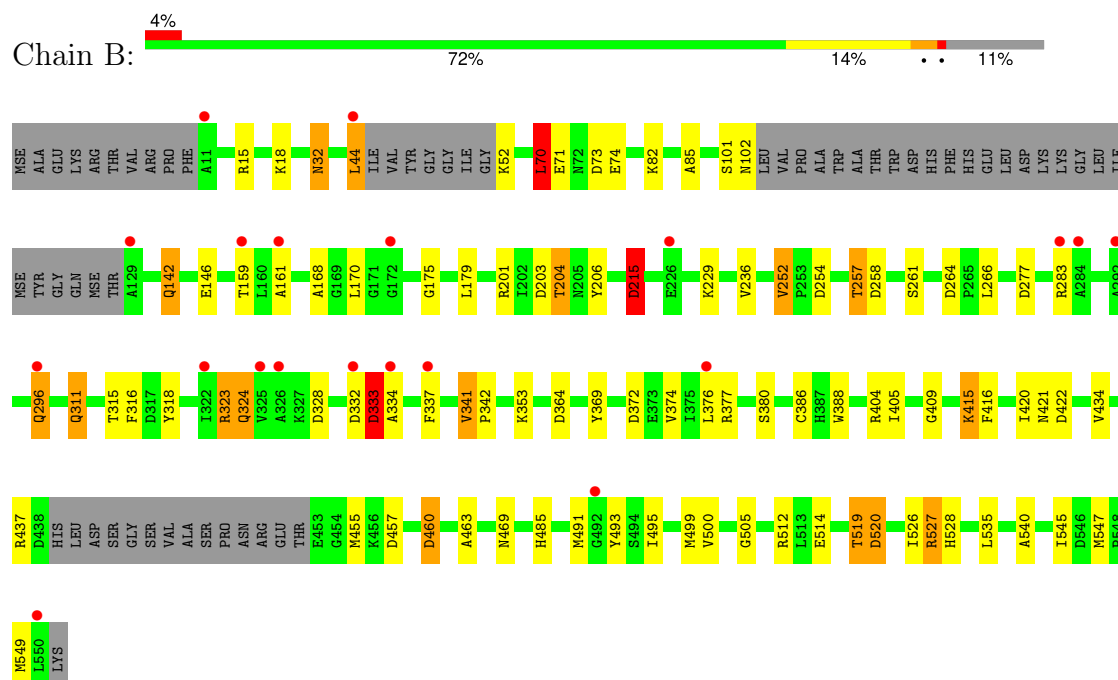
3 Residue-property plots

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 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: Urocanase protein



• Molecule 1: Urocanase protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	100.12Å 101.61Å 113.85Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.40 20.00 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.8 (20.00-2.40) 99.6 (20.00-2.40)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.12 (at 2.41Å)	Xtriage
Refinement program	REFMAC 5.1.9999	Depositor
R, R_{free}	0.187 , 0.232 0.192 , 0.238	Depositor DCC
R_{free} test set	2323 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	46.1	Xtriage
Anisotropy	0.050	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 48.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.017 for k,h,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7883	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.68% 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: 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	0.64	0/3887	1.04	18/5245 (0.3%)
1	B	0.54	0/3821	1.00	16/5147 (0.3%)
All	All	0.59	0/7708	1.02	34/10392 (0.3%)

There are no bond length outliers.

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	512	ARG	NE-CZ-NH2	-9.00	115.80	120.30
1	A	111	ASP	CB-CG-OD2	7.38	124.94	118.30
1	A	254	ASP	CB-CG-OD2	7.32	124.89	118.30
1	B	364	ASP	CB-CG-OD2	7.23	124.81	118.30
1	A	264	ASP	CB-CG-OD2	7.03	124.62	118.30
1	A	520	ASP	CB-CG-OD2	6.91	124.51	118.30
1	B	203	ASP	CB-CG-OD2	6.80	124.42	118.30
1	B	460	ASP	CB-CG-OD2	6.70	124.33	118.30
1	B	264	ASP	CB-CG-OD2	6.67	124.31	118.30
1	A	404	ARG	NE-CZ-NH2	-6.30	117.15	120.30
1	A	317	ASP	CB-CG-OD2	6.25	123.93	118.30
1	A	333	ASP	CB-CG-OD2	6.16	123.85	118.30
1	A	201	ARG	NE-CZ-NH2	-6.09	117.25	120.30
1	A	277	ASP	CB-CG-OD2	6.05	123.75	118.30
1	A	512	ARG	NE-CZ-NH1	6.05	123.33	120.30
1	B	520	ASP	CB-CG-OD2	6.03	123.73	118.30
1	A	323	ARG	NE-CZ-NH1	6.02	123.31	120.30
1	B	328	ASP	CB-CG-OD2	5.93	123.64	118.30
1	B	422	ASP	CB-CG-OD2	5.71	123.44	118.30
1	A	328	ASP	CB-CG-OD2	5.70	123.43	118.30
1	B	73	ASP	CB-CG-OD2	5.63	123.37	118.30
1	A	194	ASP	CB-CG-OD2	5.62	123.35	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	203	ASP	CB-CG-OD2	5.59	123.33	118.30
1	B	254	ASP	CB-CG-OD2	5.56	123.31	118.30
1	B	457	ASP	CB-CG-OD2	5.50	123.25	118.30
1	B	512	ARG	NE-CZ-NH2	-5.50	117.55	120.30
1	B	333	ASP	CB-CG-OD2	5.37	123.13	118.30
1	A	438	ASP	CB-CG-OD2	5.27	123.04	118.30
1	A	73	ASP	CB-CG-OD2	5.25	123.02	118.30
1	B	332	ASP	CB-CG-OD2	5.23	123.01	118.30
1	B	215	ASP	CB-CG-OD2	5.12	122.91	118.30
1	B	70	LEU	CA-CB-CG	5.11	127.05	115.30
1	B	277	ASP	CB-CG-OD2	5.09	122.88	118.30
1	A	215	ASP	CB-CG-OD2	5.04	122.83	118.30

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	3811	0	3806	56	0
1	B	3766	0	3758	58	0
2	A	44	0	26	1	0
2	B	44	0	26	0	0
3	A	156	0	0	0	0
3	B	62	0	0	3	0
All	All	7883	0	7616	104	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 7.

All (104) 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:324:GLN:HE21	1:B:324:GLN:HA	1.40	0.85
1:A:243:LEU:HG	1:A:308[B]:MSE:HE1	1.58	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:204:THR:HG23	1:A:206:TYR:HD1	1.52	0.74
1:A:150:GLU:OE1	1:A:153:ARG:NH2	2.22	0.73
1:B:463:ALA:HB1	1:B:495:ILE:HD12	1.70	0.72
1:B:415:LYS:HB3	1:B:415:LYS:NZ	2.08	0.69
1:A:526:ILE:HD13	1:B:85:ALA:HB2	1.77	0.68
1:A:461:ALA:HA	1:B:101:SER:OG	1.94	0.66
1:B:170:LEU:O	1:B:201:ARG:NH1	2.30	0.64
1:B:142:GLN:HG2	1:B:405:ILE:HG23	1.81	0.63
1:B:168:ALA:HB3	1:B:258:ASP:OD2	1.99	0.63
1:B:415:LYS:HB3	1:B:415:LYS:HZ2	1.65	0.62
1:A:317:ASP:HB2	1:A:323:ARG:HD2	1.81	0.61
1:A:374:VAL:HG22	1:A:416:PHE:HE1	1.66	0.61
1:A:526:ILE:HD13	1:B:85:ALA:CB	2.30	0.60
1:B:527:ARG:NE	3:B:856:HOH:O	2.31	0.60
1:A:142:GLN:HG2	1:A:405:ILE:HG23	1.84	0.60
1:A:464:ASP:OD2	1:B:485:HIS:NE2	2.33	0.59
1:A:28:MSE:HE2	1:A:133:ILE:HG22	1.84	0.59
1:B:416:PHE:CZ	1:B:420:ILE:HD11	2.37	0.59
1:B:266:LEU:O	1:B:283:ARG:HD3	2.02	0.59
1:A:165:THR:HB	1:A:255:VAL:HB	1.85	0.58
1:A:85:ALA:HB2	1:B:526:ILE:HD13	1.85	0.58
1:B:32:ASN:ND2	3:B:805:HOH:O	2.37	0.58
1:A:191:ILE:N	1:A:191:ILE:HD13	2.20	0.57
1:A:541:LYS:NZ	1:A:551:LYS:HG3	2.19	0.57
1:A:485[B]:HIS:CE1	1:B:495:ILE:HD11	2.40	0.56
1:A:463:ALA:HB1	1:A:495:ILE:HD12	1.88	0.56
1:A:256:LEU:HD21	1:A:308[B]:MSE:HE3	1.88	0.56
1:A:214:LEU:HD11	1:A:246:LEU:HD21	1.90	0.53
1:B:204:THR:HG23	1:B:206:TYR:HD1	1.73	0.53
1:B:257:THR:HG21	1:B:316:PHE:CE2	2.44	0.53
1:A:204:THR:HG23	1:A:206:TYR:CD1	2.38	0.53
1:A:88:ARG:NH1	1:B:514:GLU:OE2	2.42	0.52
1:B:323:ARG:HD2	1:B:337:PHE:CE1	2.44	0.52
1:A:173:MSE:HE3	2:A:600:NAD:O2N	2.10	0.52
1:B:44:LEU:HA	1:B:52:LYS:HG2	1.91	0.52
1:B:353:LYS:HA	1:B:409:GLY:HA2	1.91	0.52
1:B:159:THR:HG22	1:B:161:ALA:H	1.74	0.51
1:A:243:LEU:CG	1:A:308[B]:MSE:HE1	2.36	0.51
1:A:341:VAL:HG22	1:A:342:PRO:HD3	1.92	0.50
1:A:346:ARG:HB3	1:A:347:PRO:HD3	1.93	0.50
1:B:204:THR:HG23	1:B:206:TYR:CD1	2.47	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:460:ASP:O	1:A:461:ALA:HB2	2.12	0.50
1:B:324:GLN:HE21	1:B:324:GLN:CA	2.18	0.49
1:A:28:MSE:HE1	1:A:359:VAL:HG21	1.94	0.49
1:B:380:SER:HA	1:B:386:CYS:SG	2.52	0.49
1:A:323:ARG:HG2	1:A:334:ALA:O	2.12	0.49
1:B:215:ASP:OD2	1:B:215:ASP:N	2.45	0.49
1:A:485[B]:HIS:CE1	1:A:495:ILE:HG12	2.48	0.49
1:B:142:GLN:HE22	1:B:146:GLU:HG3	1.77	0.49
1:A:541:LYS:HZ3	1:A:551:LYS:HG3	1.78	0.48
1:B:70:LEU:HD22	1:B:74:GLU:HB2	1.95	0.48
1:A:353:LYS:HA	1:A:409:GLY:HA2	1.95	0.47
1:B:257:THR:HG22	1:B:316:PHE:O	2.14	0.47
1:A:374:VAL:HG22	1:A:416:PHE:CE1	2.49	0.47
1:B:526:ILE:HG23	1:B:547:MSE:HE3	1.97	0.47
1:B:519:THR:HG21	3:B:821:HOH:O	2.14	0.47
1:A:145:TYR:CE1	1:A:388:TRP:HB2	2.49	0.47
1:A:416:PHE:CE2	1:A:420:ILE:HD11	2.50	0.47
1:B:324:GLN:HA	1:B:324:GLN:NE2	2.19	0.47
1:B:421:ASN:ND2	1:B:505:GLY:H	2.13	0.47
1:A:177:GLN:HB2	1:A:178:PRO:HD3	1.96	0.46
1:A:193:VAL:HG21	1:A:270:ILE:HG13	1.98	0.46
1:B:296:GLN:HE21	1:B:296:GLN:HA	1.81	0.46
1:B:526:ILE:HD11	1:B:545:ILE:CG2	2.46	0.46
1:A:70:LEU:HD22	1:A:74:GLU:HB2	1.97	0.46
1:A:110:TRP:O	1:A:112:HIS:N	2.49	0.46
1:A:317:ASP:HB2	1:A:323:ARG:CD	2.45	0.46
1:B:493:TYR:HE2	1:B:495:ILE:HD11	1.81	0.45
1:A:191:ILE:HD12	1:A:235:LEU:HD23	1.98	0.45
1:B:142:GLN:HB2	1:B:388:TRP:CZ2	2.52	0.45
1:A:530:ASP:OD1	1:B:82:LYS:NZ	2.39	0.45
1:B:421:ASN:HD21	1:B:505:GLY:H	1.65	0.44
1:B:175:GLY:HA3	1:B:201:ARG:HD2	1.99	0.44
1:B:252:VAL:HG11	1:B:311:GLN:HG2	2.00	0.44
1:B:437:ARG:NE	1:B:520:ASP:OD1	2.46	0.44
1:B:437:ARG:NH2	1:B:469:ASN:HD21	2.16	0.44
1:A:102:ASN:ND2	1:A:484:HIS:NE2	2.56	0.44
1:A:434:VAL:HG13	1:A:500:VAL:HG13	1.99	0.43
1:A:142:GLN:HB3	1:A:388:TRP:CH2	2.53	0.43
1:B:341:VAL:HG22	1:B:342:PRO:HD3	2.00	0.43
1:B:333:ASP:OD1	1:B:333:ASP:N	2.41	0.43
1:B:455:MSE:HE1	1:B:528:HIS:CD2	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:526:ILE:HD11	1:A:545:ILE:CG2	2.48	0.42
1:A:281:GLU:OE1	1:A:281:GLU:C	2.57	0.42
1:A:304:ALA:O	1:A:308[B]:MSE:HG3	2.20	0.42
1:B:333:ASP:O	1:B:334:ALA:C	2.58	0.42
1:A:460:ASP:O	1:A:461:ALA:CB	2.67	0.42
1:A:463:ALA:HB3	1:B:485:HIS:CE1	2.54	0.42
1:B:369:TYR:O	1:B:372:ASP:HB2	2.20	0.41
1:B:434:VAL:HG13	1:B:500:VAL:HG13	2.02	0.41
1:B:540:ALA:HA	1:B:545:ILE:HD12	2.01	0.41
1:A:85:ALA:CB	1:B:526:ILE:HD13	2.50	0.41
1:A:540:ALA:HA	1:A:545:ILE:HD12	2.02	0.41
1:B:257:THR:CG2	1:B:316:PHE:CE2	3.02	0.41
1:B:311:GLN:HE21	1:B:311:GLN:HA	1.86	0.41
1:A:213:SER:HB3	1:A:216:ALA:HB3	2.02	0.41
1:A:317:ASP:OD2	1:A:323:ARG:HD3	2.21	0.41
1:B:257:THR:CG2	1:B:316:PHE:CD2	3.03	0.41
1:A:142:GLN:CB	1:A:388:TRP:CH2	3.04	0.41
1:A:469:ASN:HB2	1:A:523:LEU:HD23	2.02	0.40
1:A:142:GLN:HE22	1:A:407:TRP:HE1	1.68	0.40
1:B:18:LYS:NZ	1:B:71:GLU:OE2	2.47	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	489/551 (89%)	474 (97%)	14 (3%)	1 (0%)	44	59
1	B	485/551 (88%)	463 (96%)	22 (4%)	0	100	100
All	All	974/1102 (88%)	937 (96%)	36 (4%)	1 (0%)	48	65

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	461	ALA

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	389/414 (94%)	358 (92%)	31 (8%)	10	16
1	B	383/414 (92%)	349 (91%)	34 (9%)	8	12
All	All	772/828 (93%)	707 (92%)	65 (8%)	9	14

All (65) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	58	GLU
1	A	69	ARG
1	A	70	LEU
1	A	104	VAL
1	A	131	SER
1	A	142	GLN
1	A	165	THR
1	A	187	VAL
1	A	204	THR
1	A	207	LEU
1	A	223	GLN
1	A	236	VAL
1	A	257	THR
1	A	266	LEU
1	A	281	GLU
1	A	306	LEU
1	A	310	LYS
1	A	315	THR
1	A	318	TYR
1	A	327	LYS
1	A	341	VAL
1	A	346	ARG
1	A	376	LEU

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Mol	Chain	Res	Type
1	A	404	ARG
1	A	426	LYS
1	A	440	LEU
1	A	460	ASP
1	A	499	MSE
1	A	507	LYS
1	A	514	GLU
1	A	535	LEU
1	B	15	ARG
1	B	32	ASN
1	B	44	LEU
1	B	70	LEU
1	B	102	ASN
1	B	142	GLN
1	B	179	LEU
1	B	204	THR
1	B	215	ASP
1	B	229	LYS
1	B	236	VAL
1	B	252	VAL
1	B	257	THR
1	B	261	SER
1	B	296	GLN
1	B	311	GLN
1	B	315	THR
1	B	318	TYR
1	B	323	ARG
1	B	324	GLN
1	B	333	ASP
1	B	341	VAL
1	B	374	VAL
1	B	376	LEU
1	B	377	ARG
1	B	404	ARG
1	B	415	LYS
1	B	460	ASP
1	B	491	MSE
1	B	499	MSE
1	B	519	THR
1	B	527	ARG
1	B	535	LEU
1	B	549	MSE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	102	ASN
1	A	223	GLN
1	A	289	GLN
1	A	324	GLN
1	A	473	ASN
1	B	102	ASN
1	B	142	GLN
1	B	296	GLN
1	B	309	GLN
1	B	311	GLN
1	B	324	GLN
1	B	421	ASN
1	B	469	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 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	NAD	A	600	-	42,48,48	1.98	6 (14%)	50,73,73	1.43	4 (8%)
2	NAD	B	600	-	42,48,48	1.84	3 (7%)	50,73,73	1.29	4 (8%)

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	NAD	A	600	-	-	0/26/62/62	0/5/5/5
2	NAD	B	600	-	-	3/26/62/62	0/5/5/5

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	600	NAD	O7N-C7N	9.76	1.42	1.24
2	A	600	NAD	O7N-C7N	9.41	1.41	1.24
2	A	600	NAD	PN-O3	4.24	1.64	1.59
2	A	600	NAD	PA-O3	3.60	1.63	1.59
2	A	600	NAD	C2A-N3A	3.60	1.37	1.32
2	B	600	NAD	C2A-N3A	3.46	1.37	1.32
2	B	600	NAD	C2A-N1A	2.90	1.39	1.33
2	A	600	NAD	O4B-C4B	-2.47	1.39	1.45
2	A	600	NAD	C2A-N1A	2.14	1.37	1.33

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	600	NAD	N3A-C2A-N1A	-6.69	119.59	128.67
2	A	600	NAD	N3A-C2A-N1A	-6.12	120.36	128.67
2	A	600	NAD	C4B-O4B-C1B	-3.64	106.59	109.92
2	A	600	NAD	O2A-PA-O3	3.56	116.89	107.27
2	B	600	NAD	C3N-C7N-N7N	2.36	120.65	117.74
2	B	600	NAD	O2A-PA-O3	2.29	113.47	107.27
2	A	600	NAD	C2N-C3N-C4N	2.26	120.89	118.26
2	B	600	NAD	C1B-N9A-C4A	-2.23	122.72	126.64

There are no chirality outliers.

All (3) torsion outliers are listed below:

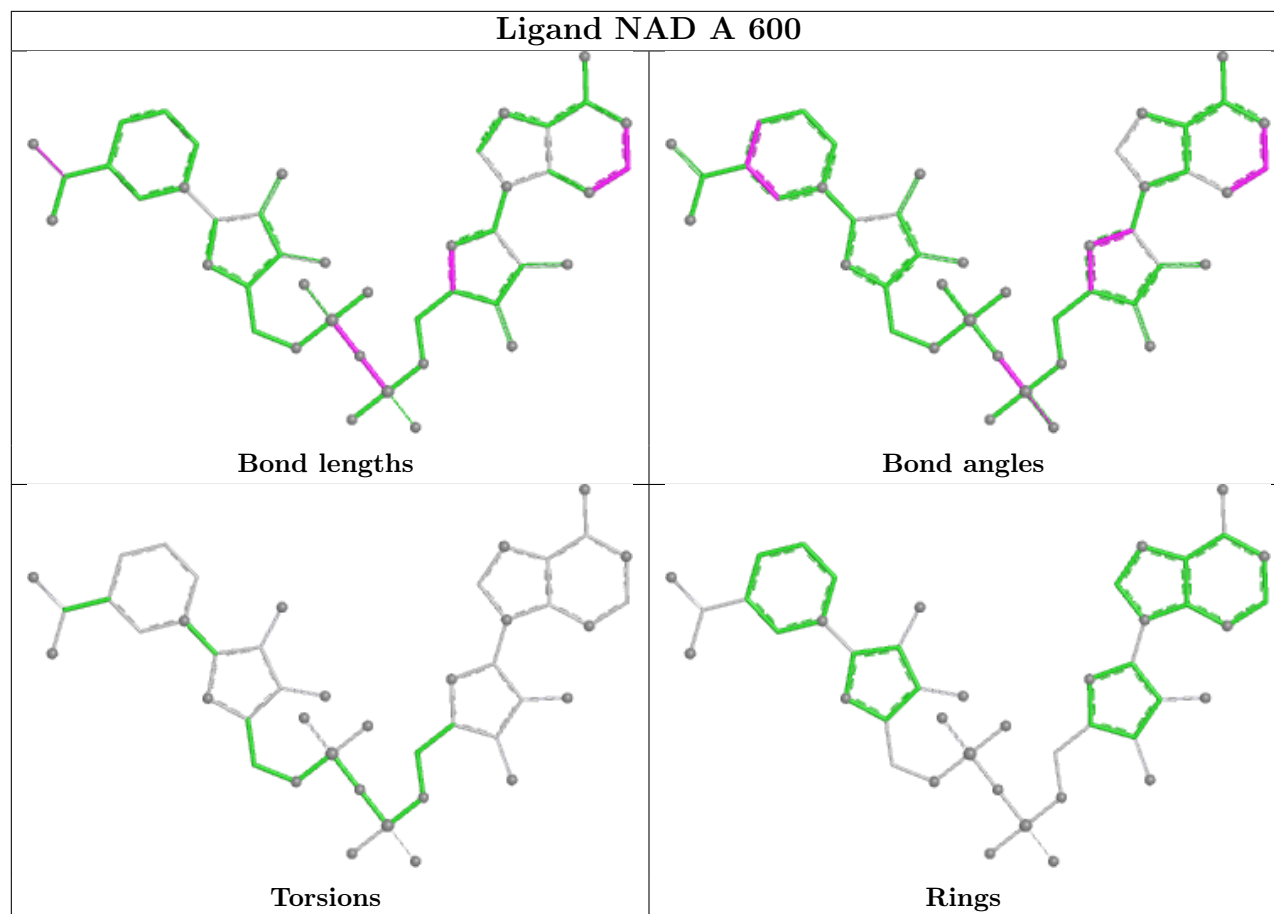
Mol	Chain	Res	Type	Atoms
2	B	600	NAD	O4D-C4D-C5D-O5D
2	B	600	NAD	C3D-C4D-C5D-O5D
2	B	600	NAD	O4B-C4B-C5B-O5B

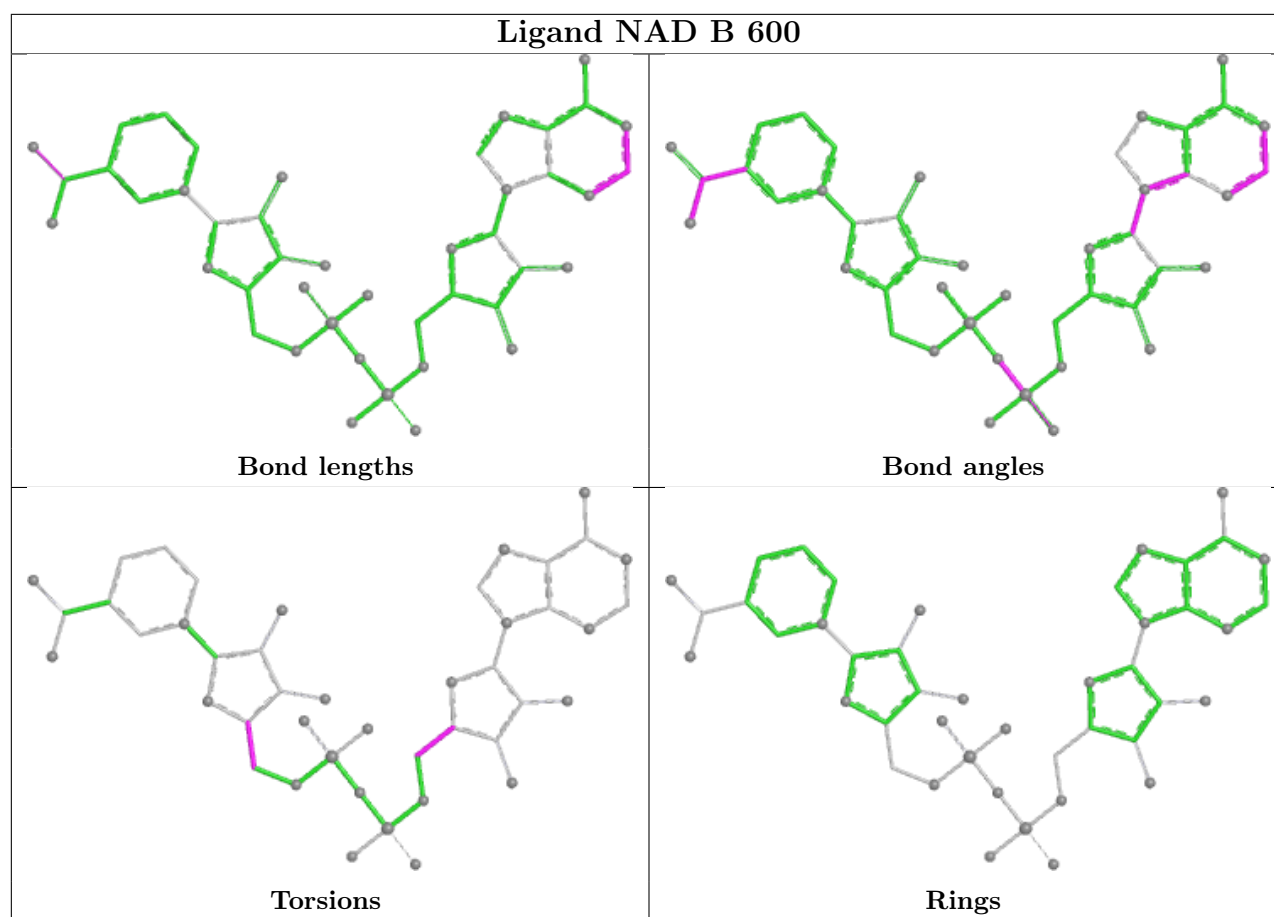
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	600	NAD	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





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

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	482/551 (87%)	-0.19	8 (1%) 69 65	31, 43, 59, 77	1 (0%)
1	B	479/551 (86%)	0.56	20 (4%) 41 38	30, 51, 70, 84	0
All	All	961/1102 (87%)	0.18	28 (2%) 54 50	30, 47, 68, 84	1 (0%)

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	129	ALA	3.5
1	B	492	GLY	3.0
1	B	44	LEU	3.0
1	B	376	LEU	3.0
1	B	334	ALA	2.9
1	B	550	LEU	2.7
1	A	110	TRP	2.6
1	B	161	ALA	2.6
1	A	36	ASP	2.6
1	A	460	ASP	2.5
1	B	284	ALA	2.5
1	B	226	GLU	2.5
1	A	37	VAL	2.4
1	A	439	HIS	2.4
1	B	11	ALA	2.4
1	B	159	THR	2.3
1	B	332	ASP	2.2
1	B	325	VAL	2.2
1	A	131	SER	2.2
1	B	326	ALA	2.2
1	A	440	LEU	2.2
1	B	296	GLN	2.1
1	B	322	ILE	2.1
1	B	172	GLY	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	292	ALA	2.0
1	A	130	GLY	2.0
1	B	337	PHE	2.0
1	B	283	ARG	2.0

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 monosaccharides in this entry.

6.4 Ligands [i](#)

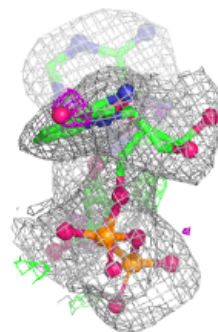
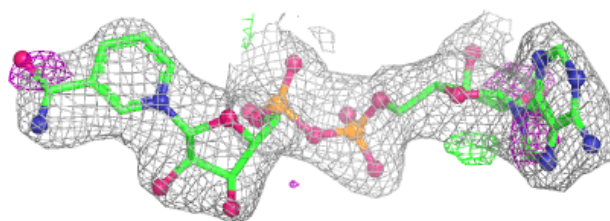
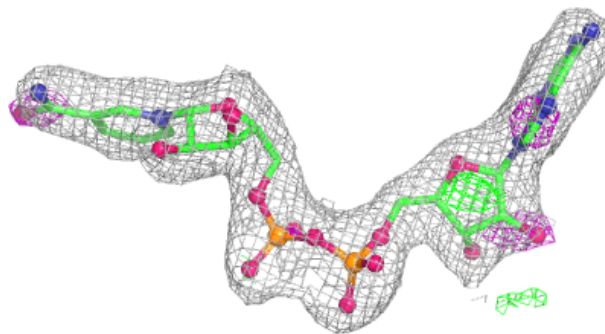
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	NAD	B	600	44/44	0.90	0.10	52,55,67,68	0
2	NAD	A	600	44/44	0.97	0.06	34,37,50,51	0

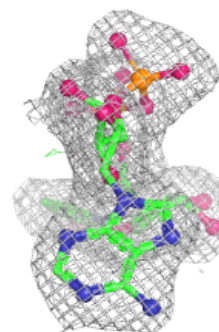
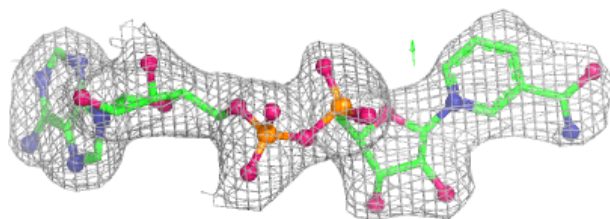
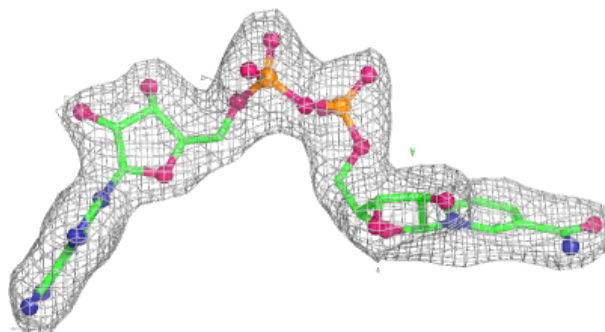
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around NAD B 600:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around NAD A 600:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.