



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 05:27 AM UTC

PDB ID : 4KQX / pdb_00004kqx
Title : Mutant Slackia exigua KARI DDV in complex with NAD and an inhibitor
Authors : Brinkmann-Chen, S.; Flock, T.; Cahn, J.K.B.; Snow, C.D.; Brustad, E.M.;
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Deposited on : 2013-05-15
Resolution : 1.80 Å(reported)

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

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

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

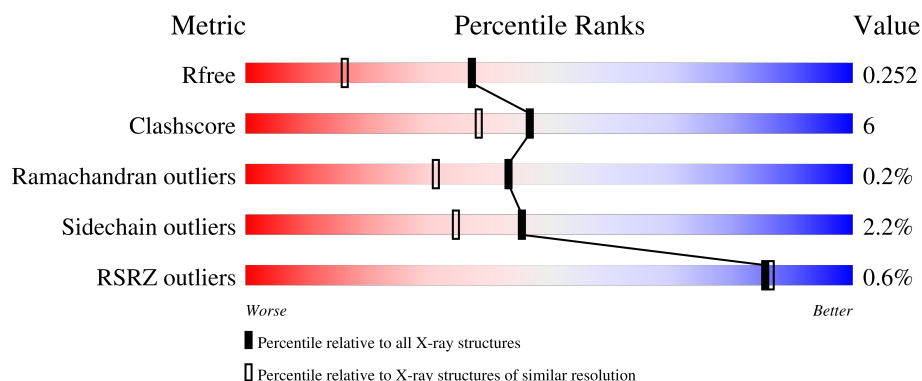
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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}	180053	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (1.80-1.80)

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	350	 73% 17% • 8%
1	B	350	 79% 15% • •

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 5497 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 Ketol-acid reductoisomerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	322	Total	C	N	O	S	0	0	0
			2491	1572	416	487	16			
1	B	335	Total	C	N	O	S	0	0	0
			2598	1641	432	508	17			

There are 22 discrepancies between the modelled and reference sequences:

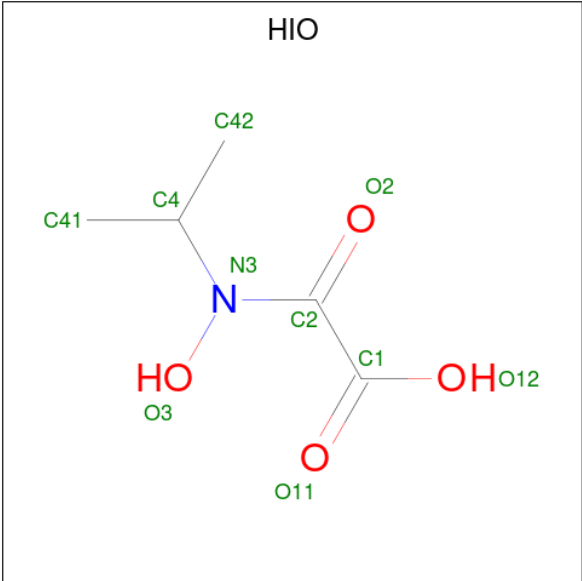
Chain	Residue	Modelled	Actual	Comment	Reference
A	61	ASP	SER	engineered mutation	UNP D0WGK0
A	63	ASP	SER	engineered mutation	UNP D0WGK0
A	95	VAL	ILE	engineered mutation	UNP D0WGK0
A	343	LEU	-	expression tag	UNP D0WGK0
A	344	GLU	-	expression tag	UNP D0WGK0
A	345	HIS	-	expression tag	UNP D0WGK0
A	346	HIS	-	expression tag	UNP D0WGK0
A	347	HIS	-	expression tag	UNP D0WGK0
A	348	HIS	-	expression tag	UNP D0WGK0
A	349	HIS	-	expression tag	UNP D0WGK0
A	350	HIS	-	expression tag	UNP D0WGK0
B	61	ASP	SER	engineered mutation	UNP D0WGK0
B	63	ASP	SER	engineered mutation	UNP D0WGK0
B	95	VAL	ILE	engineered mutation	UNP D0WGK0
B	343	LEU	-	expression tag	UNP D0WGK0
B	344	GLU	-	expression tag	UNP D0WGK0
B	345	HIS	-	expression tag	UNP D0WGK0
B	346	HIS	-	expression tag	UNP D0WGK0
B	347	HIS	-	expression tag	UNP D0WGK0
B	348	HIS	-	expression tag	UNP D0WGK0
B	349	HIS	-	expression tag	UNP D0WGK0
B	350	HIS	-	expression tag	UNP D0WGK0

- Molecule 2 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (CCD ID: NAD) (formula: $C_{21}H_{27}N_7O_{14}P_2$).



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

- Molecule 3 is N-HYDROXY-N-ISOPROPYLOXAMIC ACID (CCD ID: HIO) (formula: C₅H₉NO₄).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			10	5	1	4		

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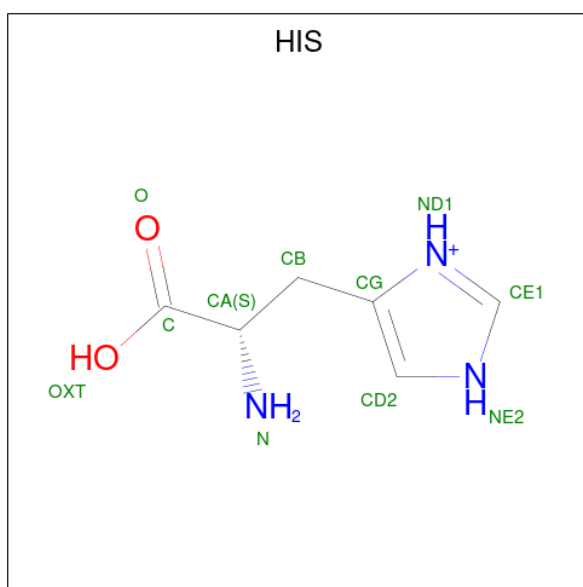
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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	1	Total	C	N	O	0	0
			10	5	1	4		

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	4	Total	Mg	0	0
			4	4		
4	B	6	Total	Mg	0	0
			6	6		

- Molecule 5 is HISTIDINE (CCD ID: HIS) (formula: C₆H₁₀N₃O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	B	1	Total	C	N	O	0	0
			10	6	3	1		

- Molecule 6 is water.

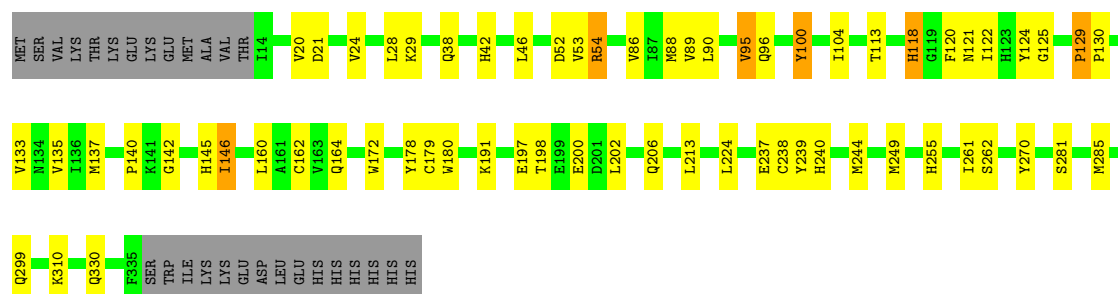
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	126	Total	O	0	0
			126	126		
6	B	154	Total	O	0	0
			154	154		

3 Residue-property plots [i](#)

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

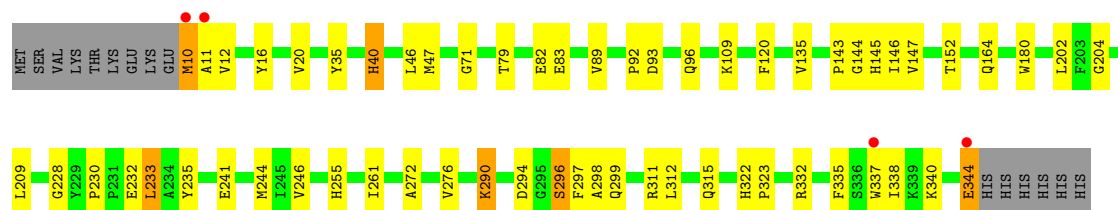
- Molecule 1: Ketol-acid reductoisomerase

Chain A: 



- Molecule 1: Ketol-acid reductoisomerase

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	49.98Å 105.35Å 122.36Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	79.83 – 1.80 79.83 – 1.80	Depositor EDS
% Data completeness (in resolution range)	98.6 (79.83-1.80) 98.6 (79.83-1.80)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.04 (at 1.81Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.201 , 0.252 0.201 , 0.252	Depositor DCC
R_{free} test set	3035 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	17.0	Xtriage
Anisotropy	0.884	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 38.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5497	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.64% 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, HIO, MG

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	1.46	18/2544 (0.7%)	1.31	11/3447 (0.3%)
1	B	1.49	12/2653 (0.5%)	1.32	9/3594 (0.3%)
All	All	1.48	30/5197 (0.6%)	1.31	20/7041 (0.3%)

All (30) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	140	PRO	C-O	8.39	1.33	1.23
1	B	135	VAL	C-O	8.24	1.33	1.24
1	B	145	HIS	CE1-NE2	7.72	1.40	1.32
1	B	145	HIS	CG-CD2	7.48	1.44	1.35
1	B	96	GLN	CA-C	6.47	1.59	1.52
1	A	42	HIS	C-O	6.39	1.31	1.24
1	B	261	ILE	CA-C	6.36	1.59	1.53
1	A	255	HIS	CG-CD2	6.20	1.42	1.35
1	A	21	ASP	C-O	6.19	1.31	1.24
1	A	129	PRO	CA-C	6.03	1.57	1.52
1	A	261	ILE	C-O	5.95	1.31	1.24
1	A	262	SER	C-O	5.95	1.31	1.23
1	A	255	HIS	ND1-CE1	5.95	1.38	1.32
1	A	240	HIS	ND1-CE1	5.85	1.38	1.32
1	B	298	ALA	N-CA	5.82	1.53	1.46
1	A	240	HIS	CG-CD2	5.76	1.42	1.35
1	B	40	HIS	N-CA	5.70	1.53	1.46
1	A	142	GLY	N-CA	5.61	1.52	1.44
1	B	322	HIS	ND1-CE1	5.58	1.38	1.32
1	B	89	VAL	CA-C	5.51	1.58	1.53
1	A	124	TYR	N-CA	5.34	1.53	1.46
1	A	172	TRP	CD2-CE2	5.18	1.50	1.41
1	A	118	HIS	CG-ND1	-5.17	1.32	1.38
1	B	152	THR	C-O	-5.10	1.17	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	118	HIS	CG-CD2	5.05	1.41	1.35
1	A	145	HIS	CG-CD2	5.01	1.41	1.35
1	A	240	HIS	N-CA	5.01	1.52	1.46
1	B	40	HIS	CE1-NE2	5.00	1.37	1.32
1	A	146	ILE	C-O	5.00	1.30	1.24
1	B	120	PHE	N-CA	5.00	1.52	1.46

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	202	LEU	N-CA-C	-7.46	103.15	111.28
1	A	238	CYS	N-CA-C	7.09	121.16	112.23
1	B	297	PHE	N-CA-C	-6.68	103.92	111.07
1	B	147	VAL	N-CA-C	-6.41	104.08	110.62
1	A	224	LEU	N-CA-C	6.38	117.90	111.07
1	A	95	VAL	N-CA-CB	-6.25	105.33	112.21
1	B	209	LEU	N-CA-C	6.05	118.83	111.82
1	A	249	MET	N-CA-C	5.87	117.68	111.28
1	A	239	TYR	N-CA-C	-5.79	104.20	112.13
1	B	276	VAL	N-CA-C	-5.78	105.45	111.58
1	B	235	TYR	N-CA-C	5.65	117.11	111.07
1	A	95	VAL	N-CA-C	-5.64	107.29	113.43
1	A	122	ILE	CB-CA-C	-5.47	104.67	112.22
1	B	228	GLY	N-CA-C	5.46	123.61	115.64
1	A	100	TYR	CA-C-N	5.30	127.38	120.28
1	A	100	TYR	C-N-CA	5.30	127.38	120.28
1	A	310	LYS	N-CA-C	5.28	116.72	111.07
1	A	121	ASN	N-CA-C	5.23	116.66	111.07
1	B	46	LEU	N-CA-C	-5.07	105.83	111.36
1	B	144	GLY	N-CA-C	-5.03	106.69	112.73

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	A	2491	0	2393	33	0
1	B	2598	0	2502	34	2
2	A	44	0	26	1	0
2	B	44	0	26	3	0
3	A	10	0	8	0	0
3	B	10	0	7	1	0
4	A	4	0	0	0	0
4	B	6	0	0	0	0
5	B	10	0	6	0	3
6	A	126	0	0	3	1
6	B	154	0	0	6	4
All	All	5497	0	4968	58	5

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 (58) 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:100:TYR:HA	1:A:104:ILE:HD12	1.65	0.77
1:A:38:GLN:HE21	2:A:401:NAD:H71N	1.41	0.68
6:A:604:HOH:O	1:B:255:HIS:HE1	1.81	0.63
1:B:79:THR:O	1:B:83:GLU:HG3	2.01	0.61
1:A:244:MET:HG2	1:B:146:ILE:HG21	1.83	0.60
1:A:89:VAL:CG1	1:A:96:GLN:HG2	2.32	0.60
1:A:86:VAL:HG22	1:A:113:THR:HB	1.84	0.59
1:A:125:GLY:HA2	6:A:552:HOH:O	2.01	0.59
1:A:129:PRO:HG2	1:A:135:VAL:HG21	1.85	0.57
1:B:10:MET:HG2	1:B:12:VAL:O	2.04	0.56
1:B:16:TYR:HB3	6:B:631:HOH:O	2.06	0.55
1:B:143:PRO:HA	2:B:402:NAD:N7N	2.22	0.55
1:A:213:LEU:HD21	1:B:246:VAL:HG22	1.89	0.55
1:A:89:VAL:HG12	1:A:96:GLN:HG2	1.91	0.53
1:A:270:TYR:CD2	1:B:315:GLN:HB2	2.43	0.52
1:A:197:GLU:O	1:A:198:THR:C	2.52	0.51
1:A:20:VAL:HG21	1:A:180:TRP:HB2	1.93	0.51
1:A:281:SER:O	1:A:285:MET:HG3	2.10	0.51
1:B:241:GLU:OE2	3:B:401:HIO:H4	2.11	0.50
1:A:146:ILE:HG21	1:B:244:MET:HG2	1.94	0.50
1:B:311:ARG:O	1:B:312:LEU:C	2.54	0.50
1:A:130:PRO:HG2	1:A:133:VAL:HG22	1.94	0.49
1:A:118:HIS:CD2	1:A:120:PHE:HB2	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:294:ASP:OD1	1:B:296:SER:OG	2.30	0.49
1:A:202:LEU:O	1:A:206:GLN:HG3	2.13	0.48
1:B:164:GLN:O	6:B:621:HOH:O	2.20	0.48
1:B:20:VAL:HG21	1:B:180:TRP:HB2	1.97	0.47
1:A:24:VAL:O	1:A:28:LEU:HD12	2.15	0.47
1:A:88:MET:HG3	1:A:90:LEU:HG	1.96	0.47
1:A:200:GLU:CD	1:B:233:LEU:HD11	2.39	0.47
2:B:402:NAD:H5N	6:B:516:HOH:O	2.15	0.47
1:B:290:LYS:HG3	6:B:637:HOH:O	2.15	0.46
1:B:344:GLU:OE2	1:B:344:GLU:HA	2.13	0.46
1:A:191:LYS:NZ	6:A:594:HOH:O	2.49	0.45
1:A:29:LYS:HG2	1:A:52:ASP:OD2	2.17	0.45
1:A:200:GLU:OE2	1:A:200:GLU:C	2.60	0.45
1:B:335:PHE:CD2	1:B:337:TRP:CZ2	3.04	0.45
1:B:255:HIS:HD2	6:B:627:HOH:O	2.00	0.45
1:B:230:PRO:HD2	1:B:233:LEU:HD23	2.00	0.44
1:B:35:TYR:CE1	1:B:40:HIS:HA	2.53	0.43
1:A:200:GLU:HG2	1:B:233:LEU:HD11	2.00	0.43
1:A:179:CYS:O	1:A:180:TRP:C	2.62	0.43
1:B:332:ARG:O	1:B:338:ILE:HD12	2.19	0.42
1:B:47:MET:HE1	1:B:71:GLY:O	2.19	0.42
1:B:311:ARG:NH1	6:B:630:HOH:O	2.37	0.42
1:A:237:GLU:OE2	1:B:204:GLY:HA3	2.20	0.42
1:B:232:GLU:OE1	1:B:232:GLU:N	2.50	0.42
1:B:82:GLU:OE2	1:B:109:LYS:NZ	2.53	0.42
1:A:206:GLN:HB3	1:B:272:ALA:HB2	2.01	0.42
1:A:53:VAL:HG12	1:A:54:ARG:N	2.35	0.42
1:B:340:LYS:HA	1:B:340:LYS:HD3	1.80	0.41
1:B:143:PRO:HA	2:B:402:NAD:H71N	1.85	0.41
1:A:46:LEU:HD11	1:A:178:TYR:CE2	2.56	0.41
1:A:137:MET:HE3	1:A:162:CYS:SG	2.60	0.41
1:A:164:GLN:HA	1:A:164:GLN:OE1	2.20	0.41
1:A:200:GLU:CG	1:B:233:LEU:HD11	2.51	0.41
1:B:92:PRO:O	1:B:93:ASP:C	2.64	0.41
1:A:244:MET:CG	1:B:146:ILE:HG21	2.50	0.40

All (5) 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)
5:B:403:HIS:C	6:A:561:HOH:O[1_655]	1.67	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:B:403:HIS:O	6:B:646:HOH:O[1_655]	1.93	0.27
1:B:344:GLU:C	6:B:548:HOH:O[1_655]	1.99	0.21
1:B:344:GLU:O	6:B:548:HOH:O[1_655]	2.11	0.09
5:B:403:HIS:N	6:B:548:HOH:O[1_655]	2.19	0.01

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	320/350 (91%)	309 (97%)	11 (3%)	0	100	100
1	B	333/350 (95%)	317 (95%)	15 (4%)	1 (0%)	36	25
All	All	653/700 (93%)	626 (96%)	26 (4%)	1 (0%)	43	31

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	11	ALA

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	261/288 (91%)	256 (98%)	5 (2%)	50	41
1	B	273/288 (95%)	266 (97%)	7 (3%)	40	28
All	All	534/576 (93%)	522 (98%)	12 (2%)	45	34

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	54	ARG
1	A	95	VAL
1	A	160	LEU
1	A	299	GLN
1	A	330	GLN
1	B	10	MET
1	B	233	LEU
1	B	290	LYS
1	B	296	SER
1	B	299	GLN
1	B	323	PRO
1	B	344	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	38	GLN
1	B	255	HIS
1	B	309	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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 15 ligands modelled in this entry, 10 are monoatomic - leaving 5 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
2	NAD	A	401	4	46,48,48	1.52	9 (19%)	64,73,73	1.79	16 (25%)
3	HIO	B	401	4	8,9,9	1.90	2 (25%)	5,12,12	3.18	4 (80%)
5	HIS	B	403	-	9,10,11	1.07	1 (11%)	8,12,14	0.97	0
2	NAD	B	402	4	46,48,48	1.77	7 (15%)	64,73,73	1.88	21 (32%)
3	HIO	A	402	4	8,9,9	1.77	2 (25%)	5,12,12	4.23	3 (60%)

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	401	4	-	7/30/62/62	0/5/5/5
3	HIO	B	401	4	-	5/9/12/12	-
5	HIS	B	403	-	-	1/5/6/8	0/1/1/1
2	NAD	B	402	4	-	3/30/62/62	0/5/5/5
3	HIO	A	402	4	-	5/9/12/12	-

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	402	NAD	O4D-C1D	6.12	1.48	1.40
2	A	401	NAD	C5A-C4A	5.70	1.49	1.39
2	B	402	NAD	C5A-C4A	4.86	1.47	1.39
2	B	402	NAD	PA-O3	4.69	1.64	1.59
2	B	402	NAD	PN-O3	3.89	1.63	1.59
3	B	401	HIO	C2-N3	3.73	1.38	1.34
2	A	401	NAD	PN-O3	3.02	1.62	1.59
2	A	401	NAD	PA-O3	3.00	1.62	1.59
3	A	402	HIO	O2-C2	2.86	1.28	1.23
2	B	402	NAD	C5A-C6A	2.84	1.48	1.41
2	B	402	NAD	O2D-C2D	2.47	1.49	1.43
2	A	401	NAD	C5D-C4D	2.32	1.58	1.51
5	B	403	HIS	CD2-CG	2.30	1.41	1.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	NAD	C4A-N9A	-2.23	1.33	1.37
2	A	401	NAD	C5A-C6A	2.22	1.47	1.41
3	A	402	HIO	O12-C1	-2.19	1.24	1.30
2	B	402	NAD	C7N-N7N	-2.18	1.28	1.33
2	A	401	NAD	C5A-N7A	-2.17	1.35	1.39
3	B	401	HIO	O2-C2	2.07	1.27	1.23
2	A	401	NAD	C2N-N1N	2.03	1.37	1.35
2	A	401	NAD	O4D-C1D	2.03	1.43	1.40

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	402	HIO	C42-C4-N3	-8.67	100.67	109.96
2	A	401	NAD	C3N-C7N-N7N	-5.03	111.54	117.74
2	B	402	NAD	C4A-N9A-C8A	4.62	110.58	105.74
3	B	401	HIO	C41-C4-N3	4.26	114.53	109.96
2	A	401	NAD	N3A-C4A-N9A	4.22	134.35	127.17
2	B	402	NAD	N3A-C4A-N9A	3.97	133.92	127.17
2	A	401	NAD	C5A-C4A-N3A	-3.91	121.34	126.72
2	A	401	NAD	C4A-N9A-C8A	3.86	109.79	105.74
2	A	401	NAD	N3A-C2A-N1A	-3.82	122.80	128.58
3	B	401	HIO	C42-C4-N3	3.81	114.05	109.96
2	B	402	NAD	N3A-C2A-N1A	-3.77	122.87	128.58
2	B	402	NAD	C5A-C4A-N3A	-3.74	121.56	126.72
2	B	402	NAD	C2A-N3A-C4A	3.57	120.55	111.83
2	A	401	NAD	O2A-PA-O1A	3.54	128.89	112.44
2	B	402	NAD	O2A-PA-O1A	3.45	128.50	112.44
2	A	401	NAD	C2A-N1A-C6A	3.43	124.36	118.73
3	B	401	HIO	C42-C4-C41	3.37	121.62	112.46
2	B	402	NAD	O7N-C7N-C3N	3.26	123.58	119.60
2	B	402	NAD	O3D-C3D-C4D	-3.21	101.88	111.08
2	B	402	NAD	C6A-C5A-N7A	3.19	138.24	132.09
2	B	402	NAD	O2A-PA-O3	-3.17	98.71	107.27
2	B	402	NAD	N9A-C8A-N7A	-3.11	109.52	113.94
2	A	401	NAD	O2N-PN-O1N	2.93	126.06	112.44
2	A	401	NAD	C2A-N3A-C4A	2.75	118.55	111.83
2	B	402	NAD	O3B-C3B-C4B	-2.74	103.22	111.08
2	B	402	NAD	C5A-N7A-C8A	2.72	107.72	103.45
3	A	402	HIO	O12-C1-O11	-2.70	117.47	123.90
2	B	402	NAD	C4A-C5A-N7A	-2.56	107.65	110.58
2	A	401	NAD	O4B-C1B-N9A	2.55	112.99	108.09
2	A	401	NAD	O2N-PN-O3	2.46	113.94	107.27

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	402	NAD	C4D-O4D-C1D	-2.33	107.80	109.92
2	A	401	NAD	C5A-N7A-C8A	2.32	107.09	103.45
3	B	401	HIO	O12-C1-O11	-2.30	118.42	123.90
2	A	401	NAD	C4A-C5A-N7A	-2.27	107.99	110.58
2	B	402	NAD	C6A-C5A-C4A	-2.25	114.10	117.18
2	B	402	NAD	O2B-C2B-C3B	2.24	119.00	111.82
2	B	402	NAD	C5D-C4D-C3D	-2.21	107.25	115.21
2	A	401	NAD	N9A-C8A-N7A	-2.20	110.81	113.94
2	A	401	NAD	O7N-C7N-N7N	2.16	125.74	122.62
2	A	401	NAD	O3-PA-O1A	-2.10	104.40	110.70
2	B	402	NAD	C4A-N9A-C1B	-2.08	121.78	126.63
3	A	402	HIO	C41-C4-N3	2.06	112.17	109.96
2	B	402	NAD	O4B-C1B-N9A	2.03	112.00	108.09
2	B	402	NAD	C2A-N1A-C6A	2.03	122.07	118.73

There are no chirality outliers.

All (21) torsion outliers are listed below:

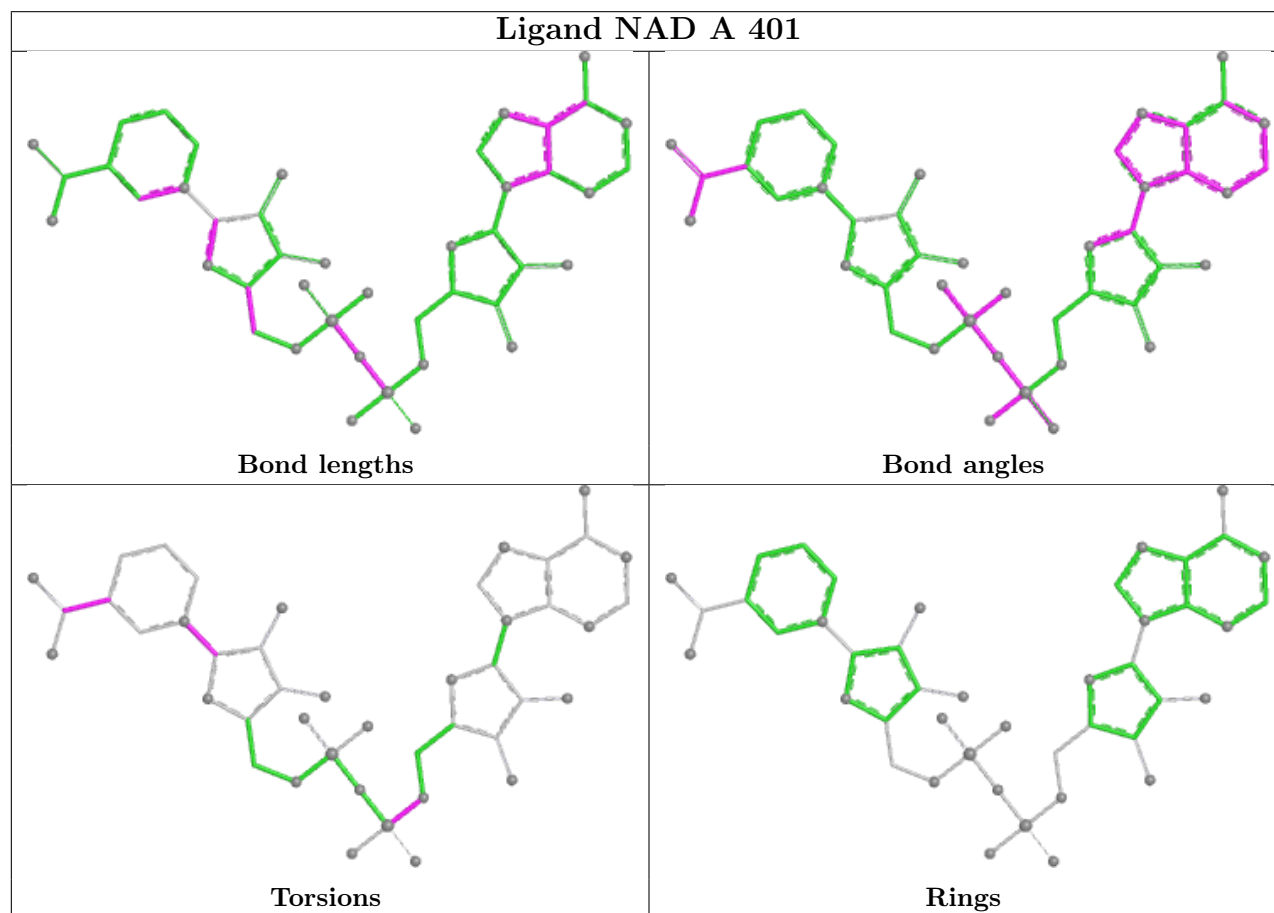
Mol	Chain	Res	Type	Atoms
2	A	401	NAD	C5B-O5B-PA-O2A
2	A	401	NAD	C5B-O5B-PA-O3
2	A	401	NAD	O4D-C1D-N1N-C2N
2	A	401	NAD	C2D-C1D-N1N-C6N
2	B	402	NAD	O4D-C1D-N1N-C2N
3	A	402	HIO	O11-C1-C2-N3
3	A	402	HIO	O12-C1-C2-O2
3	A	402	HIO	O2-C2-N3-O3
3	B	401	HIO	O12-C1-C2-O2
3	B	401	HIO	C41-C4-N3-O3
3	B	401	HIO	C42-C4-N3-O3
5	B	403	HIS	O-C-CA-CB
3	B	401	HIO	C42-C4-N3-C2
3	B	401	HIO	O2-C2-N3-O3
2	B	402	NAD	C5B-O5B-PA-O1A
3	A	402	HIO	O11-C1-C2-O2
3	A	402	HIO	O2-C2-N3-C4
2	A	401	NAD	C2N-C3N-C7N-N7N
2	A	401	NAD	C2D-C1D-N1N-C2N
2	A	401	NAD	O4D-C1D-N1N-C6N
2	B	402	NAD	O4D-C1D-N1N-C6N

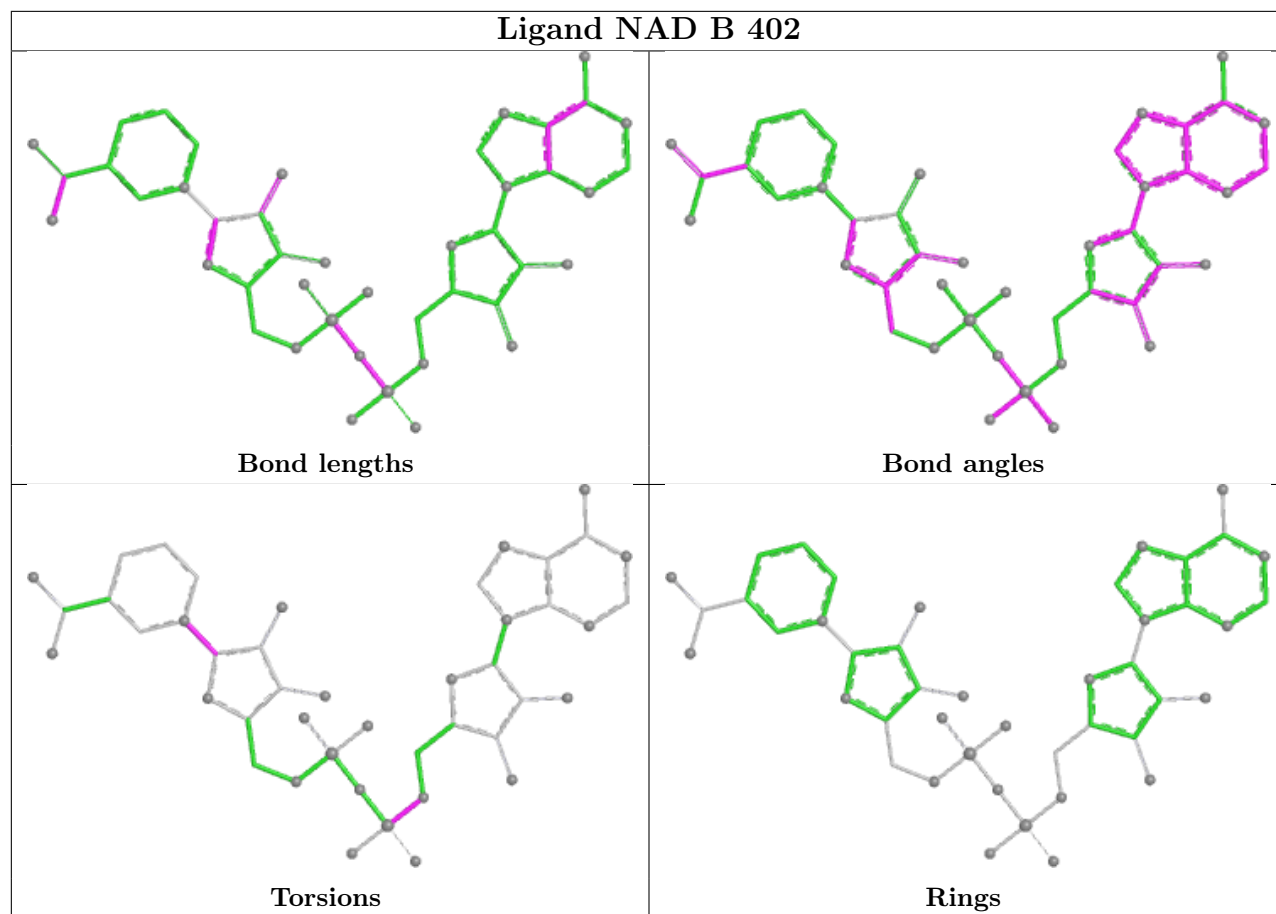
There are no ring outliers.

4 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	NAD	1	0
3	B	401	HIO	1	0
5	B	403	HIS	0	3
2	B	402	NAD	3	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 [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	322/350 (92%)	-0.08	0 100 100	10, 20, 33, 46	0
1	B	335/350 (95%)	-0.27	4 (1%) 76 77	11, 17, 28, 75	0
All	All	657/700 (93%)	-0.17	4 (0%) 85 86	10, 18, 31, 75	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	11	ALA	5.2
1	B	337	TRP	4.0
1	B	10	MET	3.1
1	B	344	GLU	2.2

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 oligosaccharides 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.

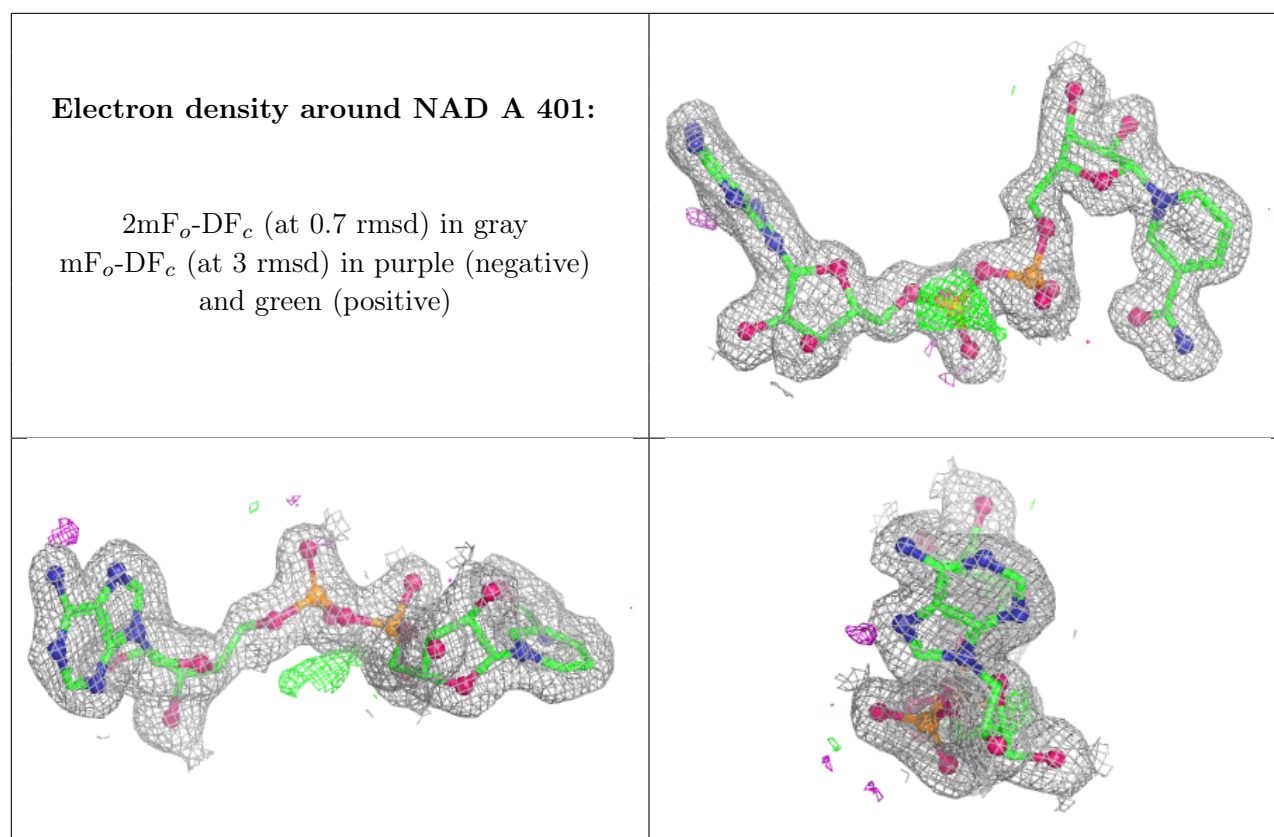
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	HIS	B	403	10/11	0.87	0.13	32,45,50,53	0

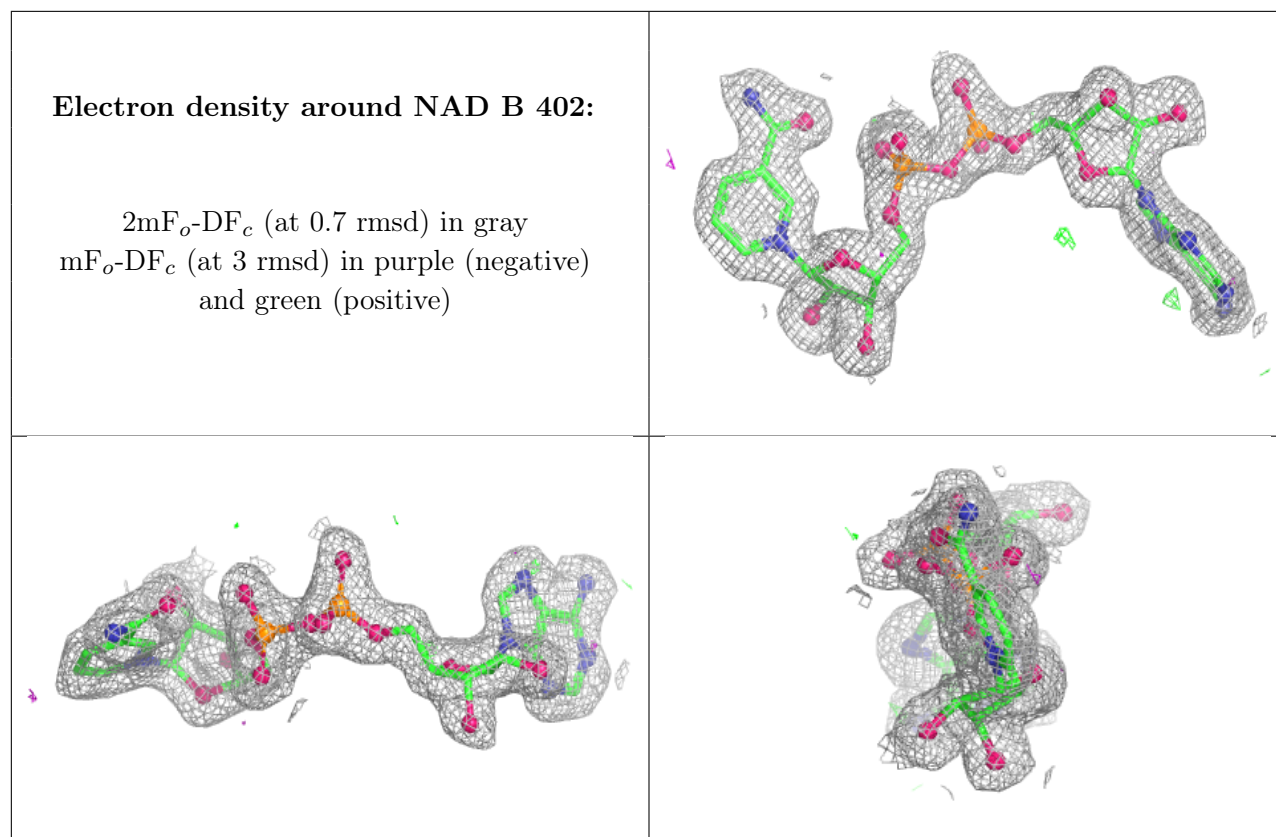
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	MG	A	406	1/1	0.93	0.05	12,12,12,12	0
3	HIO	B	401	10/10	0.93	0.08	13,18,19,20	0
3	HIO	A	402	10/10	0.94	0.07	14,16,18,23	0
2	NAD	A	401	44/44	0.96	0.07	16,21,25,26	0
4	MG	B	409	1/1	0.96	0.05	16,16,16,16	0
2	NAD	B	402	44/44	0.96	0.06	14,20,22,23	0
4	MG	A	404	1/1	0.97	0.05	11,11,11,11	0
4	MG	B	406	1/1	0.98	0.03	9,9,9,9	0
4	MG	B	407	1/1	0.98	0.04	15,15,15,15	0
4	MG	B	404	1/1	0.98	0.08	9,9,9,9	0
4	MG	B	405	1/1	0.98	0.07	8,8,8,8	0
4	MG	B	408	1/1	0.99	0.06	23,23,23,23	0
4	MG	A	405	1/1	0.99	0.02	20,20,20,20	0
4	MG	A	403	1/1	0.99	0.11	9,9,9,9	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.





6.5 Other polymers [i](#)

There are no such residues in this entry.