



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2026 – 09:16 PM UTC

PDB ID : 3FMX / pdb_00003fmx
Title : Crystal structure of Tartrate dehydrogenase from Pseudomonas putida complexed with NADH
Authors : Malik, R.; Viola, R.E.
Deposited on : 2008-12-22
Resolution : 2.95 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

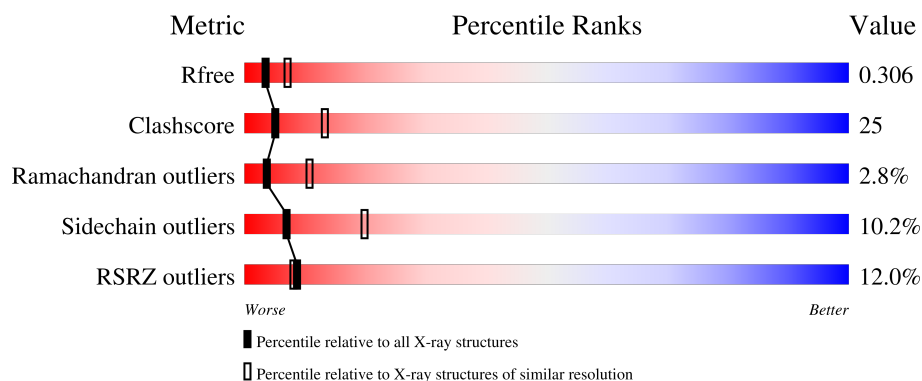
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.95 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	1130 (2.98-2.94)
Clashscore	190562	1157 (2.98-2.94)
Ramachandran outliers	187476	1101 (2.98-2.94)
Sidechain outliers	187428	1101 (2.98-2.94)
RSRZ outliers	180081	1130 (2.98-2.94)

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	X	364	

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 2875 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.

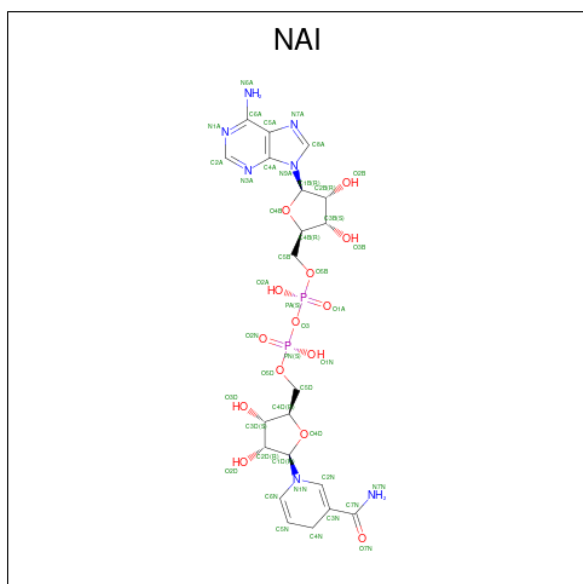
- Molecule 1 is a protein called Tartrate dehydrogenase/decarboxylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	X	359	2813	1792	484	523	14	24	0	0

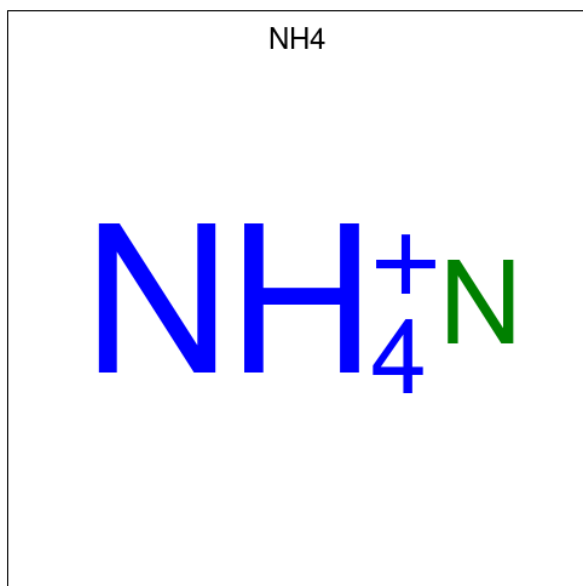
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
X	79	GLY	ASP	engineered mutation	UNP Q51945
X	?	-	VAL	deletion	UNP Q51945

- Molecule 2 is 1,4-DIHYDRONICOTINAMIDE ADENINE DINUCLEOTIDE (CCD ID: NAI) (formula: C₂₁H₂₉N₇O₁₄P₂).



- Molecule 3 is AMMONIUM ION (CCD ID: NH4) (formula: H_4N).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	X	1	Total	N	0	0
			1	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	X	17	Total	O	0	0
			17	17		

4 Data and refinement statistics

Property	Value	Source
Space group	P 6 2 2	Depositor
Cell constants a, b, c, α , β , γ	182.91Å 182.91Å 66.98Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	50.00 – 2.95 50.00 – 2.95	Depositor EDS
% Data completeness (in resolution range)	85.0 (50.00-2.95) 84.9 (50.00-2.95)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.54 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.260 , 0.297 0.259 , 0.306	Depositor DCC
R_{free} test set	696 reflections (4.14%)	wwPDB-VP
Wilson B-factor (Å ²)	52.8	Xtriage
Anisotropy	0.082	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 55.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.42$, $\langle L^2 \rangle = 0.25$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.87	EDS
Total number of atoms	2875	wwPDB-VP
Average B, all atoms (Å ²)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.37% 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: NAI, NH4

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	X	0.90	4/2882 (0.1%)	1.10	16/3911 (0.4%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	X	183	GLU	CA-CB	-27.89	1.10	1.53
1	X	124	LYS	CA-CB	-15.01	1.29	1.53
1	X	318	GLU	CA-CB	-7.46	1.41	1.53
1	X	108	ARG	CA-CB	-6.27	1.45	1.53

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	X	124	LYS	CB-CA-C	18.00	136.84	109.84
1	X	124	LYS	N-CA-CB	-13.36	88.77	109.48
1	X	183	GLU	CB-CA-C	9.34	125.55	110.88
1	X	318	GLU	N-CA-C	-6.78	104.48	112.89
1	X	122	ASN	N-CA-C	6.41	124.44	110.80
1	X	244	ALA	N-CA-C	5.50	117.21	108.79
1	X	5	SER	N-CA-CB	5.44	119.75	110.50
1	X	347	THR	N-CA-C	-5.41	106.73	113.38
1	X	176	PHE	N-CA-C	-5.40	106.88	112.93
1	X	182	ARG	N-CA-C	5.39	116.47	108.60
1	X	207	THR	N-CA-C	-5.29	105.20	110.97
1	X	183	GLU	N-CA-CB	5.21	117.56	110.01
1	X	341	THR	CA-C-N	5.17	126.30	119.84
1	X	341	THR	C-N-CA	5.17	126.30	119.84
1	X	272	ASN	CA-C-N	5.14	125.43	119.47
1	X	272	ASN	C-N-CA	5.14	125.43	119.47

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	X	2813	0	2746	137	0
2	X	44	0	27	7	0
3	X	1	0	0	1	0
4	X	17	0	0	0	0
All	All	2875	0	2773	138	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 25.

All (138) 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:X:251:ILE:HG23	1:X:252:LEU:H	0.93	1.08
1:X:21:LEU:HD12	1:X:302:MET:CE	1.86	1.05
1:X:251:ILE:CG2	1:X:252:LEU:H	1.74	0.97
1:X:251:ILE:HG23	1:X:252:LEU:N	1.73	0.97
1:X:21:LEU:HA	1:X:302:MET:HE2	1.45	0.96
1:X:21:LEU:HD12	1:X:302:MET:HE1	1.49	0.93
1:X:274:GLU:HB2	1:X:276:ASN:HD21	1.36	0.91
1:X:340:VAL:HB	1:X:344:MET:HE3	1.56	0.87
1:X:134:ARG:HG3	1:X:249:GLY:HA3	1.55	0.86
1:X:120:LEU:HG	1:X:260:ALA:HB2	1.57	0.86
1:X:74:TYR:OH	1:X:302:MET:O	1.99	0.81
1:X:309:MET:O	1:X:313:LEU:HB2	1.85	0.77
1:X:274:GLU:HB2	1:X:276:ASN:ND2	1.99	0.76
1:X:47:ALA:HB2	1:X:59:MET:CE	2.16	0.76
1:X:222:GLN:NE2	1:X:230:ARG:HD2	2.01	0.75
1:X:251:ILE:CG2	1:X:252:LEU:N	2.35	0.75
1:X:106:ASN:HB2	1:X:136:ASN:ND2	2.02	0.75
1:X:331:ILE:O	1:X:335:ILE:HG12	1.90	0.71
1:X:325:ASP:OD2	1:X:325:ASP:N	2.23	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:X:334:VAL:HG13	1:X:339:SER:HB2	1.73	0.70
1:X:130:PHE:HB2	1:X:239:ASP:O	1.92	0.70
1:X:119:ALA:HB3	1:X:260:ALA:HA	1.73	0.70
1:X:21:LEU:CD1	1:X:302:MET:HE1	2.22	0.68
1:X:123:ARG:HG2	1:X:128:ILE:HD11	1.76	0.67
1:X:165:ARG:HE	1:X:206:ARG:NH2	1.93	0.66
1:X:6:PHE:HB2	1:X:38:ALA:O	1.99	0.63
1:X:95:LEU:O	1:X:99:ARG:HB2	1.99	0.62
1:X:65:GLU:CD	1:X:65:GLU:H	2.08	0.61
1:X:231:PHE:HE2	1:X:256:GLY:HA3	1.65	0.61
1:X:21:LEU:CD1	1:X:302:MET:CE	2.73	0.61
1:X:134:ARG:CG	1:X:249:GLY:HA3	2.27	0.61
1:X:107:ILE:HD11	1:X:271:LEU:HD22	1.83	0.59
1:X:21:LEU:CA	1:X:302:MET:HE2	2.28	0.59
1:X:196:MET:HG3	1:X:200:MET:HG2	1.85	0.59
1:X:136:ASN:ND2	1:X:270:ASN:OD1	2.35	0.58
1:X:231:PHE:HA	1:X:238:PHE:HE2	1.69	0.58
2:X:401:NAI:O2A	2:X:401:NAI:H51N	2.05	0.56
1:X:107:ILE:HA	1:X:132:VAL:O	2.05	0.56
1:X:178:LEU:O	1:X:182:ARG:HD3	2.05	0.56
1:X:340:VAL:HG23	1:X:346:GLY:H	1.71	0.56
1:X:174:TYR:O	1:X:174:TYR:HD2	1.89	0.55
1:X:263:ILE:HD11	2:X:401:NAI:O4B	2.06	0.55
1:X:90:LEU:O	1:X:95:LEU:HB2	2.07	0.55
1:X:118:CYS:HA	1:X:260:ALA:HB1	1.88	0.55
1:X:112:LEU:HD21	1:X:125:VAL:HG22	1.88	0.55
1:X:333:ARG:HE	1:X:364:LEU:HD11	1.71	0.55
1:X:173:LYS:O	1:X:175:ALA:N	2.32	0.55
1:X:59:MET:HB2	1:X:60:PRO:HD2	1.88	0.54
1:X:262:THR:HG23	1:X:264:GLY:H	1.73	0.54
1:X:231:PHE:HE2	1:X:256:GLY:CA	2.20	0.54
1:X:247:LEU:C	1:X:249:GLY:H	2.15	0.54
1:X:134:ARG:NH2	1:X:250:ASP:HB2	2.22	0.54
1:X:317:ASP:HB3	1:X:319:ARG:HB3	1.89	0.53
1:X:222:GLN:HE22	1:X:230:ARG:HD2	1.73	0.53
1:X:269:ALA:HB1	1:X:271:LEU:CD1	2.39	0.53
1:X:101:PHE:CE1	1:X:278:PRO:HG2	2.44	0.52
1:X:21:LEU:HD12	1:X:302:MET:HE2	1.84	0.52
1:X:106:ASN:HB3	1:X:134:ARG:HB3	1.91	0.52
1:X:147:ILE:HD13	1:X:150:GLU:HG3	1.91	0.52
1:X:259:CYS:SG	1:X:260:ALA:N	2.83	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:X:333:ARG:HH21	1:X:364:LEU:HD21	1.75	0.51
1:X:334:VAL:HA	1:X:337:ASP:HB2	1.93	0.51
1:X:231:PHE:HA	1:X:238:PHE:CE2	2.46	0.51
1:X:340:VAL:HB	1:X:344:MET:CE	2.37	0.51
1:X:91:TRP:HA	1:X:95:LEU:HB2	1.92	0.51
1:X:244:ALA:HB1	1:X:248:PHE:HB2	1.91	0.51
1:X:249:GLY:C	1:X:251:ILE:H	2.18	0.51
1:X:267:PRO:HB3	1:X:304:TRP:HB3	1.94	0.50
1:X:174:TYR:O	1:X:174:TYR:CD2	2.64	0.50
1:X:12:PRO:O	1:X:17:GLY:HA3	2.12	0.50
1:X:263:ILE:HD11	2:X:401:NAI:C8A	2.42	0.50
1:X:231:PHE:CE2	1:X:256:GLY:HA3	2.45	0.49
1:X:247:LEU:O	1:X:251:ILE:HB	2.12	0.49
1:X:113:PHE:O	1:X:114:PRO:C	2.55	0.49
1:X:255:LEU:O	1:X:259:CYS:N	2.46	0.49
1:X:341:THR:HG23	1:X:353:VAL:HG21	1.94	0.49
1:X:63:TRP:HE3	1:X:67:LEU:HD12	1.78	0.49
1:X:63:TRP:O	1:X:66:GLN:HB3	2.13	0.48
1:X:5:SER:HA	1:X:38:ALA:HB1	1.94	0.48
1:X:97:PHE:O	1:X:101:PHE:HB2	2.13	0.48
1:X:118:CYS:C	1:X:120:LEU:H	2.22	0.48
1:X:126:GLY:C	1:X:128:ILE:H	2.22	0.48
1:X:251:ILE:O	1:X:253:SER:N	2.47	0.48
1:X:61:ASP:N	1:X:61:ASP:OD1	2.45	0.48
1:X:147:ILE:O	1:X:147:ILE:HG23	2.14	0.47
1:X:78:VAL:O	2:X:401:NAI:H1D	2.14	0.47
1:X:192:LYS:H	1:X:200:MET:HG3	1.79	0.47
1:X:230:ARG:HB3	1:X:238:PHE:HZ	1.81	0.46
1:X:231:PHE:CE2	1:X:256:GLY:CA	2.98	0.46
1:X:174:TYR:CE1	1:X:273:PRO:HG3	2.50	0.46
1:X:107:ILE:HG13	1:X:269:ALA:HB3	1.98	0.46
1:X:134:ARG:CD	1:X:249:GLY:HA3	2.45	0.46
1:X:173:LYS:C	1:X:175:ALA:H	2.19	0.46
1:X:222:GLN:HB2	1:X:227:LEU:HD13	1.97	0.46
1:X:249:GLY:C	1:X:251:ILE:N	2.74	0.46
1:X:283:PRO:O	3:X:402:NH4:N	2.49	0.46
1:X:257:PRO:HB3	1:X:266:ALA:HB2	1.99	0.45
1:X:144:LEU:HD12	1:X:162:ILE:HD11	1.99	0.45
1:X:165:ARG:O	1:X:169:ASP:HB2	2.17	0.45
1:X:286:GLY:N	2:X:401:NAI:O1A	2.35	0.45
1:X:120:LEU:HD21	1:X:259:CYS:SG	2.56	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:X:31:ALA:HA	1:X:34:LYS:HB3	1.98	0.45
1:X:113:PHE:O	1:X:116:VAL:HG23	2.17	0.44
1:X:174:TYR:HE1	1:X:273:PRO:HG3	1.83	0.44
1:X:140:GLU:HB2	1:X:247:LEU:HB2	2.00	0.44
1:X:78:VAL:O	2:X:401:NAI:H2N	2.18	0.44
1:X:251:ILE:O	1:X:252:LEU:C	2.61	0.44
1:X:175:ALA:C	1:X:177:ASP:H	2.26	0.43
1:X:169:ASP:OD2	1:X:206:ARG:NH2	2.52	0.43
1:X:16:ILE:HA	1:X:19:GLU:HB2	1.99	0.43
1:X:247:LEU:C	1:X:249:GLY:N	2.77	0.43
1:X:317:ASP:HB3	1:X:319:ARG:CB	2.48	0.43
1:X:314:GLY:C	1:X:315:GLN:HG3	2.44	0.43
1:X:106:ASN:HA	1:X:269:ALA:O	2.18	0.43
1:X:179:ALA:C	1:X:181:LYS:H	2.27	0.43
1:X:310:LEU:HD11	1:X:327:MET:HE1	2.00	0.43
1:X:256:GLY:C	1:X:258:ALA:N	2.78	0.42
1:X:221:LYS:HE2	1:X:221:LYS:HB3	1.83	0.42
1:X:95:LEU:HD23	1:X:95:LEU:HA	1.78	0.42
1:X:315:GLN:HB2	1:X:316:GLY:H	1.64	0.42
1:X:318:GLU:O	1:X:322:ARG:HG3	2.20	0.42
1:X:101:PHE:CB	1:X:103:GLN:HE21	2.33	0.42
1:X:298:ASN:HB3	2:X:401:NAI:H62A	1.84	0.42
1:X:113:PHE:HA	1:X:114:PRO:HD2	1.88	0.41
1:X:13:GLY:HA2	1:X:47:ALA:O	2.20	0.41
1:X:103:GLN:HA	1:X:271:LEU:O	2.20	0.41
1:X:196:MET:HE3	1:X:196:MET:HB3	1.97	0.41
1:X:333:ARG:HB2	1:X:360:THR:HG21	2.02	0.41
1:X:94:LEU:HG	1:X:98:ARG:HE	1.86	0.41
1:X:158:ILE:HD13	1:X:158:ILE:HA	1.84	0.41
1:X:291:ILE:O	1:X:291:ILE:CG2	2.67	0.41
1:X:255:LEU:O	1:X:258:ALA:C	2.64	0.41
1:X:303:ILE:O	1:X:304:TRP:C	2.64	0.41
1:X:300:ILE:O	1:X:304:TRP:N	2.48	0.40
1:X:267:PRO:HG3	1:X:304:TRP:CD2	2.56	0.40
1:X:262:THR:HG22	1:X:265:ILE:HG12	2.04	0.40
1:X:101:PHE:HB2	1:X:103:GLN:HE21	1.86	0.40
1:X:328:LEU:HD12	1:X:328:LEU:HA	1.96	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	X	357/364 (98%)	296 (83%)	51 (14%)	10 (3%)	4	11

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	X	122	ASN
1	X	174	TYR
1	X	251	ILE
1	X	252	LEU
1	X	253	SER
1	X	261	GLY
1	X	114	PRO
1	X	153	GLU
1	X	250	ASP
1	X	257	PRO

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	X	295/299 (99%)	265 (90%)	30 (10%)	7	20

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	X	27	VAL

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Mol	Chain	Res	Type
1	X	42	ASP
1	X	53	LEU
1	X	61	ASP
1	X	65	GLU
1	X	69	GLN
1	X	99	ARG
1	X	107	ILE
1	X	112	LEU
1	X	113	PHE
1	X	120	LEU
1	X	124	LYS
1	X	131	VAL
1	X	144	LEU
1	X	159	GLN
1	X	183	GLU
1	X	196	MET
1	X	208	GLU
1	X	218	SER
1	X	228	CYS
1	X	232	VAL
1	X	234	GLN
1	X	251	ILE
1	X	276	ASN
1	X	284	VAL
1	X	315	GLN
1	X	325	ASP
1	X	337	ASP
1	X	347	THR
1	X	352	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	X	69	GLN
1	X	222	GLN
1	X	234	GLN
1	X	276	ASN
1	X	315	GLN
1	X	321	GLN
1	X	324	HIS

5.3.3 RNA ⓘ

There are no RNA molecules 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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 2 ligands modelled in this entry, 1 is modelled with single atom - leaving 1 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	NAI	X	401	-	47,48,48	2.08	10 (21%)	64,73,73	1.67	11 (17%)

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	NAI	X	401	-	-	8/29/72/72	0/5/5/5

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	X	401	NAI	O7N-C7N	7.96	1.43	1.24
2	X	401	NAI	PA-O3	4.96	1.64	1.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	X	401	NAI	C4N-C3N	-4.73	1.41	1.50
2	X	401	NAI	C8A-N7A	3.93	1.39	1.31
2	X	401	NAI	C2A-N3A	3.24	1.39	1.33
2	X	401	NAI	C6N-C5N	3.20	1.42	1.33
2	X	401	NAI	C4N-C5N	-3.10	1.41	1.49
2	X	401	NAI	C2A-N1A	3.00	1.39	1.33
2	X	401	NAI	PN-O3	2.39	1.62	1.59
2	X	401	NAI	C2N-C3N	2.27	1.41	1.35

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	X	401	NAI	N3A-C2A-N1A	-6.36	118.95	128.58
2	X	401	NAI	C5A-C4A-N3A	-4.56	120.43	126.72
2	X	401	NAI	C2A-N3A-C4A	3.81	121.14	111.83
2	X	401	NAI	N9A-C8A-N7A	-3.14	109.49	113.94
2	X	401	NAI	O3D-C3D-C4D	-2.74	103.21	111.08
2	X	401	NAI	C5A-N7A-C8A	2.57	107.50	103.45
2	X	401	NAI	O4B-C1B-N9A	-2.54	103.22	108.09
2	X	401	NAI	C5A-C4A-N9A	2.38	108.41	105.81
2	X	401	NAI	N3A-C4A-N9A	2.34	131.14	127.17
2	X	401	NAI	O4D-C1D-N1N	2.27	112.42	108.08
2	X	401	NAI	C4A-C5A-N7A	-2.21	108.06	110.58

There are no chirality outliers.

All (8) torsion outliers are listed below:

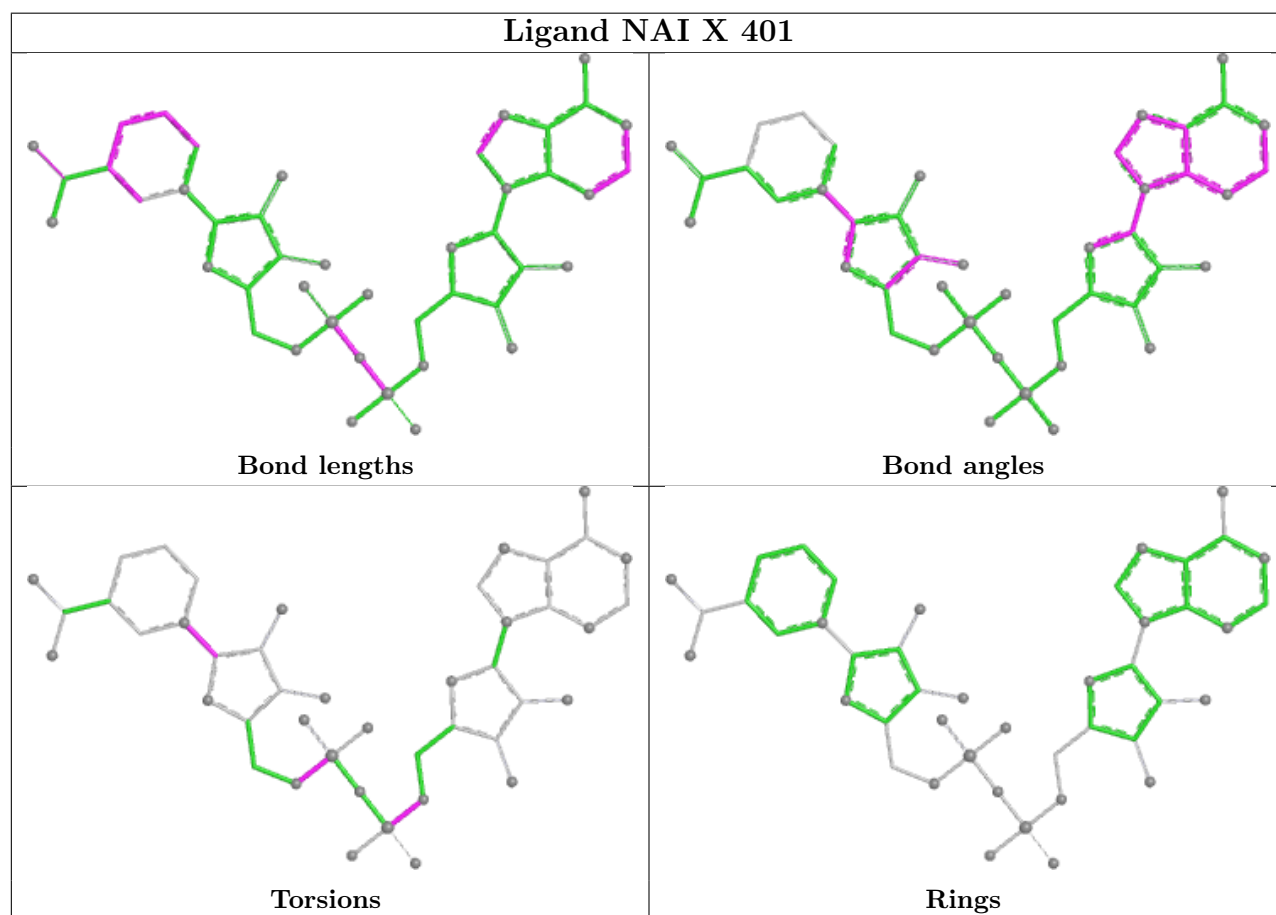
Mol	Chain	Res	Type	Atoms
2	X	401	NAI	C5D-O5D-PN-O3
2	X	401	NAI	C5D-O5D-PN-O1N
2	X	401	NAI	C5D-O5D-PN-O2N
2	X	401	NAI	C2D-C1D-N1N-C2N
2	X	401	NAI	C2D-C1D-N1N-C6N
2	X	401	NAI	O4D-C1D-N1N-C2N
2	X	401	NAI	C5B-O5B-PA-O1A
2	X	401	NAI	O4D-C1D-N1N-C6N

There are no ring outliers.

1 monomer is involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	X	401	NAI	7	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 ⓘ

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	X	359/364 (98%)	0.72	43 (11%) 9 8	21, 55, 86, 97	5 (1%)

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	X	5	SER	5.0
1	X	359	ASP	4.7
1	X	112	LEU	4.6
1	X	258	ALA	4.2
1	X	326	ASP	4.1
1	X	121	ALA	4.0
1	X	30	ALA	3.4
1	X	111	ARG	3.2
1	X	364	LEU	3.1
1	X	32	ALA	3.0
1	X	119	ALA	3.0
1	X	118	CYS	2.9
1	X	10	ALA	2.9
1	X	125	VAL	2.8
1	X	260	ALA	2.8
1	X	128	ILE	2.7
1	X	217	VAL	2.7
1	X	330	ALA	2.7
1	X	120	LEU	2.5
1	X	233	LEU	2.5
1	X	122	ASN	2.5
1	X	239	ASP	2.4
1	X	27	VAL	2.4
1	X	322	ARG	2.4
1	X	116	VAL	2.4
1	X	46	TRP	2.4
1	X	8	ILE	2.3

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Mol	Chain	Res	Type	RSRZ
1	X	13	GLY	2.3
1	X	362	ALA	2.3
1	X	340	VAL	2.3
1	X	114	PRO	2.3
1	X	257	PRO	2.2
1	X	358	SER	2.2
1	X	328	LEU	2.2
1	X	333	ARG	2.2
1	X	363	ARG	2.2
1	X	353	VAL	2.2
1	X	318	GLU	2.2
1	X	342	PRO	2.0
1	X	310	LEU	2.0
1	X	237	ARG	2.0
1	X	36	GLY	2.0
1	X	124	LYS	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

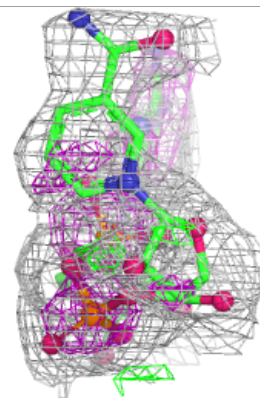
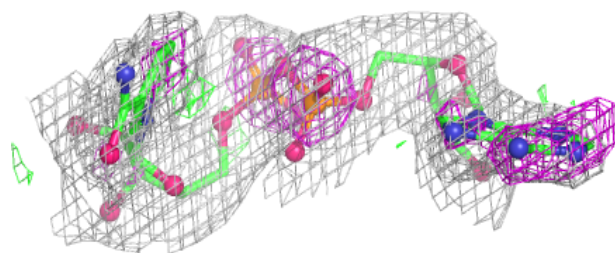
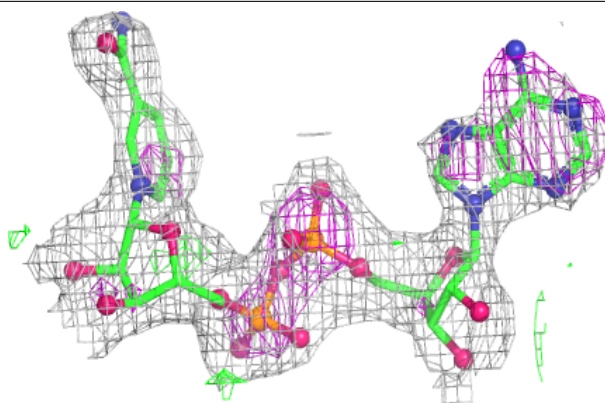
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	NH4	X	402	1/1	0.93	0.19	22,22,22,22	0
2	NAI	X	401	44/44	0.95	0.10	12,14,16,16	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 NAI X 401:

$2mF_o-DF_c$ (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.