



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 09:48 AM UTC

PDB ID : 8ZMX / pdb_00008zmx
Title : Crystal Structures of meso-diaminopimelate dehydrogenase from *Bacillus thermozeamaize*
Authors : Geng, Q.; Wei, Y.; Zhang, Z.J.
Deposited on : 2024-05-24
Resolution : 2.46 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

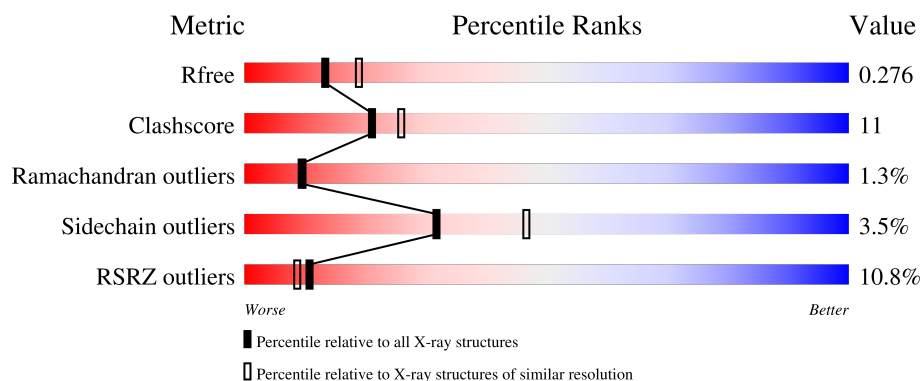
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.46 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	1190 (2.46-2.46)
Clashscore	190562	1229 (2.46-2.46)
Ramachandran outliers	187476	1218 (2.46-2.46)
Sidechain outliers	187428	1218 (2.46-2.46)
RSRZ outliers	180081	1190 (2.46-2.46)

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	313	11% 69% 22% 6%
1	B	313	10% 67% 22% 9%
1	C	313	9% 62% 25% 11%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6776 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 Meso-diaminopimelate D-dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	293	Total	C	N	O	S	0	0	0
			2265	1423	405	424	13			
1	B	286	Total	C	N	O	S	0	0	0
			2201	1377	395	416	13			
1	C	279	Total	C	N	O	S	0	0	0
			2140	1343	384	400	13			

There are 36 discrepancies between the modelled and reference sequences:

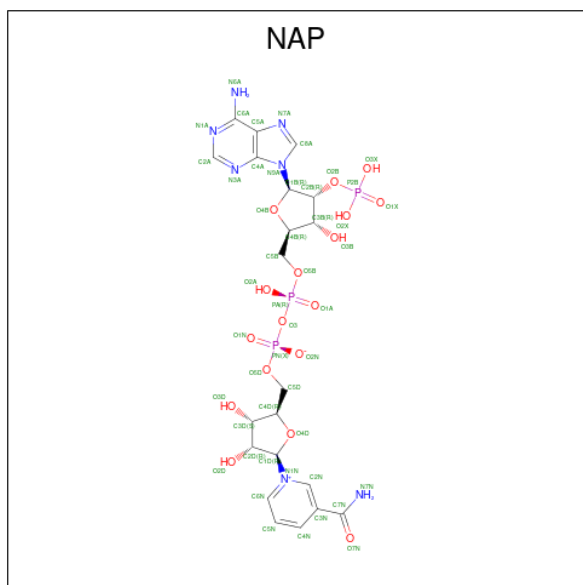
Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	MET	-	initiating methionine	UNP A0A1Y3PXT7
A	-4	HIS	-	expression tag	UNP A0A1Y3PXT7
A	-3	HIS	-	expression tag	UNP A0A1Y3PXT7
A	-2	HIS	-	expression tag	UNP A0A1Y3PXT7
A	-1	HIS	-	expression tag	UNP A0A1Y3PXT7
A	0	HIS	-	expression tag	UNP A0A1Y3PXT7
A	1	HIS	-	expression tag	UNP A0A1Y3PXT7
A	129	THR	TRP	engineered mutation	UNP A0A1Y3PXT7
A	134	CYS	ASP	engineered mutation	UNP A0A1Y3PXT7
A	154	VAL	PHE	engineered mutation	UNP A0A1Y3PXT7
A	177	ALA	SER	engineered mutation	UNP A0A1Y3PXT7
A	235	ILE	HIS	engineered mutation	UNP A0A1Y3PXT7
B	-5	MET	-	initiating methionine	UNP A0A1Y3PXT7
B	-4	HIS	-	expression tag	UNP A0A1Y3PXT7
B	-3	HIS	-	expression tag	UNP A0A1Y3PXT7
B	-2	HIS	-	expression tag	UNP A0A1Y3PXT7
B	-1	HIS	-	expression tag	UNP A0A1Y3PXT7
B	0	HIS	-	expression tag	UNP A0A1Y3PXT7
B	1	HIS	-	expression tag	UNP A0A1Y3PXT7
B	129	THR	TRP	engineered mutation	UNP A0A1Y3PXT7
B	134	CYS	ASP	engineered mutation	UNP A0A1Y3PXT7
B	154	VAL	PHE	engineered mutation	UNP A0A1Y3PXT7
B	177	ALA	SER	engineered mutation	UNP A0A1Y3PXT7

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Chain	Residue	Modelled	Actual	Comment	Reference
B	235	ILE	HIS	engineered mutation	UNP A0A1Y3PXT7
C	-5	MET	-	initiating methionine	UNP A0A1Y3PXT7
C	-4	HIS	-	expression tag	UNP A0A1Y3PXT7
C	-3	HIS	-	expression tag	UNP A0A1Y3PXT7
C	-2	HIS	-	expression tag	UNP A0A1Y3PXT7
C	-1	HIS	-	expression tag	UNP A0A1Y3PXT7
C	0	HIS	-	expression tag	UNP A0A1Y3PXT7
C	1	HIS	-	expression tag	UNP A0A1Y3PXT7
C	129	THR	TRP	engineered mutation	UNP A0A1Y3PXT7
C	134	CYS	ASP	engineered mutation	UNP A0A1Y3PXT7
C	154	VAL	PHE	engineered mutation	UNP A0A1Y3PXT7
C	177	ALA	SER	engineered mutation	UNP A0A1Y3PXT7
C	235	ILE	HIS	engineered mutation	UNP A0A1Y3PXT7

- Molecule 2 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (CCD ID: NAP) (formula: $C_{21}H_{28}N_7O_{17}P_3$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 48	C 21	N 7	O 17	P 3	0	0
2	B	1	Total 48	C 21	N 7	O 17	P 3	0	0
2	C	1	Total 48	C 21	N 7	O 17	P 3	0	0

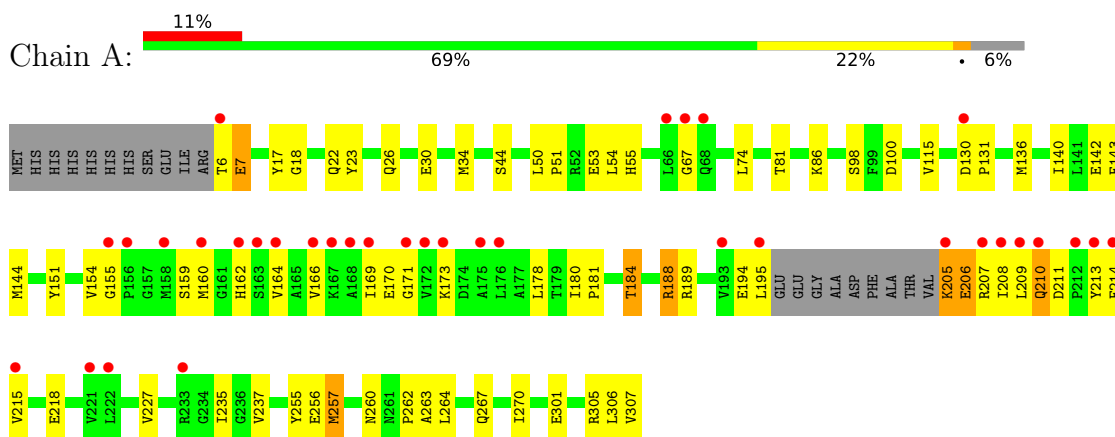
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	8	Total 8	O 8	0	0
3	B	7	Total 7	O 7	0	0
3	C	11	Total 11	O 11	0	0

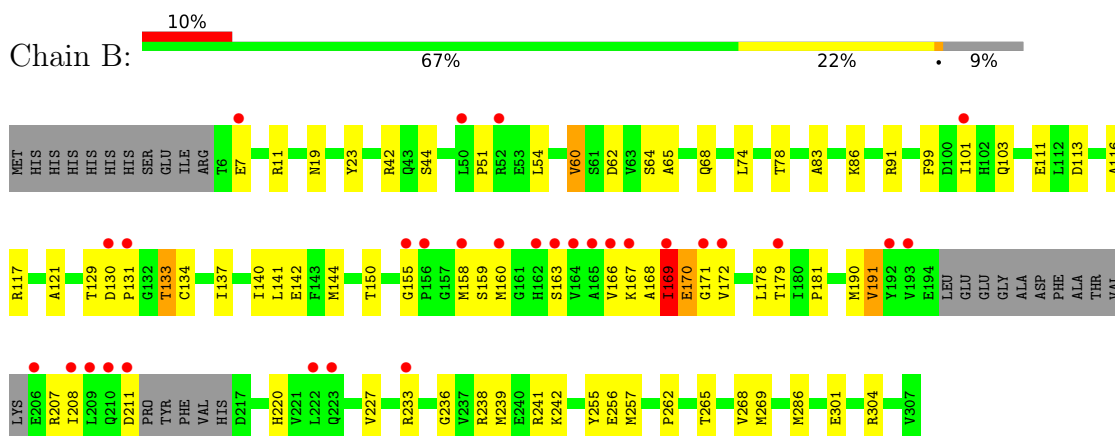
3 Residue-property plots [i](#)

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.

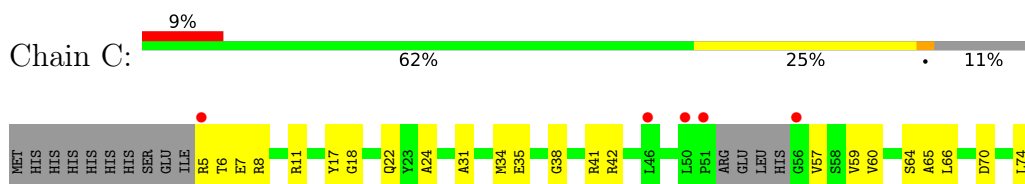
- Molecule 1: Meso-diaminopimelate D-dehydrogenase

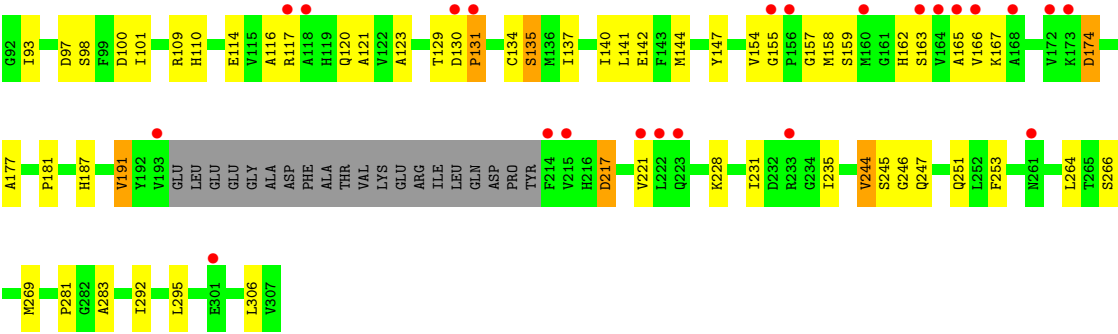


- Molecule 1: Meso-diaminopimelate D-dehydrogenase



- Molecule 1: Meso-diaminopimelate D-dehydrogenase





4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	157.09Å 157.09Å 111.45Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.81 – 2.46 49.81 – 2.46	Depositor EDS
% Data completeness (in resolution range)	99.9 (49.81-2.46) 99.9 (49.81-2.46)	Depositor EDS
R_{merge}	0.22	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.11 (at 2.45Å)	Xtriage
Refinement program	PHENIX 1.19.2_4158	Depositor
R, R_{free}	0.225 , 0.275 0.233 , 0.276	Depositor DCC
R_{free} test set	2560 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	64.6	Xtriage
Anisotropy	0.262	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 40.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6776	wwPDB-VP
Average B, all atoms (Å ²)	70.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.81% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: NAP

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	0.42	0/2300	0.62	0/3115
1	B	0.41	0/2231	0.67	2/3019 (0.1%)
1	C	0.40	0/2171	0.62	3/2939 (0.1%)
All	All	0.41	0/6702	0.64	5/9073 (0.1%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	129	THR	CA-C-N	-5.67	107.97	121.80
1	B	129	THR	C-N-CA	-5.67	107.97	121.80
1	C	131	PRO	N-CA-C	-5.64	100.85	112.47
1	C	129	THR	CA-C-N	-5.57	108.22	121.80
1	C	129	THR	C-N-CA	-5.57	108.22	121.80

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	2265	0	2307	54	0
1	B	2201	0	2241	46	0
1	C	2140	0	2183	61	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	48	0	25	1	0
2	B	48	0	25	3	0
2	C	48	0	25	3	0
3	A	8	0	0	0	0
3	B	7	0	0	0	0
3	C	11	0	0	0	0
All	All	6776	0	6806	149	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 11.

All (149) 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:C:155:GLY:HA2	1:C:181:PRO:HG3	1.54	0.90
1:C:147:TYR:HB2	1:C:244:VAL:HG13	1.56	0.84
1:A:306:LEU:HD21	1:C:142:GLU:HG3	1.60	0.83
1:A:142:GLU:HG3	1:C:306:LEU:HD21	1.65	0.77
1:C:140:ILE:HG22	1:C:144:MET:HE2	1.66	0.76
1:C:38:GLY:HA3	1:C:66:LEU:HD22	1.70	0.73
1:B:166:VAL:HA	1:B:169:ILE:HD11	1.71	0.71
1:B:155:GLY:HA2	1:B:181:PRO:HG3	1.73	0.70
1:A:255:TYR:OH	1:A:257:MET:HE3	1.91	0.70
1:B:103:GLN:CD	1:B:103:GLN:H	2.00	0.69
1:A:166:VAL:HA	1:A:169:ILE:CD1	2.24	0.68
1:C:130:ASP:HA	1:C:134:CYS:HB2	1.76	0.67
1:C:91:ARG:HB3	1:C:93:ILE:HD12	1.77	0.66
1:C:167:LYS:NZ	1:C:174:ASP:OD1	2.28	0.66
1:B:140:ILE:HG22	1:B:144:MET:HE2	1.78	0.66
1:C:228:LYS:HA	1:C:231:ILE:HD12	1.78	0.65
1:A:173:LYS:CD	1:A:195:LEU:HA	2.30	0.62
1:B:301:GLU:OE1	1:B:304:ARG:NH1	2.31	0.62
1:A:155:GLY:HA2	1:A:181:PRO:HG3	1.83	0.61
1:A:173:LYS:HD3	1:A:195:LEU:HA	1.83	0.60
1:B:255:TYR:OH	1:B:257:MET:HE3	2.02	0.60
1:A:307:VAL:HA	1:C:135:SER:HB3	1.82	0.60
1:C:11:ARG:HD2	1:C:35:GLU:OE2	2.02	0.60
1:B:86:LYS:NZ	1:B:111:GLU:OE1	2.27	0.59
1:C:116:ALA:HB1	1:C:121:ALA:O	2.02	0.59
1:A:209:LEU:HD12	1:A:210:GLN:N	2.19	0.58
1:C:79:ARG:NH1	1:C:101:ILE:HD11	2.18	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:166:VAL:HA	1:A:169:ILE:HD12	1.83	0.58
1:A:207:ARG:NE	1:A:207:ARG:HA	2.18	0.58
1:A:211:ASP:HB2	1:A:215:VAL:HB	1.86	0.58
1:B:170:GLU:O	1:B:172:VAL:N	2.29	0.58
1:B:78:THR:HG21	1:B:101:ILE:HG12	1.86	0.57
1:C:191:VAL:HG22	1:C:221:VAL:HG22	1.86	0.57
1:C:159:SER:HB3	1:C:162:HIS:HB2	1.88	0.55
1:A:30:GLU:HG2	1:C:247:GLN:HG3	1.88	0.55
1:A:18:GLY:O	1:A:22:GLN:HG3	2.07	0.54
1:A:162:HIS:HB3	1:A:189:ARG:NH1	2.23	0.54
1:B:233:ARG:HG3	1:B:262:PRO:CD	2.37	0.53
1:C:18:GLY:O	1:C:22:GLN:HG3	2.08	0.53
1:A:171:GLY:O	1:A:195:LEU:HD21	2.08	0.53
1:B:268:VAL:HG13	1:B:286:MET:HE1	1.91	0.53
1:A:142:GLU:HG3	1:C:306:LEU:CD2	2.38	0.53
1:B:130:ASP:HB3	1:B:131:PRO:HD3	1.90	0.53
1:A:143:PHE:HB3	1:C:295:LEU:HD23	1.90	0.53
1:A:213:TYR:C	1:A:215:VAL:H	2.17	0.53
1:B:265:THR:O	1:B:269:MET:HG3	2.10	0.52
1:C:42:ARG:NH2	2:C:401:NAP:H2B	2.24	0.52
1:C:31:ALA:HB3	1:C:34:MET:HB2	1.90	0.52
1:B:78:THR:N	2:B:401:NAP:O3D	2.43	0.51
1:C:154:VAL:HG12	1:C:187:HIS:NE2	2.24	0.51
1:C:98:SER:O	2:C:401:NAP:H2N	2.11	0.51
1:B:170:GLU:C	1:B:172:VAL:H	2.16	0.51
1:C:140:ILE:O	1:C:144:MET:HG3	2.10	0.51
1:A:50:LEU:HD23	1:A:55:HIS:HD2	1.76	0.50
1:A:159:SER:OG	1:A:162:HIS:ND1	2.43	0.50
1:A:237:VAL:O	1:A:256:GLU:HA	2.11	0.50
1:B:60:VAL:HG21	1:B:65:ALA:HB3	1.93	0.50
1:C:74:LEU:O	1:C:269:MET:HE1	2.12	0.50
1:C:158:MET:HE1	1:C:231:ILE:HG23	1.93	0.50
1:A:209:LEU:HD12	1:A:210:GLN:H	1.77	0.49
1:C:79:ARG:NH1	1:C:101:ILE:CD1	2.75	0.49
1:B:238:ARG:HG3	1:B:256:GLU:HG2	1.94	0.49
1:A:178:LEU:HB2	1:A:180:ILE:HD11	1.94	0.49
1:A:207:ARG:C	1:A:209:LEU:H	2.21	0.48
1:B:42:ARG:NH2	2:B:401:NAP:H2B	2.27	0.48
1:A:205:LYS:O	1:A:207:ARG:N	2.46	0.48
1:B:62:ASP:OD1	1:B:64:SER:OG	2.24	0.48
1:B:238:ARG:HG3	1:B:256:GLU:CG	2.44	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:86:LYS:HG2	1:A:115:VAL:HG21	1.94	0.48
1:B:113:ASP:O	1:B:117:ARG:HG3	2.15	0.47
1:A:260:ASN:HB3	1:A:263:ALA:HB3	1.97	0.47
1:B:239:MET:HB3	1:B:255:TYR:HB3	1.96	0.47
1:B:11:ARG:NH1	1:B:68:GLN:O	2.47	0.47
1:A:169:ILE:HG22	1:A:170:GLU:H	1.79	0.46
1:C:8:ARG:HH22	1:C:70:ASP:CG	2.22	0.46
1:B:191:VAL:HG23	1:B:220:HIS:O	2.15	0.46
1:A:301:GLU:OE2	1:A:305:ARG:NH2	2.44	0.46
1:B:170:GLU:OE1	1:B:208:ILE:HD13	2.15	0.46
1:C:245:SER:OG	1:C:251:GLN:HG3	2.16	0.46
1:B:23:TYR:CD2	1:B:262:PRO:HB2	2.51	0.46
1:A:34:MET:HE1	1:A:270:ILE:O	2.16	0.45
1:A:98:SER:O	2:A:401:NAP:H2N	2.15	0.45
1:A:264:LEU:HD11	1:C:144:MET:HB3	1.98	0.45
1:C:97:ASP:OD1	1:C:97:ASP:C	2.58	0.45
1:A:50:LEU:HD23	1:A:55:HIS:CD2	2.51	0.45
1:C:57:VAL:O	1:C:59:VAL:HG23	2.17	0.45
1:A:23:TYR:CD2	1:A:262:PRO:HB2	2.52	0.45
1:C:60:VAL:HG21	1:C:65:ALA:HB3	1.99	0.45
1:B:64:SER:HB3	1:B:91:ARG:HH12	1.82	0.45
1:B:208:ILE:HD12	1:B:208:ILE:H	1.82	0.45
1:B:51:PRO:HD2	1:B:54:LEU:HD12	1.99	0.44
1:C:217:ASP:N	1:C:217:ASP:OD1	2.50	0.44
1:A:151:TYR:CE1	1:A:184:THR:HG22	2.52	0.44
1:A:206:GLU:HG2	1:A:207:ARG:HE	1.83	0.44
1:C:157:GLY:HA3	1:C:235:ILE:HG13	1.98	0.44
1:A:26:GLN:HB3	1:C:247:GLN:HB2	2.00	0.44
1:C:24:ALA:HA	1:C:266:SER:HB2	1.98	0.44
1:A:17:TYR:OH	1:A:53:GLU:HG3	2.17	0.44
1:C:110:HIS:O	1:C:114:GLU:HG3	2.18	0.44
1:B:150:THR:OG1	1:B:241:ARG:HD2	2.18	0.44
1:B:141:LEU:HA	1:B:144:MET:HE3	1.99	0.44
1:C:100:ASP:HA	1:C:131:PRO:HG2	1.99	0.43
1:B:130:ASP:HA	1:B:134:CYS:HB2	1.99	0.43
1:B:83:ALA:HB2	1:C:5:ARG:O	2.18	0.43
1:A:194:GLU:HB2	1:A:227:VAL:HG13	2.01	0.43
1:B:158:MET:HA	1:B:178:LEU:HD22	2.00	0.43
1:C:123:ALA:O	1:C:283:ALA:HA	2.19	0.43
1:A:257:MET:HE2	1:C:253:PHE:CD2	2.54	0.43
1:B:116:ALA:HB1	1:B:121:ALA:O	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:78:THR:N	2:C:401:NAP:O3D	2.52	0.43
1:A:162:HIS:O	1:A:166:VAL:HG23	2.18	0.43
1:A:51:PRO:HD2	1:A:54:LEU:HD12	1.99	0.42
1:B:242:LYS:HB3	1:B:242:LYS:HE3	1.89	0.42
1:A:136:MET:O	1:A:140:ILE:HG13	2.20	0.42
1:A:237:VAL:HB	1:A:257:MET:HG3	2.00	0.42
1:A:74:LEU:HD23	1:A:81:THR:HG23	2.02	0.42
1:B:207:ARG:HH11	1:B:207:ARG:HG2	1.85	0.42
1:C:137:ILE:O	1:C:141:LEU:HG	2.19	0.42
1:C:141:LEU:HA	1:C:144:MET:HE3	2.01	0.42
1:C:162:HIS:O	1:C:165:ALA:HB3	2.20	0.42
1:B:133:THR:O	1:B:137:ILE:HG12	2.20	0.42
1:C:120:GLN:HA	1:C:281:PRO:HG3	2.01	0.42
1:C:158:MET:HE3	1:C:158:MET:HB2	1.72	0.42
1:B:7:GLU:OE2	1:B:7:GLU:HA	2.20	0.42
1:C:228:LYS:HB2	1:C:228:LYS:HE3	1.69	0.42
1:C:159:SER:OG	1:C:162:HIS:ND1	2.51	0.41
1:A:213:TYR:O	1:A:214:PHE:HB2	2.20	0.41
1:C:6:THR:OG1	1:C:7:GLU:N	2.53	0.41
1:C:163:SER:HB3	1:C:177:ALA:H	1.85	0.41
1:A:144:MET:HG2	1:C:264:LEU:HD11	2.02	0.41
1:A:130:ASP:HB3	1:A:131:PRO:HD3	2.03	0.41
1:B:19:ASN:HB3	2:B:401:NAP:O1N	2.20	0.41
1:C:79:ARG:HA	1:C:79:ARG:HD2	1.75	0.41
1:A:160:MET:O	1:A:164:VAL:HG22	2.21	0.41
1:A:100:ASP:HA	1:A:131:PRO:HG2	2.02	0.41
1:B:169:ILE:HB	1:B:170:GLU:H	1.52	0.41
1:A:6:THR:O	1:A:7:GLU:HB2	2.21	0.41
1:B:170:GLU:C	1:B:172:VAL:N	2.78	0.41
1:B:236:GLY:HA2	1:B:257:MET:O	2.20	0.41
1:C:41:ARG:HE	1:C:41:ARG:HB3	1.66	0.41
1:C:162:HIS:O	1:C:166:VAL:HG23	2.21	0.41
1:C:292:ILE:HA	1:C:295:LEU:HD12	2.03	0.41
1:A:267:GLN:OE1	1:C:246:GLY:N	2.43	0.40
1:B:74:LEU:O	1:B:269:MET:HE1	2.22	0.40
1:B:159:SER:OG	1:B:160:MET:N	2.54	0.40
1:B:190:MET:HE3	1:B:190:MET:HB3	1.83	0.40
1:C:17:TYR:CE2	1:C:22:GLN:HG2	2.56	0.40
1:A:188:ARG:HD3	1:A:218:GLU:OE2	2.22	0.40
1:C:8:ARG:NH2	1:C:70:ASP:OD2	2.55	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	289/313 (92%)	267 (92%)	17 (6%)	5 (2%)	7	6
1	B	280/313 (90%)	256 (91%)	19 (7%)	5 (2%)	6	5
1	C	273/313 (87%)	255 (93%)	17 (6%)	1 (0%)	30	37
All	All	842/939 (90%)	778 (92%)	53 (6%)	11 (1%)	9	9

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	7	GLU
1	A	67	GLY
1	A	206	GLU
1	A	210	GLN
1	B	168	ALA
1	B	169	ILE
1	C	174	ASP
1	A	208	ILE
1	B	170	GLU
1	B	171	GLY
1	B	99	PHE

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	244/261 (94%)	237 (97%)	7 (3%)	37	52

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	237/261 (91%)	226 (95%)	11 (5%)	24	36
1	C	230/261 (88%)	223 (97%)	7 (3%)	36	51
All	All	711/783 (91%)	686 (96%)	25 (4%)	32	46

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	44	SER
1	A	154	VAL
1	A	184	THR
1	A	188	ARG
1	A	205	LYS
1	A	235	ILE
1	A	257	MET
1	B	44	SER
1	B	60	VAL
1	B	133	THR
1	B	142	GLU
1	B	163	SER
1	B	167	LYS
1	B	169	ILE
1	B	179	THR
1	B	191	VAL
1	B	211	ASP
1	B	227	VAL
1	C	64	SER
1	C	109	ARG
1	C	117	ARG
1	C	135	SER
1	C	191	VAL
1	C	217	ASP
1	C	244	VAL

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

Mol	Chain	Res	Type
1	A	43	GLN
1	A	55	HIS
1	A	68	GLN
1	A	216	HIS

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Mol	Chain	Res	Type
1	B	110	HIS
1	C	249	GLN
1	C	261	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](#)

3 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	NAP	C	401	-	50,52,52	0.64	1 (2%)	71,80,80	0.93	3 (4%)
2	NAP	B	401	-	50,52,52	0.59	1 (2%)	71,80,80	0.93	2 (2%)
2	NAP	A	401	-	50,52,52	0.63	1 (2%)	71,80,80	0.86	4 (5%)

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	NAP	C	401	-	-	0/35/67/67	0/5/5/5
2	NAP	B	401	-	-	6/35/67/67	0/5/5/5
2	NAP	A	401	-	-	2/35/67/67	0/5/5/5

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	NAP	C2N-N1N	2.88	1.38	1.35
2	B	401	NAP	C2N-N1N	2.71	1.38	1.35
2	C	401	NAP	C2N-N1N	2.42	1.37	1.35

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	401	NAP	P2B-O2B-C2B	-5.07	109.89	123.43
2	C	401	NAP	P2B-O2B-C2B	-4.69	110.89	123.43
2	A	401	NAP	P2B-O2B-C2B	-3.64	113.71	123.43
2	B	401	NAP	C6N-N1N-C2N	-3.06	119.27	121.88
2	C	401	NAP	C6N-N1N-C2N	-2.71	119.57	121.88
2	A	401	NAP	C6N-N1N-C2N	-2.61	119.66	121.88
2	A	401	NAP	C4D-O4D-C1D	-2.24	107.87	109.92
2	C	401	NAP	O2B-C2B-C3B	-2.12	104.06	111.68
2	A	401	NAP	O2B-C2B-C3B	-2.07	104.27	111.68

There are no chirality outliers.

All (8) torsion outliers are listed below:

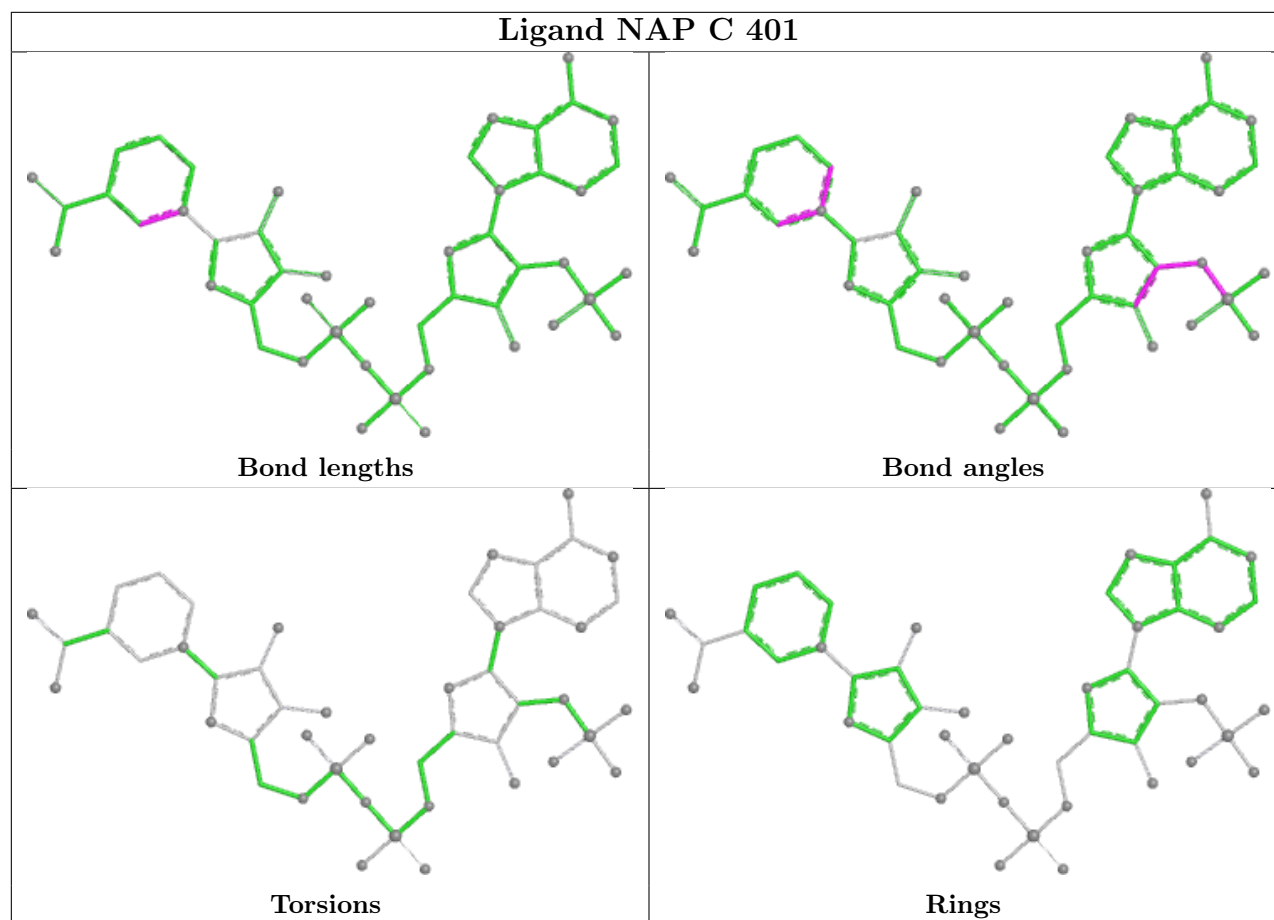
Mol	Chain	Res	Type	Atoms
2	B	401	NAP	C5D-O5D-PN-O3
2	B	401	NAP	C5D-O5D-PN-O1N
2	B	401	NAP	C3D-C4D-C5D-O5D
2	B	401	NAP	O4D-C4D-C5D-O5D
2	B	401	NAP	C5D-O5D-PN-O2N
2	B	401	NAP	O4B-C4B-C5B-O5B
2	A	401	NAP	O4B-C4B-C5B-O5B
2	A	401	NAP	O4B-C1B-N9A-C8A

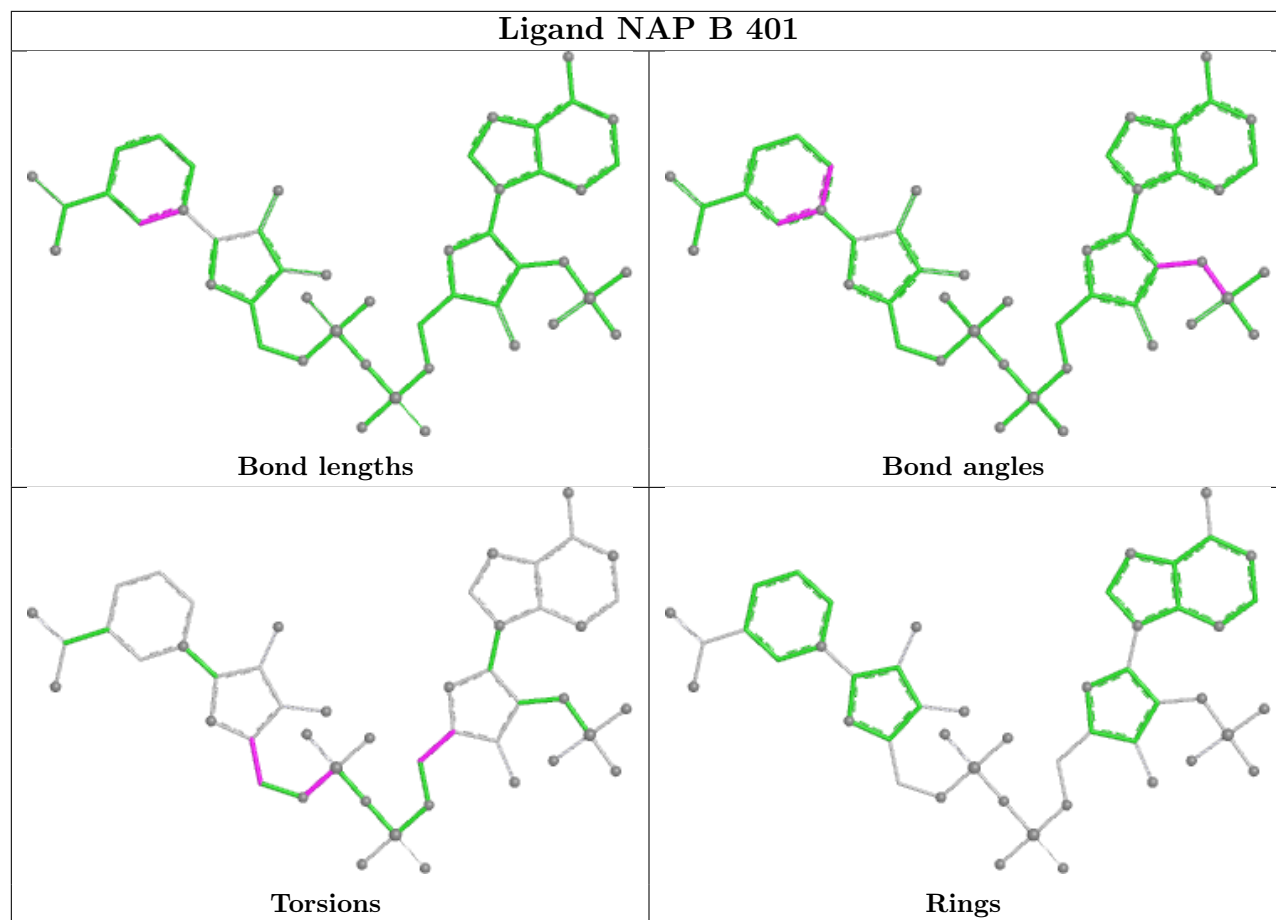
There are no ring outliers.

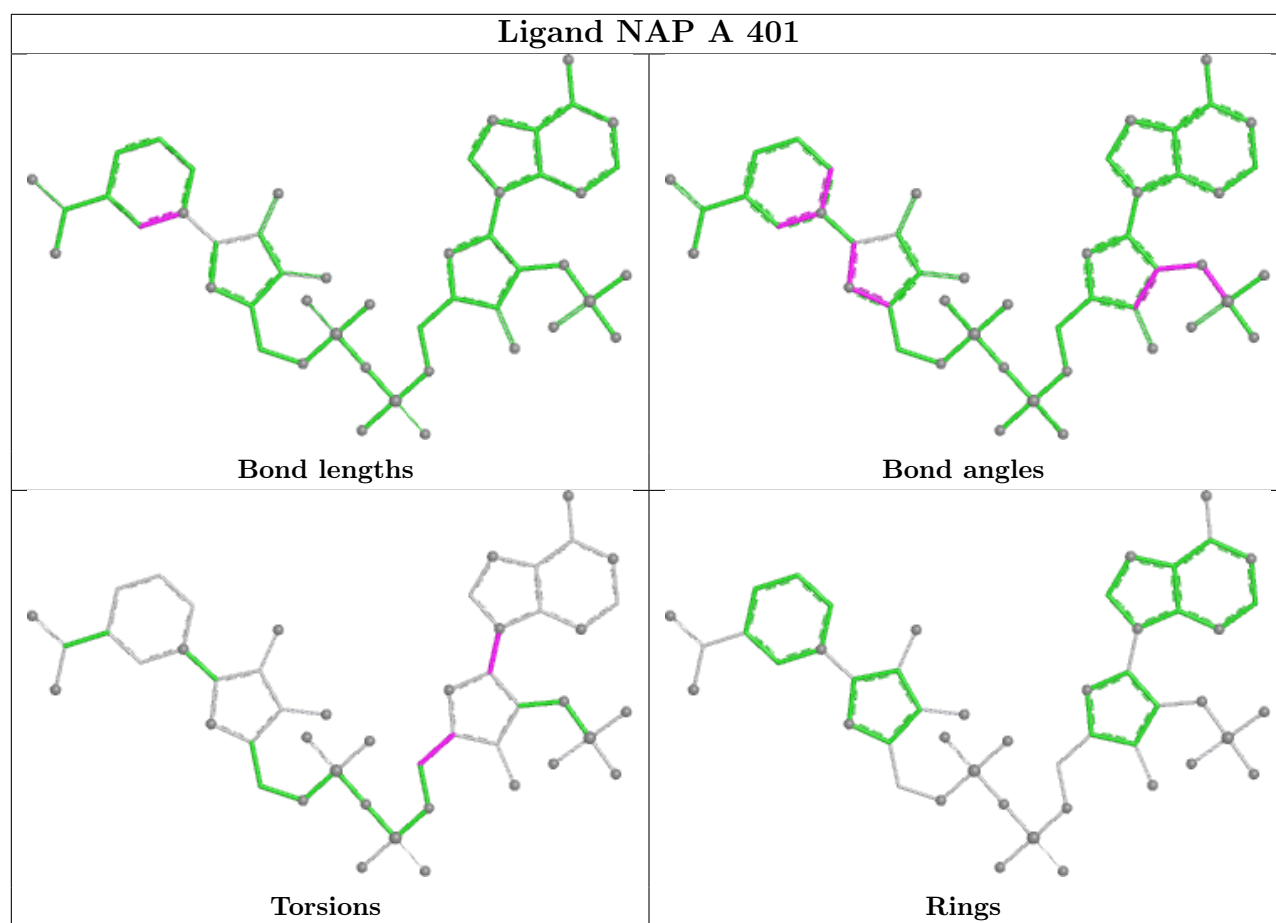
3 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	401	NAP	3	0
2	B	401	NAP	3	0
2	A	401	NAP	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	293/313 (93%)	0.60	35 (11%) 9 7	49, 62, 111, 130	0
1	B	286/313 (91%)	0.57	30 (10%) 11 9	30, 65, 112, 128	0
1	C	279/313 (89%)	0.72	28 (10%) 12 10	30, 68, 111, 134	0
All	All	858/939 (91%)	0.63	93 (10%) 11 9	30, 65, 112, 134	0

All (93) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	214	PHE	7.5
1	C	172	VAL	6.6
1	A	209	LEU	6.5
1	C	117	ARG	5.2
1	C	155	GLY	5.1
1	A	208	ILE	4.8
1	A	213	TYR	4.7
1	B	130	ASP	4.6
1	B	52	ARG	4.6
1	A	171	GLY	4.4
1	A	166	VAL	4.4
1	C	164	VAL	4.3
1	B	208	ILE	4.3
1	B	209	LEU	4.2
1	C	193	VAL	4.2
1	C	215	VAL	4.1
1	C	130	ASP	4.0
1	B	211	ASP	4.0
1	B	164	VAL	4.0
1	C	56	GLY	3.8
1	B	233	ARG	3.8
1	B	160	MET	3.8
1	C	51	PRO	3.7

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Mol	Chain	Res	Type	RSRZ
1	B	172	VAL	3.6
1	C	166	VAL	3.6
1	A	130	ASP	3.6
1	A	162	HIS	3.5
1	A	205	LYS	3.5
1	B	166	VAL	3.5
1	A	164	VAL	3.5
1	A	155	GLY	3.5
1	B	171	GLY	3.5
1	A	172	VAL	3.4
1	A	210	GLN	3.4
1	B	167	LYS	3.3
1	B	193	VAL	3.3
1	B	131	PRO	3.3
1	A	158	MET	3.2
1	C	160	MET	3.2
1	B	158	MET	3.0
1	A	195	LEU	3.0
1	C	173	LYS	3.0
1	B	162	HIS	2.9
1	B	7	GLU	2.9
1	A	163	SER	2.9
1	B	210	GLN	2.9
1	B	206	GLU	2.9
1	C	156	PRO	2.9
1	B	222	LEU	2.8
1	C	222	LEU	2.8
1	A	173	LYS	2.8
1	A	215	VAL	2.8
1	C	168	ALA	2.8
1	A	67	GLY	2.8
1	C	5	ARG	2.7
1	A	233	ARG	2.6
1	A	212	PRO	2.6
1	A	66	LEU	2.6
1	A	214	PHE	2.5
1	C	221	VAL	2.5
1	B	169	ILE	2.5
1	A	193	VAL	2.5
1	B	101	ILE	2.5
1	B	165	ALA	2.5
1	A	169	ILE	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	68	GLN	2.4
1	B	163	SER	2.4
1	B	192	TYR	2.4
1	C	301	GLU	2.4
1	C	223	GLN	2.4
1	A	160	MET	2.4
1	C	261	ASN	2.4
1	C	165	ALA	2.4
1	C	131	PRO	2.3
1	A	221	VAL	2.3
1	A	6	THR	2.3
1	A	176	LEU	2.3
1	A	175	ALA	2.2
1	B	223	GLN	2.2
1	A	167	LYS	2.2
1	B	50	LEU	2.2
1	A	207	ARG	2.2
1	B	155	GLY	2.2
1	C	46	LEU	2.1
1	C	163	SER	2.1
1	B	156	PRO	2.1
1	C	118	ALA	2.1
1	A	168	ALA	2.0
1	A	222	LEU	2.0
1	C	233	ARG	2.0
1	C	50	LEU	2.0
1	B	179	THR	2.0
1	A	156	PRO	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.

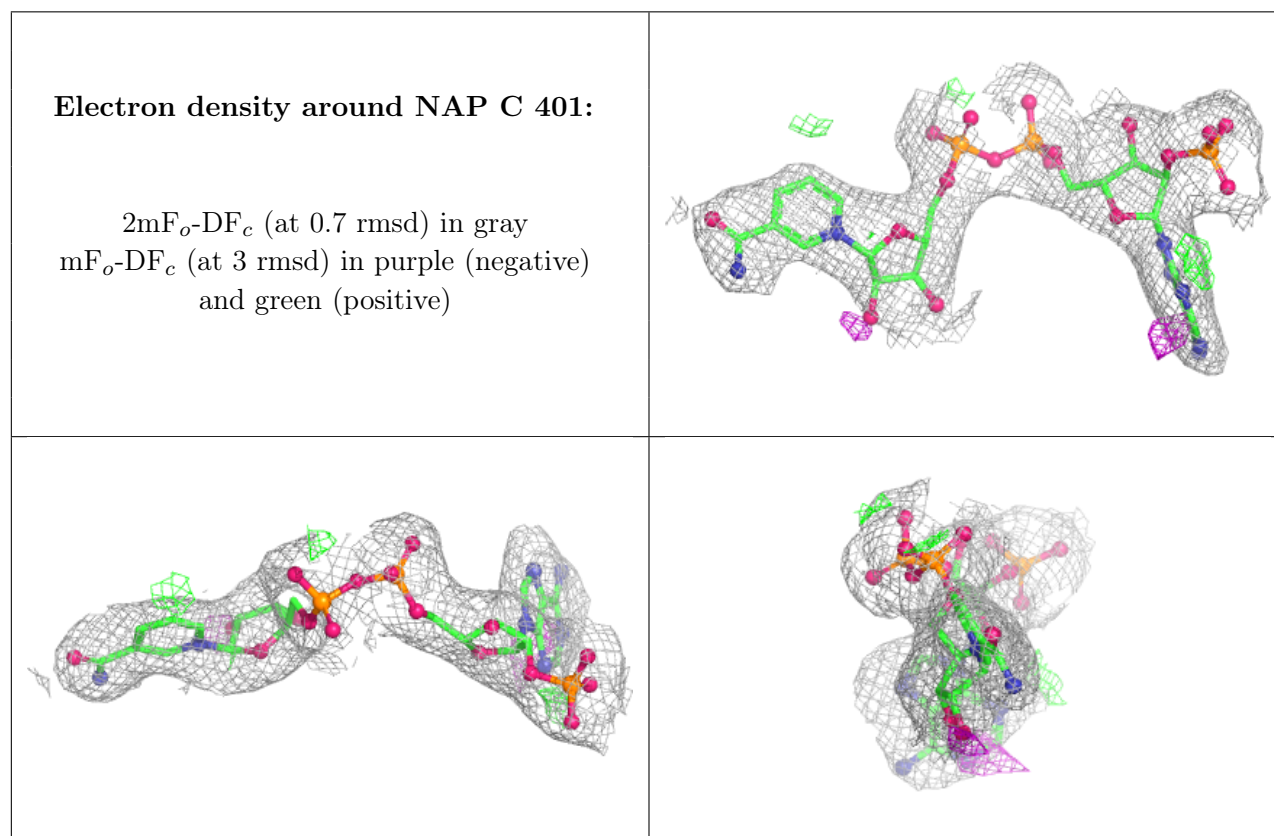
6.4 Ligands [i](#)

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.

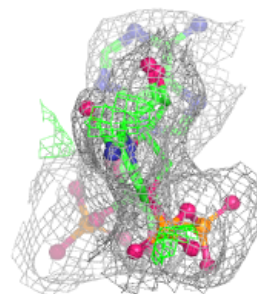
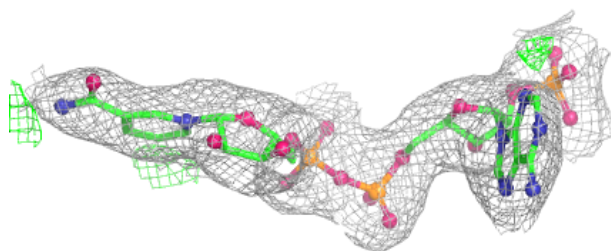
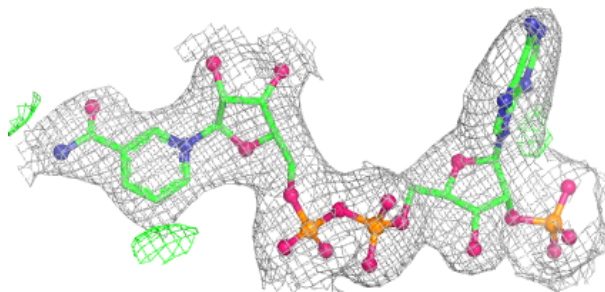
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	NAP	C	401	48/48	0.92	0.10	59,70,79,82	0
2	NAP	B	401	48/48	0.94	0.08	57,66,76,82	0
2	NAP	A	401	48/48	0.95	0.07	57,65,72,74	0

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.

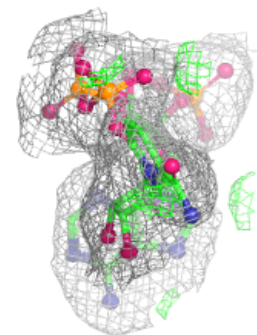
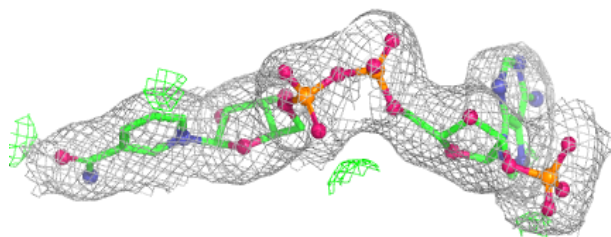
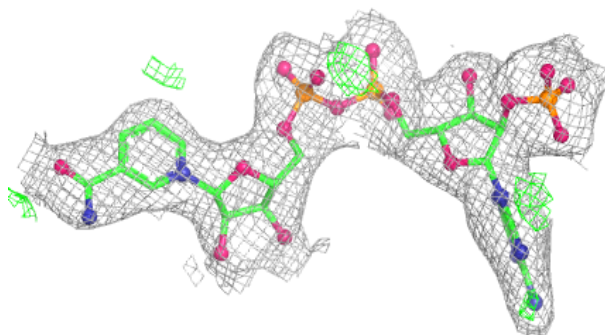


Electron density around NAP B 401:

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

**Electron density around NAP A 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.