



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 21, 2025 – 04:29 pm BST

PDB ID : 9H4F / pdb_00009h4f
Title : Structure of Imine Reductase 361 from Micromonospora sp. mutant M125W /I127F/L179V/H250L
Authors : Ho, E.; Domenech, J.; Crossley, A.; Green, A.P.; Grogan, G.
Deposited on : 2024-10-18
Resolution : 2.77 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 Phenix2.0rc1
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	2.0rc1
EDS	:	3.0
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.006 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.45.1

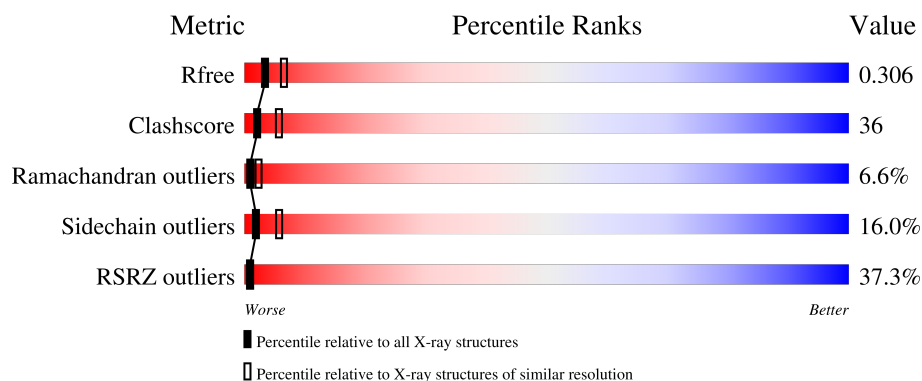
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.77 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	4924 (2.80-2.76)
Clashscore	180529	5458 (2.80-2.76)
Ramachandran outliers	177936	5386 (2.80-2.76)
Sidechain outliers	177891	5388 (2.80-2.76)
RSRZ outliers	164620	4926 (2.80-2.76)

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	297	<div> <div>29%</div> <div>39%</div> <div>51%</div> <div>8%</div> <div>.</div> </div>
1	B	297	<div> <div>41%</div> <div>49%</div> <div>39%</div> <div>8%</div> <div>..</div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3785 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 6-phosphogluconate dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	291	Total	C	N	O	S	0	0	0
			1960	1246	331	376	7			
1	B	288	Total	C	N	O	S	78	0	0
			1777	1101	312	359	5			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	125	TRP	MET	engineered mutation	UNP A0A1S8Y2S4
A	127	PHE	ILE	engineered mutation	UNP A0A1S8Y2S4
A	179	VAL	LEU	engineered mutation	UNP A0A1S8Y2S4
A	250	LEU	HIS	engineered mutation	UNP A0A1S8Y2S4
B	125	TRP	MET	engineered mutation	UNP A0A1S8Y2S4
B	127	PHE	ILE	engineered mutation	UNP A0A1S8Y2S4
B	179	VAL	LEU	engineered mutation	UNP A0A1S8Y2S4
B	250	LEU	HIS	engineered mutation	UNP A0A1S8Y2S4

- Molecule 2 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (CCD ID: NAP) (formula: C₂₁H₂₈N₇O₁₇P₃) (labeled as "Ligand of Interest" by depositor).

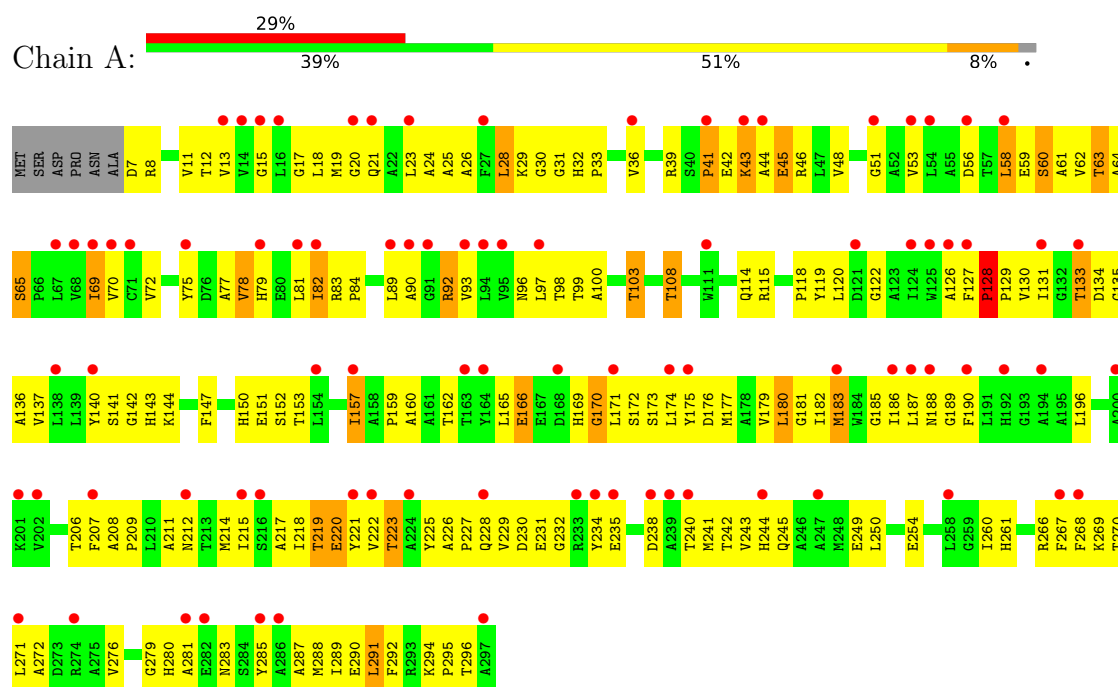


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			48	21	7	17	3		

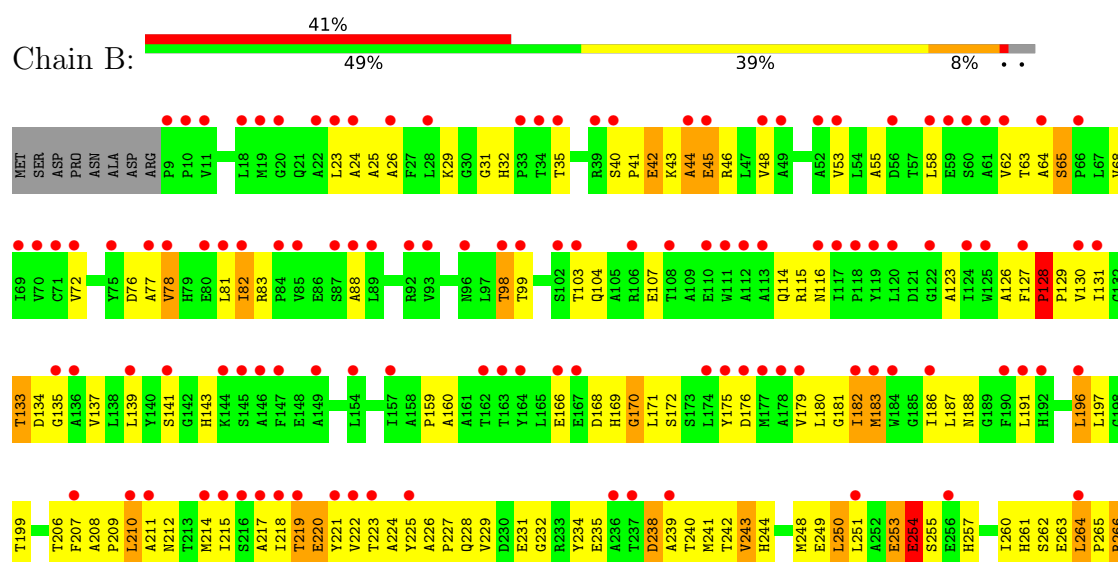
3 Residue-property plots [i](#)

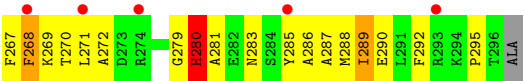
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: 6-phosphogluconate dehydrogenase



• Molecule 1: 6-phosphogluconate dehydrogenase





4 Data and refinement statistics

Property	Value	Source
Space group	I 21 3	Depositor
Cell constants a, b, c, α , β , γ	184.27 Å 184.27 Å 184.27 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	130.30 – 2.77 130.30 – 2.77	Depositor EDS
% Data completeness (in resolution range)	59.6 (130.30-2.77) 59.6 (130.30-2.77)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.18 (at 2.77 Å)	Xtriage
Refinement program	REFMAC 5.8.0430	Depositor
R, R_{free}	0.226 , 0.304 0.233 , 0.306	Depositor DCC
R_{free} test set	708 reflections (2.67%)	wwPDB-VP
Wilson B-factor (Å ²)	127.2	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 246.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.027 for -l,-k,-h	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3785	wwPDB-VP
Average B, all atoms (Å ²)	134.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.76% 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: 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	0.65	0/2003	1.23	5/2767 (0.2%)
1	B	0.75	4/1806 (0.2%)	1.21	8/2496 (0.3%)
All	All	0.70	4/3809 (0.1%)	1.22	13/5263 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	55	ALA	C-N	-10.69	1.20	1.33
1	B	143	HIS	C-N	7.62	1.44	1.33
1	B	68	VAL	C-N	-5.54	1.26	1.33
1	B	114	GLN	C-N	5.52	1.42	1.33

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	55	ALA	CA-C-N	8.48	133.61	120.89
1	B	55	ALA	C-N-CA	8.48	133.61	120.89
1	A	128	PRO	CB-CA-C	6.44	118.78	110.92
1	B	128	PRO	CB-CA-C	6.29	118.60	110.92
1	A	266	ARG	CB-CA-C	6.03	120.46	110.81
1	A	63	THR	CA-CB-OG1	-5.94	100.69	109.60
1	B	76	ASP	CA-CB-CG	5.60	118.20	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	133	THR	CA-CB-OG1	-5.54	101.30	109.60
1	B	266	ARG	CB-CA-C	5.53	119.66	110.81
1	B	253	GLU	CB-CA-C	5.38	119.42	110.81
1	A	147	PHE	CA-CB-CG	-5.16	108.64	113.80
1	B	133	THR	CA-CB-OG1	-5.12	101.92	109.60
1	B	254	GLU	CB-CG-CD	5.12	121.30	112.60

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	115	ARG	Sidechain
1	A	92	ARG	Sidechain

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	1960	0	1750	148	0
1	B	1777	0	1382	124	0
2	A	48	0	25	7	0
All	All	3785	0	3157	249	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 36.

All (249) 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:130:VAL:O	1:A:133:THR:HG22	1.51	1.11
1:B:130:VAL:O	1:B:133:THR:HG22	1.54	1.06
1:A:219:THR:HG22	1:B:215:ILE:CB	2.03	0.88
1:A:296:THR:HG21	1:B:263:GLU:CB	2.12	0.79
1:A:241:MET:HE1	1:A:285:TYR:HA	1.67	0.76
1:B:241:MET:HE1	1:B:285:TYR:HA	1.68	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:19:MET:HE1	1:A:131:ILE:HG13	1.71	0.72
1:A:234:TYR:HB2	1:A:283:ASN:HB3	1.70	0.72
1:A:25:ALA:O	1:A:29:LYS:N	2.24	0.71
1:B:25:ALA:O	1:B:29:LYS:N	2.25	0.70
1:A:190:PHE:HA	1:B:182:ILE:HD11	1.74	0.69
1:A:19:MET:O	1:A:19:MET:HE3	1.92	0.69
1:A:177:MET:HG3	1:B:254:GLU:HG3	1.75	0.69
1:A:212:ASN:HD22	1:B:219:THR:HB	1.58	0.68
1:A:177:MET:HA	1:A:180:LEU:HD12	1.75	0.68
1:B:128:PRO:HB2	1:B:129:PRO:HD3	1.76	0.67
1:B:210:LEU:HD12	1:B:210:LEU:O	1.94	0.67
1:B:77:ALA:O	1:B:81:LEU:HD22	1.94	0.67
1:A:77:ALA:O	1:A:81:LEU:HD22	1.95	0.66
1:A:17:GLY:HA3	2:A:301:NAP:O2N	1.97	0.64
1:A:291:LEU:HD23	1:A:291:LEU:C	2.22	0.64
1:A:93:VAL:HA	1:A:118:PRO:HG2	1.81	0.63
1:A:133:THR:HG23	1:A:135:GLY:H	1.64	0.63
1:A:174:LEU:HG	1:B:197:LEU:HD23	1.80	0.63
1:A:208:ALA:HB3	1:A:209:PRO:HD3	1.81	0.62
1:B:234:TYR:HB2	1:B:283:ASN:HB3	1.80	0.62
1:A:294:LYS:O	1:B:261:HIS:HB2	1.99	0.62
1:A:165:LEU:O	1:A:166:GLU:O	2.18	0.61
1:B:133:THR:HG23	1:B:135:GLY:H	1.65	0.61
1:A:226:ALA:HB3	1:A:227:PRO:CD	2.30	0.61
1:A:179:VAL:C	1:A:180:LEU:O	2.43	0.60
1:B:226:ALA:HB3	1:B:227:PRO:HD3	1.83	0.60
1:A:128:PRO:HB2	1:A:129:PRO:CD	2.31	0.60
1:A:13:VAL:HG22	1:A:36:VAL:HG12	1.84	0.59
1:A:127:PHE:O	1:A:128:PRO:C	2.46	0.59
1:B:264:LEU:C	1:B:264:LEU:HD23	2.27	0.59
1:B:226:ALA:HB3	1:B:227:PRO:CD	2.33	0.59
1:B:127:PHE:O	1:B:128:PRO:C	2.46	0.58
1:A:72:VAL:HB	2:A:301:NAP:H8A	1.85	0.58
1:A:25:ALA:O	1:A:26:ALA:C	2.47	0.58
1:A:120:LEU:HD23	1:A:142:GLY:HA3	1.86	0.57
1:A:108:THR:HG22	1:A:119:TYR:CZ	2.39	0.57
1:A:169:HIS:O	1:A:170:GLY:C	2.47	0.57
1:B:169:HIS:O	1:B:170:GLY:C	2.47	0.57
1:A:179:VAL:O	1:A:180:LEU:C	2.48	0.57
1:A:127:PHE:O	1:A:130:VAL:N	2.37	0.57
1:B:208:ALA:O	1:B:209:PRO:C	2.48	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:250:LEU:O	1:B:254:GLU:N	2.27	0.56
1:B:127:PHE:O	1:B:130:VAL:N	2.38	0.56
1:B:264:LEU:HB3	1:B:265:PRO:CD	2.36	0.56
1:A:234:TYR:CB	1:A:283:ASN:HB3	2.36	0.56
1:A:189:GLY:C	1:B:182:ILE:HG13	2.30	0.56
1:A:179:VAL:O	1:A:182:ILE:HB	2.05	0.55
1:A:296:THR:HB	1:B:262:SER:H	1.71	0.55
1:A:7:ASP:O	1:A:8:ARG:C	2.49	0.55
1:A:208:ALA:O	1:A:209:PRO:C	2.49	0.55
1:B:264:LEU:C	1:B:264:LEU:CD2	2.80	0.55
1:B:269:LYS:O	1:B:270:THR:C	2.51	0.54
1:B:141:SER:OG	1:B:169:HIS:HA	2.07	0.54
1:A:291:LEU:HD23	1:A:292:PHE:N	2.23	0.54
1:A:81:LEU:O	1:A:84:PRO:HD2	2.07	0.54
1:B:25:ALA:O	1:B:26:ALA:C	2.49	0.54
1:A:97:LEU:CD2	1:A:122:GLY:HA3	2.37	0.54
1:A:17:GLY:O	1:A:21:GLN:HG3	2.07	0.54
1:A:41:PRO:O	1:A:43:LYS:N	2.39	0.54
1:A:261:HIS:N	1:B:292:PHE:O	2.36	0.53
1:B:179:VAL:O	1:B:180:LEU:C	2.52	0.53
1:A:267:PHE:O	1:A:271:LEU:HD12	2.09	0.53
1:B:44:ALA:O	1:B:46:ARG:N	2.42	0.53
1:B:225:TYR:O	1:B:228:GLN:N	2.41	0.53
1:A:261:HIS:HA	1:B:295:PRO:HA	1.91	0.53
1:A:223:THR:O	1:A:227:PRO:HD2	2.09	0.53
1:B:267:PHE:O	1:B:268:PHE:C	2.50	0.53
1:B:240:THR:O	1:B:243:VAL:HG23	2.09	0.53
1:A:108:THR:HG22	1:A:119:TYR:CE1	2.45	0.52
1:B:240:THR:O	1:B:244:HIS:HD2	1.93	0.52
1:A:69:ILE:CD1	1:A:69:ILE:N	2.72	0.52
1:A:150:HIS:O	1:A:151:GLU:C	2.53	0.52
1:A:171:LEU:O	1:A:172:SER:C	2.53	0.52
1:A:292:PHE:HA	1:B:261:HIS:HB3	1.91	0.52
1:A:225:TYR:O	1:A:228:GLN:N	2.43	0.52
1:A:28:LEU:HD21	1:A:51:GLY:C	2.35	0.52
1:A:44:ALA:O	1:A:46:ARG:N	2.42	0.52
1:B:243:VAL:O	1:B:244:HIS:C	2.52	0.52
1:A:32:HIS:CE1	1:A:153:THR:HA	2.45	0.51
1:A:12:THR:HG21	1:A:61:ALA:O	2.10	0.51
1:A:240:THR:O	1:A:244:HIS:HD2	1.92	0.51
1:A:243:VAL:O	1:A:244:HIS:C	2.52	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:128:PRO:HB2	1:B:129:PRO:CD	2.40	0.51
1:A:77:ALA:O	1:A:78:VAL:C	2.53	0.51
1:A:186:ILE:O	1:A:187:LEU:C	2.54	0.51
1:B:234:TYR:CB	1:B:283:ASN:HB3	2.40	0.51
1:B:35:THR:HA	1:B:53:VAL:O	2.11	0.51
1:B:64:ALA:O	1:B:65:SER:CB	2.59	0.51
1:B:280:HIS:CD2	1:B:287:ALA:HB1	2.46	0.51
1:A:177:MET:O	1:A:180:LEU:HB2	2.11	0.51
1:B:77:ALA:O	1:B:78:VAL:C	2.53	0.51
1:A:77:ALA:HB1	2:A:301:NAP:N7A	2.27	0.50
1:B:186:ILE:O	1:B:187:LEU:C	2.54	0.50
1:B:234:TYR:O	1:B:235:GLU:C	2.54	0.50
1:A:82:ILE:O	1:A:83:ARG:C	2.53	0.50
1:A:127:PHE:CD2	1:B:221:TYR:CD2	3.00	0.50
1:A:181:GLY:O	1:A:182:ILE:C	2.55	0.50
1:A:126:ALA:HB2	1:A:136:ALA:HA	1.94	0.50
1:B:41:PRO:O	1:B:42:GLU:CB	2.59	0.50
1:A:226:ALA:HB3	1:A:227:PRO:HD3	1.93	0.49
1:A:208:ALA:O	1:A:211:ALA:HB3	2.12	0.49
1:B:181:GLY:O	1:B:182:ILE:C	2.56	0.49
1:B:218:ILE:O	1:B:221:TYR:N	2.46	0.49
1:A:44:ALA:O	1:A:45:GLU:C	2.56	0.49
1:B:62:VAL:HB	1:B:88:ALA:HB1	1.93	0.49
1:B:271:LEU:O	1:B:272:ALA:C	2.55	0.49
1:A:79:HIS:NE2	1:A:108:THR:OG1	2.45	0.49
2:A:301:NAP:N3A	2:A:301:NAP:O2X	2.46	0.48
1:B:82:ILE:O	1:B:83:ARG:C	2.55	0.48
1:B:279:GLY:O	1:B:281:ALA:N	2.45	0.48
1:A:134:ASP:HA	1:A:160:ALA:HB2	1.96	0.48
1:B:289:ILE:O	1:B:292:PHE:N	2.42	0.48
1:A:128:PRO:HD2	1:B:238:ASP:HA	1.94	0.48
1:A:141:SER:OG	1:A:169:HIS:HA	2.14	0.48
1:B:239:ALA:HB1	1:B:243:VAL:HG21	1.95	0.48
1:B:248:MET:HE2	1:B:268:PHE:HB3	1.94	0.48
1:A:64:ALA:O	1:A:65:SER:CB	2.61	0.48
1:B:214:MET:HA	1:B:217:ALA:HB3	1.96	0.48
1:A:171:LEU:HD23	1:A:171:LEU:HA	1.62	0.48
1:A:219:THR:CG2	1:B:215:ILE:CB	2.86	0.48
1:B:171:LEU:O	1:B:172:SER:C	2.54	0.48
1:A:261:HIS:CB	1:B:292:PHE:HA	2.44	0.48
1:A:75:TYR:CD1	1:A:100:ALA:HB2	2.48	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:295:PRO:HA	1:B:261:HIS:N	2.29	0.47
1:B:208:ALA:O	1:B:211:ALA:HB3	2.13	0.47
1:B:267:PHE:O	1:B:271:LEU:HD12	2.13	0.47
1:A:196:LEU:HD12	1:A:260:ILE:HD11	1.96	0.47
1:B:23:LEU:O	1:B:24:ALA:C	2.57	0.47
1:B:133:THR:O	1:B:159:PRO:HD2	2.15	0.47
1:A:23:LEU:O	1:A:24:ALA:C	2.58	0.47
1:A:218:ILE:O	1:A:221:TYR:N	2.48	0.47
1:B:72:VAL:HG22	1:B:98:THR:OG1	2.14	0.47
1:A:271:LEU:O	1:A:272:ALA:C	2.56	0.47
1:B:104:GLN:HA	1:B:107:GLU:HG3	1.96	0.47
1:A:289:ILE:HG23	1:A:290:GLU:N	2.29	0.47
1:B:249:GLU:O	1:B:250:LEU:C	2.58	0.47
1:B:287:ALA:O	1:B:288:MET:C	2.57	0.47
1:A:279:GLY:O	1:A:281:ALA:N	2.48	0.47
1:A:128:PRO:O	1:A:129:PRO:C	2.58	0.46
1:A:208:ALA:HB3	1:A:209:PRO:CD	2.46	0.46
1:B:249:GLU:OE1	1:B:253:GLU:CB	2.62	0.46
1:B:43:LYS:O	1:B:44:ALA:C	2.58	0.46
1:A:288:MET:HB2	1:B:191:LEU:HD13	1.98	0.46
1:B:180:LEU:O	1:B:183:MET:HB3	2.15	0.46
1:A:249:GLU:O	1:A:250:LEU:C	2.58	0.46
1:A:103:THR:HG21	1:B:257:HIS:CE1	2.51	0.46
1:A:287:ALA:O	1:A:288:MET:C	2.59	0.46
1:B:196:LEU:O	1:B:199:THR:OG1	2.31	0.46
1:A:219:THR:O	1:B:212:ASN:ND2	2.48	0.46
1:A:72:VAL:HG22	1:A:98:THR:OG1	2.16	0.45
1:A:250:LEU:O	1:A:254:GLU:N	2.33	0.45
1:A:31:GLY:O	1:A:32:HIS:C	2.59	0.45
1:B:241:MET:CE	1:B:285:TYR:HA	2.43	0.45
1:A:75:TYR:CE2	1:A:108:THR:HB	2.51	0.45
1:A:78:VAL:O	1:A:82:ILE:HG12	2.17	0.45
1:B:208:ALA:HB3	1:B:209:PRO:HD3	1.98	0.45
1:B:128:PRO:O	1:B:129:PRO:C	2.59	0.45
1:A:269:LYS:O	1:A:270:THR:C	2.60	0.45
1:B:81:LEU:O	1:B:82:ILE:C	2.59	0.45
1:B:264:LEU:HB3	1:B:265:PRO:HD3	1.98	0.45
1:A:32:HIS:CE1	1:A:153:THR:HG23	2.52	0.45
1:B:221:TYR:O	1:B:224:ALA:HB3	2.16	0.45
1:A:39:ARG:CB	2:A:301:NAP:C2A	2.95	0.45
1:A:128:PRO:HB2	1:A:129:PRO:HD3	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:285:TYR:CZ	1:B:288:MET:HG3	2.52	0.45
1:A:133:THR:HG23	1:A:135:GLY:N	2.30	0.45
1:A:225:TYR:O	1:A:226:ALA:C	2.60	0.45
1:A:241:MET:CE	1:A:285:TYR:HA	2.44	0.45
1:A:12:THR:OG1	1:A:65:SER:HB2	2.17	0.44
1:B:261:HIS:CE1	1:B:264:LEU:H	2.36	0.44
1:B:289:ILE:HG23	1:B:290:GLU:N	2.31	0.44
1:B:225:TYR:O	1:B:226:ALA:C	2.59	0.44
1:A:215:ILE:HA	1:A:218:ILE:HD12	1.98	0.44
1:A:222:VAL:HB	1:B:212:ASN:HD21	1.82	0.44
1:A:77:ALA:CB	2:A:301:NAP:N7A	2.80	0.44
1:B:222:VAL:O	1:B:225:TYR:HB2	2.18	0.44
1:B:179:VAL:O	1:B:182:ILE:HB	2.17	0.44
1:B:77:ALA:O	1:B:81:LEU:CD2	2.66	0.44
1:B:133:THR:HG23	1:B:135:GLY:N	2.31	0.44
1:B:175:TYR:O	1:B:176:ASP:C	2.61	0.44
1:B:31:GLY:O	1:B:32:HIS:C	2.61	0.44
1:A:81:LEU:O	1:A:82:ILE:C	2.60	0.43
1:A:82:ILE:HG23	1:A:89:LEU:HD13	1.99	0.43
1:A:140:TYR:OH	1:A:157:ILE:HD11	2.18	0.43
1:A:245:GLN:OE1	1:A:276:VAL:HG21	2.18	0.43
1:B:208:ALA:HB3	1:B:209:PRO:CD	2.48	0.43
1:A:289:ILE:O	1:A:292:PHE:N	2.48	0.43
1:A:39:ARG:CB	2:A:301:NAP:H2A	2.49	0.43
1:A:260:ILE:O	1:A:261:HIS:C	2.61	0.43
1:B:44:ALA:O	1:B:45:GLU:C	2.61	0.43
1:A:207:PHE:O	1:A:208:ALA:C	2.61	0.43
1:A:261:HIS:N	1:B:295:PRO:HA	2.34	0.43
1:B:128:PRO:CB	1:B:129:PRO:CD	2.97	0.43
1:A:214:MET:O	1:A:217:ALA:HB3	2.19	0.42
1:A:218:ILE:O	1:A:219:THR:C	2.62	0.42
1:A:234:TYR:O	1:A:235:GLU:C	2.61	0.42
1:B:218:ILE:C	1:B:220:GLU:N	2.77	0.42
1:A:17:GLY:O	1:A:18:LEU:C	2.62	0.42
1:A:133:THR:O	1:A:159:PRO:HD2	2.18	0.42
1:A:89:LEU:O	1:A:90:ALA:C	2.63	0.42
1:B:218:ILE:O	1:B:220:GLU:N	2.52	0.42
1:A:175:TYR:O	1:A:176:ASP:C	2.61	0.42
1:A:285:TYR:CZ	1:A:288:MET:HG3	2.55	0.42
1:B:238:ASP:OD1	1:B:238:ASP:N	2.52	0.42
1:B:115:ARG:O	1:B:116:ASN:C	2.62	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:210:LEU:HD12	1:B:210:LEU:C	2.43	0.42
1:A:289:ILE:HB	1:B:191:LEU:HD22	2.00	0.42
1:B:239:ALA:HB1	1:B:243:VAL:CG2	2.50	0.42
1:B:268:PHE:HA	1:B:271:LEU:HD13	2.02	0.42
1:A:15:GLY:O	1:A:20:GLY:HA3	2.20	0.42
1:A:272:ALA:N	1:A:288:MET:HE1	2.35	0.42
1:B:207:PHE:O	1:B:208:ALA:C	2.63	0.42
1:A:133:THR:OG1	1:A:134:ASP:N	2.53	0.41
1:A:267:PHE:O	1:A:268:PHE:C	2.61	0.41
1:B:260:ILE:O	1:B:261:HIS:C	2.63	0.41
1:A:58:LEU:O	1:A:62:VAL:HG23	2.20	0.41
1:B:168:ASP:OD1	1:B:170:GLY:N	2.53	0.41
1:A:75:TYR:HA	1:A:78:VAL:HB	2.02	0.41
1:B:285:TYR:O	1:B:286:ALA:C	2.63	0.41
1:A:32:HIS:HA	1:A:33:PRO:HD2	1.95	0.41
1:A:70:VAL:HG13	1:A:96:ASN:HA	2.02	0.41
1:A:218:ILE:C	1:A:220:GLU:N	2.76	0.41
1:B:58:LEU:O	1:B:62:VAL:HG23	2.21	0.41
1:A:295:PRO:HA	1:B:260:ILE:C	2.45	0.41
1:B:218:ILE:O	1:B:219:THR:C	2.63	0.41
1:A:77:ALA:O	1:A:81:LEU:CD2	2.67	0.41
1:A:122:GLY:HA2	1:A:140:TYR:HA	2.03	0.41
1:A:143:HIS:O	1:A:144:LYS:C	2.64	0.41
1:A:218:ILE:O	1:A:220:GLU:N	2.54	0.41
1:B:196:LEU:HD12	1:B:260:ILE:HD11	2.02	0.41
1:B:134:ASP:HA	1:B:160:ALA:HB2	2.03	0.41
1:A:229:VAL:O	1:A:230:ASP:C	2.64	0.40
1:A:183:MET:HB3	1:A:183:MET:HE2	1.81	0.40
1:B:126:ALA:HB3	1:B:131:ILE:HD13	2.03	0.40
1:A:64:ALA:O	1:A:65:SER:OG	2.38	0.40
1:A:186:ILE:C	1:A:188:ASN:N	2.80	0.40
1:B:265:PRO:O	1:B:266:ARG:C	2.65	0.40
1:B:123:ALA:N	1:B:139:LEU:O	2.52	0.40
1:A:28:LEU:C	1:A:30:GLY:H	2.30	0.40
1:A:59:GLU:O	1:A:60:SER:C	2.65	0.40
1:B:179:VAL:C	1:B:181:GLY:N	2.80	0.40
1:B:186:ILE:O	1:B:188:ASN:N	2.54	0.40
1:B:226:ALA:CB	1:B:227:PRO:CD	2.99	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	289/297 (97%)	210 (73%)	59 (20%)	20 (7%)	1	2
1	B	286/297 (96%)	208 (73%)	60 (21%)	18 (6%)	1	3
All	All	575/594 (97%)	418 (73%)	119 (21%)	38 (7%)	1	2

All (38) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	41	PRO
1	A	43	LYS
1	A	45	GLU
1	A	166	GLU
1	A	180	LEU
1	A	231	GLU
1	B	42	GLU
1	B	45	GLU
1	B	166	GLU
1	B	231	GLU
1	A	42	GLU
1	A	56	ASP
1	A	82	ILE
1	A	170	GLY
1	B	44	ALA
1	B	82	ILE
1	A	63	THR
1	A	219	THR
1	B	40	SER
1	B	63	THR
1	B	65	SER
1	B	78	VAL
1	B	219	THR
1	A	65	SER
1	A	220	GLU

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Mol	Chain	Res	Type
1	B	182	ILE
1	B	268	PHE
1	B	280	HIS
1	A	78	VAL
1	A	128	PRO
1	A	280	HIS
1	B	128	PRO
1	B	170	GLY
1	B	220	GLU
1	B	232	GLY
1	A	238	ASP
1	A	232	GLY
1	A	185	GLY

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	164/227 (72%)	141 (86%)	23 (14%)	3	8
1	B	111/227 (49%)	90 (81%)	21 (19%)	1	3
All	All	275/454 (61%)	231 (84%)	44 (16%)	2	6

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	VAL
1	A	28	LEU
1	A	48	VAL
1	A	53	VAL
1	A	58	LEU
1	A	60	SER
1	A	69	ILE
1	A	92	ARG
1	A	99	THR
1	A	103	THR

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Mol	Chain	Res	Type
1	A	108	THR
1	A	114	GLN
1	A	128	PRO
1	A	137	VAL
1	A	152	SER
1	A	157	ILE
1	A	162	THR
1	A	173	SER
1	A	183	MET
1	A	206	THR
1	A	223	THR
1	A	242	THR
1	A	291	LEU
1	B	48	VAL
1	B	98	THR
1	B	99	THR
1	B	103	THR
1	B	137	VAL
1	B	183	MET
1	B	196	LEU
1	B	206	THR
1	B	210	LEU
1	B	223	THR
1	B	229	VAL
1	B	238	ASP
1	B	242	THR
1	B	243	VAL
1	B	250	LEU
1	B	251	LEU
1	B	254	GLU
1	B	255	SER
1	B	264	LEU
1	B	280	HIS
1	B	289	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	212	ASN
1	A	244	HIS
1	B	212	ASN
1	B	257	HIS

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Mol	Chain	Res	Type
1	B	261	HIS
1	B	280	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 ⓘ

1 ligand is 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	A	301	-	45,52,52	0.86	3 (6%)	56,80,80	0.92	3 (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	A	301	-	-	4/31/67/67	0/5/5/5

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	301	NAP	C2N-N1N	3.03	1.38	1.35
2	A	301	NAP	P2B-O2B	2.51	1.64	1.59
2	A	301	NAP	C8A-N7A	-2.19	1.30	1.34

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	301	NAP	C6N-N1N-C2N	-2.46	119.73	121.97
2	A	301	NAP	C1B-N9A-C4A	2.45	130.94	126.64
2	A	301	NAP	O2N-PN-O1N	2.20	123.10	112.24

There are no chirality outliers.

All (4) torsion outliers are listed below:

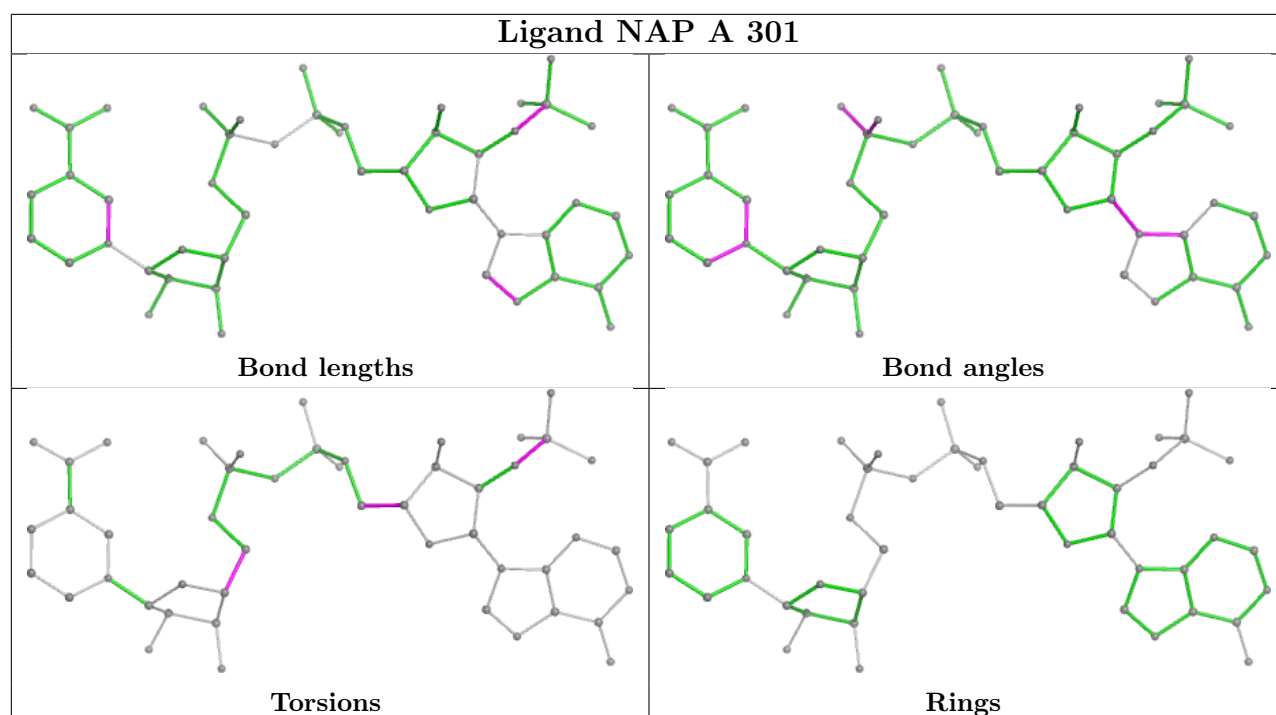
Mol	Chain	Res	Type	Atoms
2	A	301	NAP	O4B-C4B-C5B-O5B
2	A	301	NAP	O4D-C4D-C5D-O5D
2	A	301	NAP	C3D-C4D-C5D-O5D
2	A	301	NAP	C2B-O2B-P2B-O2X

There are no ring outliers.

1 monomer is involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	301	NAP	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 [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	291/297 (97%)	1.45	87 (29%) 1 1	74, 105, 159, 216	0
1	B	272/297 (91%)	2.19	123 (45%) 1 0	56, 169, 239, 269	0
All	All	563/594 (94%)	1.81	210 (37%) 1 1	56, 122, 229, 269	0

All (210) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	93	VAL	27.8
1	B	118	PRO	9.7
1	B	127	PHE	8.9
1	B	175	TYR	8.7
1	B	147	PHE	8.4
1	A	91	GLY	8.2
1	B	125	TRP	7.5
1	B	145	SER	7.3
1	B	117	ILE	7.2
1	B	225	TYR	7.2
1	B	39	ARG	6.9
1	B	66	PRO	6.7
1	B	78	VAL	6.6
1	B	82	ILE	6.5
1	A	36	VAL	6.4
1	A	23	LEU	6.3
1	B	135	GLY	6.2
1	B	52	ALA	6.1
1	A	175	TYR	6.1
1	B	11	VAL	5.9
1	B	89	LEU	5.9
1	A	267	PHE	5.8
1	B	24	ALA	5.8
1	A	190	PHE	5.8

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Mol	Chain	Res	Type	RSRZ
1	A	125	TRP	5.7
1	A	127	PHE	5.6
1	A	14	VAL	5.6
1	A	239	ALA	5.6
1	B	219	THR	5.5
1	B	218	ILE	5.4
1	A	44	ALA	5.3
1	B	92	ARG	5.2
1	B	69	ILE	5.2
1	A	215	ILE	5.1
1	B	144	LYS	5.0
1	A	53	VAL	4.9
1	B	72	VAL	4.9
1	B	84	PRO	4.8
1	B	149	ALA	4.8
1	B	215	ILE	4.7
1	A	94	LEU	4.7
1	A	71	CYS	4.5
1	B	71	CYS	4.5
1	A	89	LEU	4.5
1	B	271	LEU	4.5
1	A	235	GLU	4.5
1	A	43	LYS	4.4
1	A	268	PHE	4.4
1	B	256	GLU	4.4
1	B	210	LEU	4.4
1	B	176	ASP	4.4
1	B	23	LEU	4.3
1	B	19	MET	4.3
1	A	69	ILE	4.2
1	B	183	MET	4.2
1	B	112	ALA	4.2
1	A	187	LEU	4.1
1	B	274	ARG	4.1
1	A	68	VAL	4.1
1	B	70	VAL	4.1
1	A	67	LEU	4.0
1	B	59	GLU	4.0
1	B	177	MET	4.0
1	B	157	ILE	4.0
1	A	13	VAL	4.0
1	B	108	THR	3.9

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Mol	Chain	Res	Type	RSRZ
1	B	53	VAL	3.9
1	A	124	ILE	3.8
1	B	119	TYR	3.8
1	A	238	ASP	3.8
1	B	106	ARG	3.8
1	B	62	VAL	3.7
1	A	234	TYR	3.7
1	B	35	THR	3.7
1	B	190	PHE	3.7
1	A	282	GLU	3.6
1	A	54	LEU	3.6
1	B	20	GLY	3.6
1	A	95	VAL	3.6
1	B	22	ALA	3.6
1	A	140	TYR	3.6
1	A	70	VAL	3.5
1	A	157	ILE	3.5
1	A	79	HIS	3.5
1	A	154	LEU	3.5
1	A	222	VAL	3.5
1	B	48	VAL	3.5
1	B	164	TYR	3.5
1	B	221	TYR	3.4
1	B	85	VAL	3.4
1	B	18	LEU	3.4
1	B	81	LEU	3.4
1	B	182	ILE	3.4
1	A	82	ILE	3.3
1	A	201	LYS	3.3
1	B	124	ILE	3.3
1	B	237	THR	3.3
1	A	192	HIS	3.2
1	B	154	LEU	3.2
1	B	33	PRO	3.2
1	B	49	ALA	3.2
1	B	10	PRO	3.2
1	B	34	THR	3.2
1	A	81	LEU	3.2
1	B	186	ILE	3.1
1	A	297	ALA	3.1
1	A	93	VAL	3.1
1	B	222	VAL	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	75	TYR	3.1
1	A	171	LEU	3.1
1	B	120	LEU	3.1
1	A	233	ARG	3.1
1	A	285	TYR	3.1
1	A	186	ILE	3.0
1	A	271	LEU	3.0
1	A	240	THR	3.0
1	B	61	ALA	3.0
1	B	88	ALA	3.0
1	B	146	ALA	3.0
1	B	178	ALA	3.0
1	A	58	LEU	2.9
1	A	200	ALA	2.9
1	B	216	SER	2.9
1	A	97	LEU	2.9
1	A	194	ALA	2.9
1	B	191	LEU	2.9
1	B	99	THR	2.9
1	A	202	VAL	2.9
1	B	26	ALA	2.9
1	B	60	SER	2.9
1	B	75	TYR	2.9
1	A	247	ALA	2.9
1	B	40	SER	2.8
1	B	166	GLU	2.8
1	A	126	ALA	2.8
1	A	258	LEU	2.8
1	B	131	ILE	2.8
1	A	27	PHE	2.7
1	A	16	LEU	2.7
1	B	80	GLU	2.7
1	A	133	THR	2.7
1	B	141	SER	2.7
1	A	183	MET	2.7
1	B	236	ALA	2.7
1	B	44	ALA	2.6
1	B	268	PHE	2.6
1	A	188	ASN	2.6
1	A	224	ALA	2.6
1	B	192	HIS	2.6
1	B	58	LEU	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	56	ASP	2.5
1	B	45	GLU	2.5
1	B	285	TYR	2.5
1	A	111	TRP	2.4
1	A	21	GLN	2.4
1	A	168	ASP	2.4
1	A	138	LEU	2.4
1	B	264	LEU	2.4
1	A	207	PHE	2.4
1	A	281	ALA	2.4
1	A	15	GLY	2.4
1	B	167	GLU	2.4
1	B	251	LEU	2.4
1	A	131	ILE	2.4
1	B	64	ALA	2.3
1	B	239	ALA	2.3
1	B	122	GLY	2.3
1	A	51	GLY	2.3
1	B	116	ASN	2.3
1	B	293	ARG	2.3
1	B	136	ALA	2.3
1	A	274	ARG	2.2
1	A	216	SER	2.2
1	B	110	GLU	2.2
1	B	98	THR	2.2
1	B	162	THR	2.2
1	A	212	ASN	2.2
1	A	228	GLN	2.2
1	B	179	VAL	2.2
1	A	90	ALA	2.2
1	B	217	ALA	2.2
1	B	9	PRO	2.2
1	B	28	LEU	2.2
1	B	102	SER	2.2
1	B	103	THR	2.2
1	B	77	ALA	2.2
1	B	113	ALA	2.2
1	B	139	LEU	2.2
1	B	211	ALA	2.2
1	B	184	TRP	2.1
1	A	164	TYR	2.1
1	B	130	VAL	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	207	PHE	2.1
1	B	163	THR	2.1
1	B	214	MET	2.1
1	A	163	THR	2.1
1	B	87	SER	2.1
1	A	56	ASP	2.1
1	B	96	ASN	2.1
1	A	121	ASP	2.1
1	A	221	TYR	2.1
1	A	20	GLY	2.1
1	A	41	PRO	2.0
1	B	111	TRP	2.0
1	B	174	LEU	2.0
1	A	286	ALA	2.0
1	A	244	HIS	2.0
1	B	223	THR	2.0
1	A	174	LEU	2.0
1	B	196	LEU	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](#)

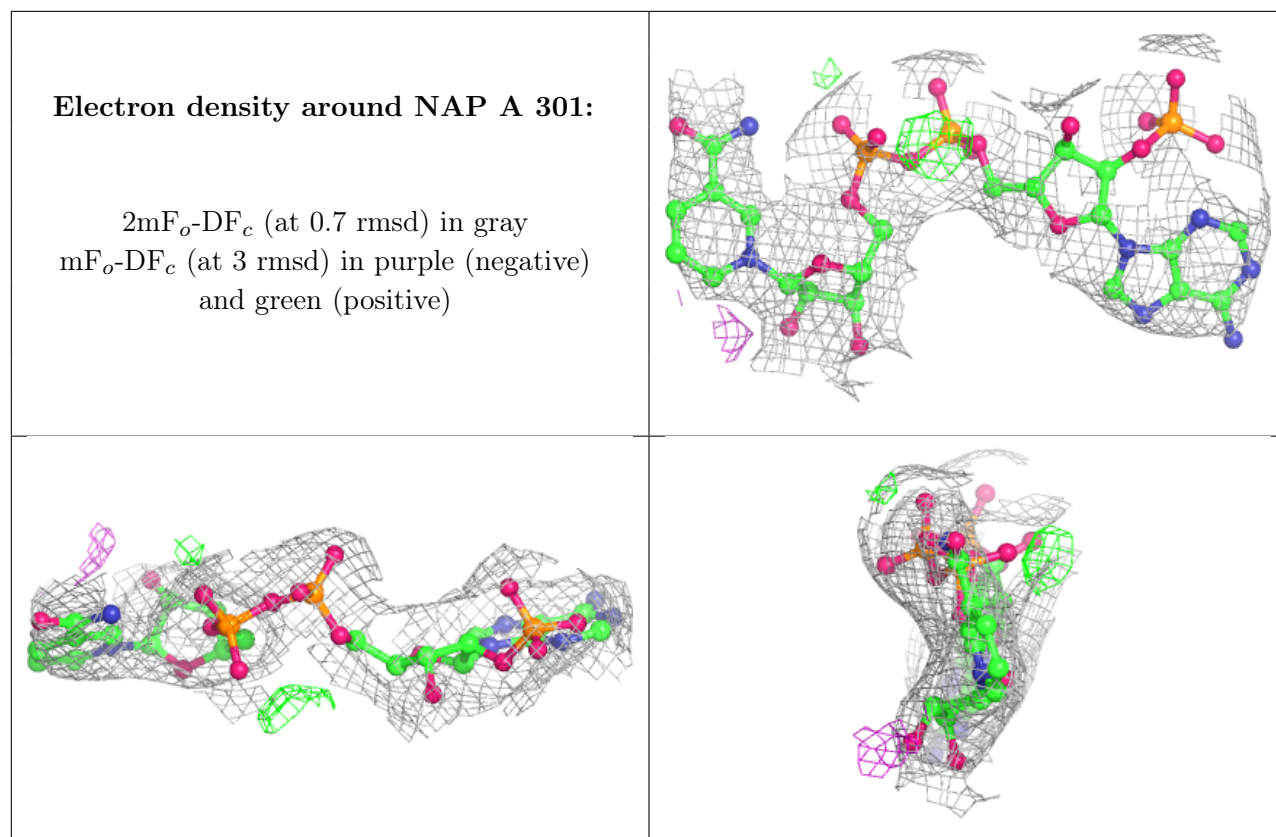
There are no oligosaccharides in this entry.

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(\AA^2)	Q<0.9
2	NAP	A	301	48/48	0.90	0.13	87,107,157,192	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.