



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 01:55 AM UTC

PDB ID : 5COB / pdb_00005cob
Title : Crystal structure of iridoid synthase in complex with NADP⁺ and 8-oxogeranial at 2.65-angstrom resolution
Authors : Qin, L.; Zhu, Y.; Ding, Z.; Zhang, X.; Ye, S.; Zhang, R.
Deposited on : 2015-07-20
Resolution : 2.65 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

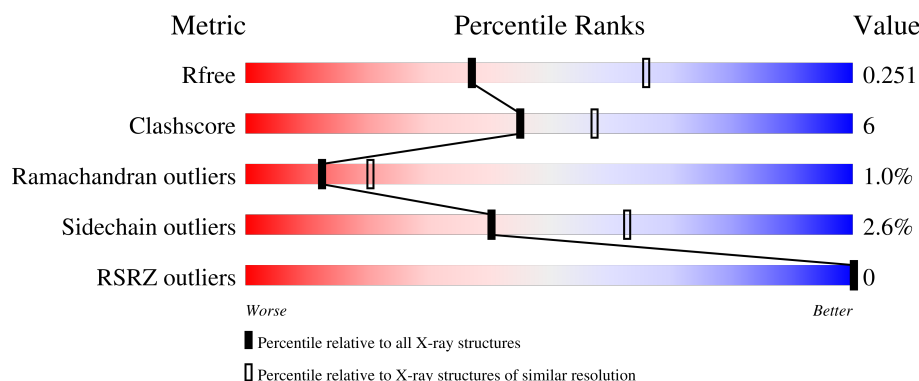
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

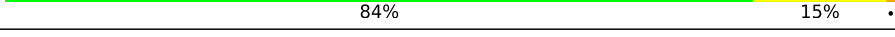
The reported resolution of this entry is 2.65 Å.

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	1110 (2.66-2.66)
Clashscore	190562	1141 (2.66-2.66)
Ramachandran outliers	187476	1126 (2.66-2.66)
Sidechain outliers	187428	1126 (2.66-2.66)
RSRZ outliers	180081	1110 (2.66-2.66)

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	365	 85% 14% .
1	B	365	 81% 19%
1	C	365	 85% 14% .
1	D	365	 84% 15% .

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 12059 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 Iridoid synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	365	Total	C	N	O	S	0	0	0
			2895	1864	477	534	20			
1	B	365	Total	C	N	O	S	0	0	0
			2895	1864	477	534	20			
1	C	365	Total	C	N	O	S	0	0	0
			2895	1864	477	534	20			
1	D	365	Total	C	N	O	S	0	0	0
			2895	1864	477	534	20			

There are 8 discrepancies between the modelled and reference sequences:

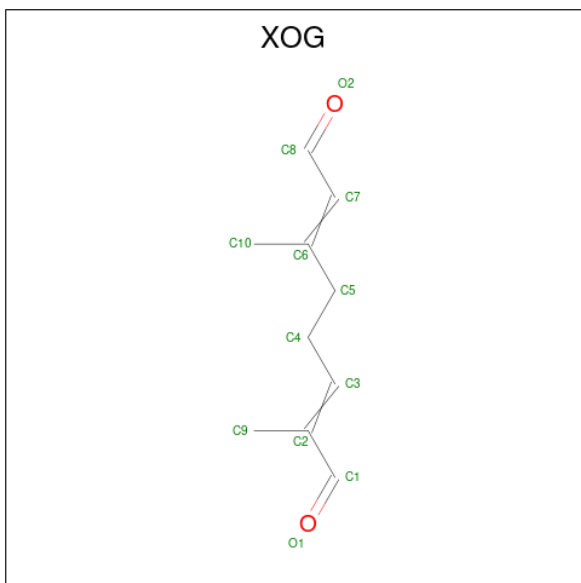
Chain	Residue	Modelled	Actual	Comment	Reference
A	24	MET	-	expression tag	UNP K7WDL7
A	25	ALA	-	expression tag	UNP K7WDL7
B	24	MET	-	expression tag	UNP K7WDL7
B	25	ALA	-	expression tag	UNP K7WDL7
C	24	MET	-	expression tag	UNP K7WDL7
C	25	ALA	-	expression tag	UNP K7WDL7
D	24	MET	-	expression tag	UNP K7WDL7
D	25	ALA	-	expression tag	UNP K7WDL7

- Molecule 2 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (CCD ID: NAP) (formula: C₂₁H₂₈N₇O₁₇P₃).



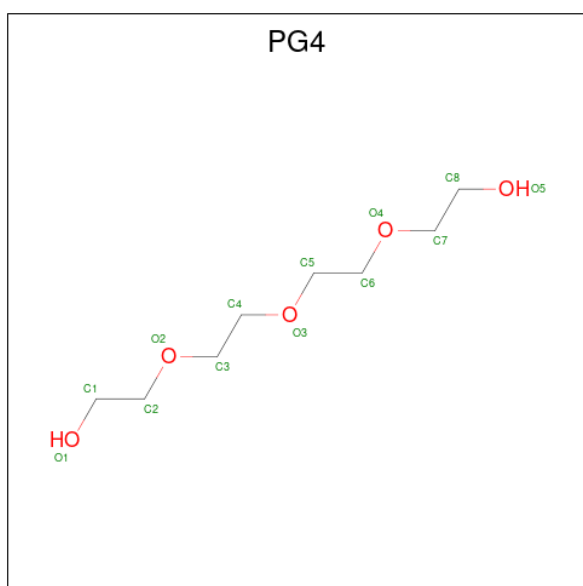
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 48	C 21	N 7	O 17	P 3	0	0
2	B	1	Total 48	C 21	N 7	O 17	P 3	0	0
2	C	1	Total 48	C 21	N 7	O 17	P 3	0	0
2	D	1	Total 48	C 21	N 7	O 17	P 3	0	0

- Molecule 3 is (2E,6E)-2,6-dimethylocta-2,6-dienedial (CCD ID: XOG) (formula: C₁₀H₁₄O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			12	10	2		
3	B	1	Total	C	O	0	0
			12	10	2		
3	C	1	Total	C	O	0	0
			12	10	2		
3	D	1	Total	C	O	0	0
			12	10	2		

- Molecule 4 is TETRAETHYLENE GLYCOL (CCD ID: PG4) (formula: $C_8H_{18}O_5$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			13	8	5		
4	B	1	Total	C	O	0	0
			13	8	5		
4	C	1	Total	C	O	0	0
			13	8	5		
4	D	1	Total	C	O	0	0
			13	8	5		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	48	Total	O	0	0
			48	48		
5	B	52	Total	O	0	0
			52	52		

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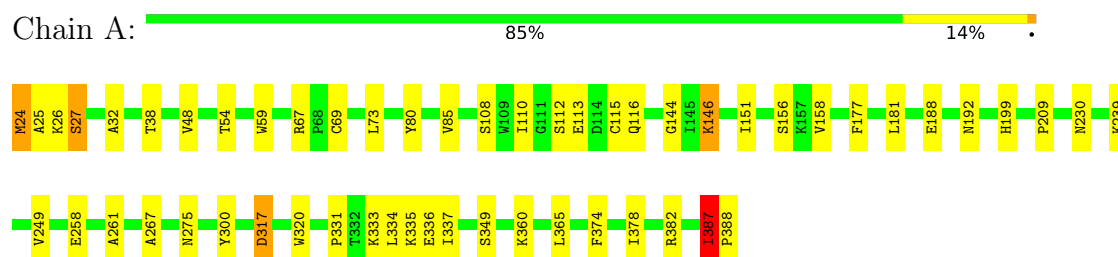
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	C	41	Total	O	0	0
			41	41		
5	D	46	Total	O	0	0
			46	46		

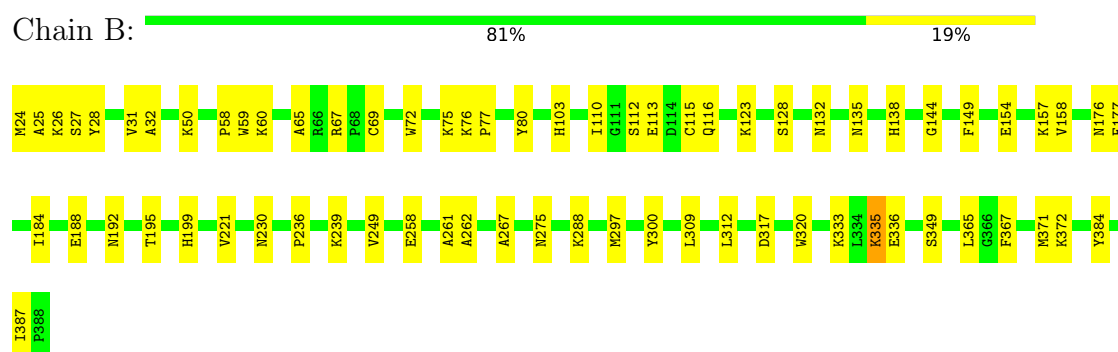
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

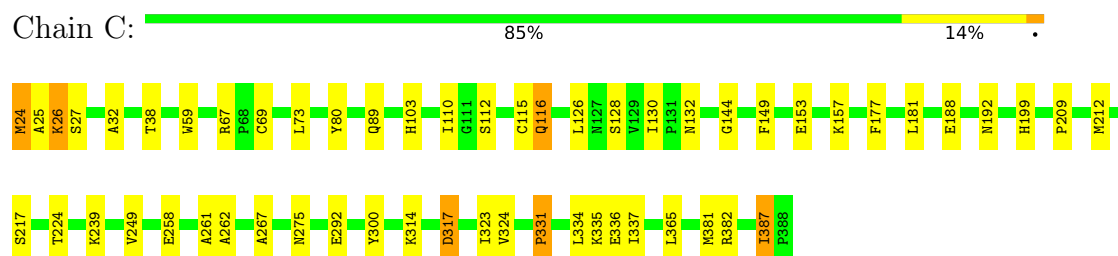
- Molecule 1: Iridoid synthase



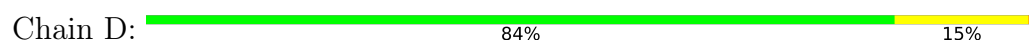
- Molecule 1: Iridoid synthase



- Molecule 1: Iridoid synthase



- Molecule 1: Iridoid synthase



K24	A25	K26	S27	V31	T38	D53	K60	R67	P69	C69	L73	Y80	I110	G111	S112	C115	Q116	K123	S128	N132	G144	I145	K146	F149	G150	S156	K157	W158	V159	P160	F177	Y178	I184	E188	T189	N192	W196	H199
P209	C210	N230	G237	S238	K239	V249	E258	A261	D264	P265	K266	A267	N275	K288	Y300	D317	L334	K335	E336	I337	F347	G348	S349	L365	M371	R362	L387	P388										

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	65.99Å 171.91Å 66.03Å 90.00° 89.72° 90.00°	Depositor
Resolution (Å)	36.14 – 2.65 36.14 – 2.65	Depositor EDS
% Data completeness (in resolution range)	98.6 (36.14-2.65) 93.4 (36.14-2.65)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.09 (at 2.65Å)	Xtriage
Refinement program	PHENIX 1.8.2_1309	Depositor
R, R_{free}	0.194 , 0.245 0.202 , 0.251	Depositor DCC
R_{free} test set	2114 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å ²)	41.3	Xtriage
Anisotropy	0.755	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 36.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.417 for -l,k,h 0.418 for h,-k,-l 0.477 for -l,-k,-h	Xtriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	12059	wwPDB-VP
Average B, all atoms (Å ²)	61.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.22% 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, PG4, XOG

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	0.28	0/2970	0.69	1/4031 (0.0%)
1	B	0.28	0/2970	0.70	0/4031
1	C	0.28	0/2970	0.72	1/4031 (0.0%)
1	D	0.28	0/2970	0.71	1/4031 (0.0%)
All	All	0.28	0/11880	0.71	3/16124 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	387	ILE	N-CA-C	-10.93	100.22	109.19
1	A	387	ILE	N-CA-C	5.21	120.12	108.88
1	D	387	ILE	N-CA-C	5.01	119.70	108.88

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	2895	0	2866	35	0
1	B	2895	0	2866	42	0
1	C	2895	0	2866	35	0
1	D	2895	0	2866	33	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	48	0	23	4	0
2	B	48	0	23	2	0
2	C	48	0	23	3	0
2	D	48	0	23	1	0
3	A	12	0	14	4	0
3	B	12	0	14	3	0
3	C	12	0	14	3	0
3	D	12	0	14	5	0
4	A	13	0	18	2	0
4	B	13	0	18	1	0
4	C	13	0	18	0	0
4	D	13	0	18	1	0
5	A	48	0	0	2	0
5	B	52	0	0	1	0
5	C	41	0	0	1	0
5	D	46	0	0	1	0
All	All	12059	0	11684	139	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 6.

All (139) 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:67:ARG:NH2	2:C:1001:NAP:O2X	2.22	0.72
1:D:249:VAL:HB	1:D:275:ASN:HA	1.70	0.72
1:A:249:VAL:HB	1:A:275:ASN:HA	1.74	0.70
1:D:67:ARG:NH1	1:D:336:GLU:OE2	2.25	0.70
1:D:67:ARG:NH2	2:D:1001:NAP:O2X	2.24	0.69
1:C:67:ARG:NH1	1:C:336:GLU:OE2	2.24	0.69
1:B:236:PRO:HG3	1:B:309:LEU:HD13	1.76	0.68
1:B:24:MET:N	1:B:28:TYR:O	2.27	0.67
1:B:249:VAL:HB	1:B:275:ASN:HA	1.77	0.65
1:C:249:VAL:HB	1:C:275:ASN:HA	1.79	0.64
1:C:365:LEU:HD13	1:D:365:LEU:HD13	1.80	0.63
1:B:72:TRP:O	1:C:116:GLN:NE2	2.32	0.63
1:A:113:GLU:OE2	1:A:146:LYS:NZ	2.31	0.63
1:A:67:ARG:NH1	1:A:336:GLU:OE2	2.32	0.63
1:B:67:ARG:NH1	1:B:336:GLU:OE2	2.33	0.62
1:C:188:GLU:O	1:C:192:ASN:ND2	2.31	0.62
1:B:67:ARG:NH2	2:B:1001:NAP:O2X	2.26	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:317:ASP:OD1	1:A:317:ASP:N	2.33	0.61
1:D:188:GLU:O	1:D:192:ASN:ND2	2.31	0.61
1:B:76:LYS:HD3	1:C:116:GLN:HE21	1.65	0.60
1:B:24:MET:HE3	1:B:135:ASN:HD22	1.66	0.60
1:D:110:ILE:HG23	1:D:112:SER:H	1.67	0.58
1:A:144:GLY:HA3	3:A:1002:XOG:H2	1.86	0.57
1:B:309:LEU:HD12	1:B:312:LEU:HD11	1.87	0.57
1:B:115:CYS:HA	1:B:177:PHE:HB3	1.86	0.57
1:A:67:ARG:NH2	2:A:1001:NAP:O2X	2.29	0.56
1:C:317:ASP:OD1	1:C:317:ASP:N	2.31	0.56
1:B:110:ILE:HG23	1:B:112:SER:H	1.70	0.56
1:B:333:LYS:HG2	1:B:335:LYS:HD3	1.88	0.56
1:B:188:GLU:O	1:B:192:ASN:ND2	2.35	0.56
1:D:382:ARG:HA	1:D:388:PRO:HD3	1.88	0.55
1:A:110:ILE:HG23	1:A:112:SER:H	1.70	0.55
1:C:25:ALA:O	1:C:27:SER:N	2.41	0.54
1:A:382:ARG:HA	1:A:388:PRO:HD3	1.90	0.54
1:D:239:LYS:HB2	1:D:300:TYR:CZ	2.43	0.53
1:B:288:LYS:NZ	5:B:1105:HOH:O	2.42	0.53
1:D:144:GLY:HA3	3:D:1002:XOG:H2	1.90	0.53
1:C:110:ILE:HG23	1:C:112:SER:H	1.73	0.52
1:B:75:LYS:HB3	1:C:115:CYS:HB2	1.89	0.52
1:B:144:GLY:HA3	3:B:1002:XOG:H2	1.91	0.52
1:D:146:LYS:HB2	3:D:1002:XOG:O2	2.09	0.52
1:C:144:GLY:HA3	3:C:1002:XOG:H2	1.92	0.52
1:D:31:VAL:HG22	1:D:60:LYS:HB3	1.93	0.51
1:C:199:HIS:CE1	1:C:258:GLU:HA	2.45	0.51
1:C:128:SER:O	1:C:132:ASN:ND2	2.44	0.51
1:D:199:HIS:CE1	1:D:258:GLU:HA	2.46	0.50
1:A:25:ALA:O	1:A:27:SER:N	2.45	0.49
1:C:224:THR:HG22	1:C:323:ILE:HD13	1.93	0.49
1:C:103:HIS:CD2	1:C:262:ALA:HB1	2.48	0.49
1:D:25:ALA:O	1:D:27:SER:N	2.45	0.49
1:B:349:SER:OG	3:B:1002:XOG:O1	2.27	0.49
1:B:50:LYS:HG3	1:B:77:PRO:HG2	1.94	0.48
1:C:217:SER:HA	1:C:381:MET:HE1	1.95	0.48
1:A:349:SER:OG	3:A:1002:XOG:O1	2.23	0.48
1:B:31:VAL:HG22	1:B:60:LYS:HB3	1.95	0.48
1:D:115:CYS:HA	1:D:177:PHE:HB3	1.96	0.48
1:C:103:HIS:HD2	1:C:262:ALA:HB1	1.80	0.47
1:B:128:SER:O	1:B:132:ASN:ND2	2.45	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:387:ILE:H	1:A:388:PRO:CD	2.28	0.47
1:A:188:GLU:O	1:A:192:ASN:ND2	2.41	0.47
1:A:115:CYS:HA	1:A:177:PHE:HB3	1.95	0.47
1:A:69:CYS:HB2	1:A:80:TYR:CD2	2.50	0.46
1:A:146:LYS:HB2	3:A:1002:XOG:O2	2.14	0.46
1:A:209:PRO:HA	1:A:337:ILE:HG13	1.98	0.46
1:A:108:SER:HB2	2:A:1001:NAP:O4B	2.15	0.46
1:A:317:ASP:OD2	5:A:1101:HOH:O	2.21	0.46
1:B:149:PHE:HZ	3:B:1002:XOG:H14	1.79	0.46
1:B:384:TYR:CE1	1:C:116:GLN:HB2	2.51	0.46
1:C:149:PHE:HZ	3:C:1002:XOG:H14	1.81	0.46
1:B:239:LYS:HB2	1:B:300:TYR:CZ	2.51	0.45
1:A:199:HIS:CE1	1:A:258:GLU:HA	2.52	0.45
3:A:1002:XOG:H8	4:A:1003:PG4:H82	1.98	0.45
1:B:28:TYR:HA	1:B:58:PRO:HG2	1.99	0.45
1:B:138:HIS:ND1	1:B:195:THR:OG1	2.38	0.45
1:C:69:CYS:HB2	1:C:80:TYR:CD2	2.52	0.45
1:D:116:GLN:H	1:D:116:GLN:HG2	1.48	0.45
1:A:38:THR:HG22	1:A:73:LEU:HD12	1.99	0.45
1:C:382:ARG:HA	1:C:387:ILE:O	2.17	0.45
1:A:24:MET:HB3	1:A:25:ALA:H	1.59	0.45
1:A:67:ARG:HD2	2:A:1001:NAP:O2X	2.17	0.44
1:A:333:LYS:HG2	1:A:335:LYS:HD3	1.99	0.44
1:B:25:ALA:O	1:B:27:SER:N	2.50	0.44
1:C:209:PRO:HA	1:C:337:ILE:HG13	1.98	0.44
1:D:387:ILE:H	1:D:388:PRO:CD	2.30	0.44
1:A:261:ALA:HA	1:A:267:ALA:HB3	1.99	0.44
1:D:288:LYS:HG3	1:D:371:MET:SD	2.58	0.44
1:D:189:THR:HG21	1:D:196:TRP:HB3	1.99	0.44
1:B:103:HIS:CD2	1:B:138:HIS:HB3	2.53	0.44
1:C:292:GLU:O	1:C:382:ARG:NH2	2.35	0.44
1:C:126:LEU:HD22	1:C:130:ILE:HD11	2.00	0.44
1:C:38:THR:HG21	1:C:67:ARG:HD3	1.99	0.44
1:B:199:HIS:CE1	1:B:258:GLU:HA	2.53	0.43
1:B:103:HIS:CD2	1:B:262:ALA:HB1	2.53	0.43
1:D:349:SER:OG	3:D:1002:XOG:O1	2.26	0.43
1:A:334:LEU:HB3	5:A:1101:HOH:O	2.18	0.43
1:B:67:ARG:HD2	2:B:1001:NAP:O2X	2.19	0.43
1:D:150:GLY:HA3	4:D:1003:PG4:H12	2.00	0.43
1:B:221:VAL:HG22	1:B:320:TRP:CH2	2.54	0.43
1:B:69:CYS:HB2	1:B:80:TYR:CD2	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:334:LEU:HB3	5:C:1106:HOH:O	2.19	0.42
1:A:177:PHE:O	1:A:181:LEU:HG	2.19	0.42
1:A:374:PHE:O	1:A:378:ILE:HG13	2.20	0.42
1:D:123:LYS:HG3	1:D:184:ILE:HD13	2.02	0.42
1:D:69:CYS:HB2	1:D:80:TYR:CD2	2.54	0.42
1:C:239:LYS:HB2	1:C:300:TYR:CZ	2.55	0.42
1:A:365:LEU:HD13	1:B:365:LEU:HD13	2.00	0.42
1:C:177:PHE:O	1:C:181:LEU:HG	2.19	0.42
1:D:149:PHE:HZ	3:D:1002:XOG:H14	1.83	0.42
1:A:239:LYS:HB2	1:A:300:TYR:CZ	2.54	0.42
1:B:32:ALA:HB2	1:B:59:TRP:CZ3	2.54	0.42
1:B:123:LYS:HG3	1:B:184:ILE:HD13	2.01	0.42
2:C:1001:NAP:C2N	3:C:1002:XOG:H7	2.50	0.42
1:D:178:TYR:OH	3:D:1002:XOG:O2	2.31	0.42
1:A:85:VAL:HG23	2:A:1001:NAP:H2A	2.01	0.42
1:C:32:ALA:HB2	1:C:59:TRP:CE3	2.55	0.42
1:B:372:LYS:O	1:C:89:GLN:NE2	2.52	0.42
1:C:24:MET:HB3	1:C:25:ALA:H	1.61	0.42
1:D:38:THR:HG22	1:D:73:LEU:HD12	2.02	0.42
1:D:317:ASP:OD1	1:D:317:ASP:N	2.34	0.42
1:B:261:ALA:HA	1:B:267:ALA:HB3	2.01	0.41
1:B:288:LYS:HG3	1:B:371:MET:SD	2.59	0.41
1:A:48:VAL:O	1:A:54:THR:OG1	2.31	0.41
1:D:261:ALA:HA	1:D:267:ALA:HB3	2.01	0.41
1:D:264:ASP:HA	1:D:265:PRO:HD3	1.86	0.41
1:D:159:VAL:HA	1:D:160:PRO:HD2	1.95	0.41
1:A:32:ALA:HB2	1:A:59:TRP:CZ3	2.56	0.41
1:C:261:ALA:HA	1:C:267:ALA:HB3	2.03	0.41
1:A:360:LYS:HB3	1:B:367:PHE:CZ	2.56	0.41
1:B:335:LYS:HE2	1:B:335:LYS:HB2	1.89	0.41
1:D:334:LEU:HB3	5:D:1108:HOH:O	2.21	0.41
1:B:176:ASN:OD1	4:B:1003:PG4:H32	2.21	0.41
1:C:324:VAL:HG11	1:C:331:PRO:HA	2.03	0.41
1:B:65:ALA:HB3	1:B:80:TYR:OH	2.21	0.40
1:C:67:ARG:HD2	2:C:1001:NAP:O2X	2.20	0.40
1:D:128:SER:O	1:D:132:ASN:ND2	2.49	0.40
1:D:237:GLY:HA3	1:D:347:PHE:HB3	2.04	0.40
1:A:320:TRP:HH2	1:A:387:ILE:HD12	1.86	0.40
1:D:209:PRO:HA	1:D:337:ILE:HG13	2.03	0.40
1:A:146:LYS:HD2	4:A:1003:PG4:H42	2.02	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	363/365 (100%)	339 (93%)	19 (5%)	5 (1%)	9	14
1	B	363/365 (100%)	345 (95%)	16 (4%)	2 (1%)	21	34
1	C	363/365 (100%)	343 (94%)	17 (5%)	3 (1%)	16	27
1	D	363/365 (100%)	344 (95%)	15 (4%)	4 (1%)	11	19
All	All	1452/1460 (100%)	1371 (94%)	67 (5%)	14 (1%)	12	20

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	387	ILE
1	D	156	SER
1	A	26	LYS
1	B	26	LYS
1	C	26	LYS
1	D	26	LYS
1	D	387	ILE
1	A	156	SER
1	A	158	VAL
1	B	158	VAL
1	D	157	LYS
1	C	157	LYS
1	A	331	PRO
1	C	331	PRO

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	323/323 (100%)	315 (98%)	8 (2%)	42	64
1	B	323/323 (100%)	314 (97%)	9 (3%)	38	60
1	C	323/323 (100%)	314 (97%)	9 (3%)	38	60
1	D	323/323 (100%)	316 (98%)	7 (2%)	45	67
All	All	1292/1292 (100%)	1259 (97%)	33 (3%)	40	63

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	24	MET
1	A	27	SER
1	A	116	GLN
1	A	146	LYS
1	A	151	ILE
1	A	230	ASN
1	A	317	ASP
1	A	387	ILE
1	B	113	GLU
1	B	116	GLN
1	B	154	GLU
1	B	157	LYS
1	B	230	ASN
1	B	297	MET
1	B	317	ASP
1	B	335	LYS
1	B	387	ILE
1	C	24	MET
1	C	26	LYS
1	C	73	LEU
1	C	116	GLN
1	C	153	GLU
1	C	212	MET
1	C	314	LYS
1	C	317	ASP
1	C	335	LYS
1	D	53	ASP
1	D	73	LEU
1	D	116	GLN
1	D	210	CYS
1	D	230	ASN
1	D	317	ASP
1	D	387	ILE

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	228	HIS
1	B	135	ASN
1	B	228	HIS
1	B	243	ASN
1	C	116	GLN
1	C	243	ASN
1	C	327	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](#)

12 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	$\# Z > 2$
3	XOG	C	1002	-	11,11,11	3.92	4 (36%)	12,12,12	5.90	4 (33%)
4	PG4	C	1003	-	12,12,12	0.45	0	11,11,11	0.29	0
2	NAP	A	1001	-	50,52,52	1.57	6 (12%)	71,80,80	1.56	12 (16%)
4	PG4	A	1003	-	12,12,12	0.46	0	11,11,11	0.28	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NAP	D	1001	-	50,52,52	1.57	6 (12%)	71,80,80	1.56	12 (16%)
2	NAP	C	1001	-	50,52,52	1.58	6 (12%)	71,80,80	1.55	12 (16%)
3	XOG	D	1002	-	11,11,11	3.96	4 (36%)	12,12,12	5.71	3 (25%)
4	PG4	D	1003	-	12,12,12	0.45	0	11,11,11	0.29	0
3	XOG	A	1002	-	11,11,11	3.93	4 (36%)	12,12,12	5.78	3 (25%)
4	PG4	B	1003	-	12,12,12	0.45	0	11,11,11	0.30	0
3	XOG	B	1002	-	11,11,11	3.96	4 (36%)	12,12,12	5.76	3 (25%)
2	NAP	B	1001	-	50,52,52	1.57	6 (12%)	71,80,80	1.56	11 (15%)

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
3	XOG	C	1002	-	-	3/11/11/11	-
4	PG4	C	1003	-	-	0/10/10/10	-
2	NAP	A	1001	-	-	11/35/67/67	0/5/5/5
4	PG4	A	1003	-	-	0/10/10/10	-
2	NAP	D	1001	-	-	15/35/67/67	0/5/5/5
2	NAP	C	1001	-	-	9/35/67/67	0/5/5/5
3	XOG	D	1002	-	-	3/11/11/11	-
4	PG4	D	1003	-	-	0/10/10/10	-
3	XOG	A	1002	-	-	3/11/11/11	-
4	PG4	B	1003	-	-	0/10/10/10	-
3	XOG	B	1002	-	-	3/11/11/11	-
2	NAP	B	1001	-	-	9/35/67/67	0/5/5/5

All (40) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1002	XOG	O1-C1	9.84	1.43	1.22
3	C	1002	XOG	O1-C1	9.82	1.43	1.22
3	D	1002	XOG	O1-C1	9.80	1.43	1.22
3	A	1002	XOG	O1-C1	9.75	1.43	1.22
3	D	1002	XOG	O2-C8	7.10	1.43	1.22
3	A	1002	XOG	O2-C8	7.09	1.43	1.22
3	B	1002	XOG	O2-C8	7.06	1.43	1.22
3	C	1002	XOG	O2-C8	7.02	1.43	1.22

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	1001	NAP	C7N-N7N	-5.75	1.22	1.33
2	B	1001	NAP	C7N-N7N	-5.74	1.22	1.33
2	A	1001	NAP	C7N-N7N	-5.71	1.22	1.33
2	D	1001	NAP	C7N-N7N	-5.68	1.22	1.33
2	D	1001	NAP	O7N-C7N	5.19	1.33	1.24
2	B	1001	NAP	O7N-C7N	5.18	1.33	1.24
2	A	1001	NAP	O7N-C7N	5.16	1.33	1.24
2	C	1001	NAP	O7N-C7N	5.14	1.33	1.24
2	C	1001	NAP	C5A-C4A	4.76	1.47	1.39
2	D	1001	NAP	C5A-C4A	4.75	1.47	1.39
2	A	1001	NAP	C5A-C4A	4.70	1.47	1.39
2	B	1001	NAP	C5A-C4A	4.67	1.47	1.39
3	B	1002	XOG	C1-C2	4.09	1.51	1.46
3	D	1002	XOG	C1-C2	4.04	1.51	1.46
3	A	1002	XOG	C1-C2	3.96	1.50	1.46
3	C	1002	XOG	C1-C2	3.84	1.50	1.46
3	D	1002	XOG	C7-C8	3.04	1.51	1.44
3	A	1002	XOG	C7-C8	2.96	1.51	1.44
3	B	1002	XOG	C7-C8	2.93	1.51	1.44
3	C	1002	XOG	C7-C8	2.89	1.51	1.44
2	C	1001	NAP	C5A-C6A	2.80	1.48	1.41
2	B	1001	NAP	C5A-C6A	2.79	1.48	1.41
2	D	1001	NAP	C5A-C6A	2.77	1.48	1.41
2	A	1001	NAP	C5A-C6A	2.73	1.48	1.41
2	B	1001	NAP	C8A-N7A	2.34	1.36	1.31
2	C	1001	NAP	C8A-N7A	2.33	1.36	1.31
2	D	1001	NAP	C8A-N7A	2.28	1.36	1.31
2	A	1001	NAP	C8A-N7A	2.23	1.36	1.31
2	D	1001	NAP	C5A-N7A	-2.19	1.35	1.39
2	C	1001	NAP	C5A-N7A	-2.17	1.35	1.39
2	A	1001	NAP	C5A-N7A	-2.14	1.35	1.39
2	B	1001	NAP	C5A-N7A	-2.13	1.35	1.39

All (60) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	1002	XOG	O1-C1-C2	-17.88	106.85	125.11
3	A	1002	XOG	O1-C1-C2	-17.18	107.57	125.11
3	B	1002	XOG	O1-C1-C2	-17.17	107.58	125.11
3	D	1002	XOG	O1-C1-C2	-17.06	107.69	125.11
3	A	1002	XOG	O2-C8-C7	-9.34	111.13	123.68
3	B	1002	XOG	O2-C8-C7	-9.19	111.34	123.68

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	1002	XOG	O2-C8-C7	-9.03	111.56	123.68
3	C	1002	XOG	O2-C8-C7	-8.74	111.94	123.68
2	A	1001	NAP	C5A-C4A-N3A	-5.50	119.15	126.72
2	B	1001	NAP	C5A-C4A-N3A	-5.49	119.15	126.72
2	C	1001	NAP	C5A-C4A-N3A	-5.42	119.26	126.72
2	D	1001	NAP	C5A-C4A-N3A	-5.36	119.34	126.72
2	B	1001	NAP	N3A-C4A-N9A	4.43	134.71	127.17
2	A	1001	NAP	N3A-C4A-N9A	4.43	134.70	127.17
2	D	1001	NAP	N3A-C4A-N9A	4.34	134.55	127.17
2	C	1001	NAP	N3A-C4A-N9A	4.33	134.53	127.17
2	B	1001	NAP	C2A-N3A-C4A	3.63	120.70	111.83
2	A	1001	NAP	C2A-N3A-C4A	3.58	120.58	111.83
2	C	1001	NAP	C4A-C5A-N7A	-3.56	106.51	110.58
2	C	1001	NAP	C2A-N3A-C4A	3.56	120.52	111.83
2	D	1001	NAP	C2A-N3A-C4A	3.51	120.40	111.83
2	A	1001	NAP	C4A-C5A-N7A	-3.51	106.57	110.58
2	B	1001	NAP	C4A-C5A-N7A	-3.51	106.57	110.58
2	D	1001	NAP	C4A-C5A-N7A	-3.49	106.59	110.58
2	B	1001	NAP	N3A-C2A-N1A	-3.35	123.51	128.58
2	A	1001	NAP	N3A-C2A-N1A	-3.33	123.54	128.58
2	C	1001	NAP	N3A-C2A-N1A	-3.27	123.63	128.58
2	D	1001	NAP	N3A-C2A-N1A	-3.21	123.72	128.58
2	B	1001	NAP	C4A-N9A-C8A	2.99	108.88	105.74
2	D	1001	NAP	C4A-N9A-C8A	2.98	108.87	105.74
2	C	1001	NAP	C4A-N9A-C8A	2.94	108.83	105.74
2	A	1001	NAP	C4A-N9A-C8A	2.86	108.74	105.74
2	B	1001	NAP	C5A-N7A-C8A	2.65	107.62	103.45
3	C	1002	XOG	C10-C6-C5	2.64	119.81	115.23
2	C	1001	NAP	C5A-N7A-C8A	2.64	107.60	103.45
2	A	1001	NAP	P2B-O2B-C2B	-2.63	116.41	123.43
2	A	1001	NAP	C5A-N7A-C8A	2.62	107.57	103.45
2	D	1001	NAP	C5A-N7A-C8A	2.59	107.52	103.45
2	B	1001	NAP	P2B-O2B-C2B	-2.53	116.68	123.43
3	B	1002	XOG	C10-C6-C5	2.51	119.59	115.23
3	A	1002	XOG	C10-C6-C5	2.47	119.51	115.23
2	D	1001	NAP	P2B-O2B-C2B	-2.37	117.09	123.43
3	D	1002	XOG	C10-C6-C5	2.32	119.25	115.23
2	A	1001	NAP	O3X-P2B-O2X	2.31	116.45	107.80
2	C	1001	NAP	P2B-O2B-C2B	-2.27	117.36	123.43
2	C	1001	NAP	C6A-C5A-N7A	2.23	136.39	132.09
2	B	1001	NAP	N9A-C8A-N7A	-2.21	110.80	113.94
2	B	1001	NAP	O3X-P2B-O2X	2.19	116.02	107.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	1001	NAP	O3X-P2B-O2X	2.19	116.02	107.80
2	D	1001	NAP	C6A-C5A-N7A	2.19	136.31	132.09
2	B	1001	NAP	C6A-C5A-N7A	2.18	136.30	132.09
2	C	1001	NAP	N9A-C8A-N7A	-2.17	110.85	113.94
2	C	1001	NAP	O3X-P2B-O2X	2.16	115.90	107.80
2	A	1001	NAP	C6A-C5A-N7A	2.14	136.22	132.09
2	D	1001	NAP	N9A-C8A-N7A	-2.13	110.91	113.94
2	A	1001	NAP	N9A-C8A-N7A	-2.09	110.97	113.94
2	A	1001	NAP	C2A-N1A-C6A	2.04	122.08	118.73
2	D	1001	NAP	C2A-N1A-C6A	2.03	122.07	118.73
3	C	1002	XOG	C9-C2-C3	-2.02	120.23	123.73
2	C	1001	NAP	C2A-N1A-C6A	2.02	122.05	118.73

There are no chirality outliers.

All (56) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1001	NAP	C5D-O5D-PN-O3
2	A	1001	NAP	C5D-O5D-PN-O1N
2	A	1001	NAP	C5D-O5D-PN-O2N
2	A	1001	NAP	O4D-C1D-N1N-C2N
2	A	1001	NAP	O4D-C1D-N1N-C6N
2	A	1001	NAP	C2D-C1D-N1N-C2N
2	A	1001	NAP	C2D-C1D-N1N-C6N
2	B	1001	NAP	C5D-O5D-PN-O3
2	B	1001	NAP	C5D-O5D-PN-O1N
2	B	1001	NAP	C5D-O5D-PN-O2N
2	B	1001	NAP	O4D-C1D-N1N-C2N
2	B	1001	NAP	O4D-C1D-N1N-C6N
2	B	1001	NAP	C2D-C1D-N1N-C2N
2	B	1001	NAP	C2D-C1D-N1N-C6N
2	C	1001	NAP	C5D-O5D-PN-O3
2	C	1001	NAP	C5D-O5D-PN-O1N
2	C	1001	NAP	C5D-O5D-PN-O2N
2	C	1001	NAP	O4D-C1D-N1N-C2N
2	C	1001	NAP	O4D-C1D-N1N-C6N
2	C	1001	NAP	C2D-C1D-N1N-C2N
2	C	1001	NAP	C2D-C1D-N1N-C6N
2	D	1001	NAP	C5B-O5B-PA-O2A
2	D	1001	NAP	C5D-O5D-PN-O3
2	D	1001	NAP	C5D-O5D-PN-O1N
2	D	1001	NAP	C5D-O5D-PN-O2N

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Mol	Chain	Res	Type	Atoms
2	D	1001	NAP	O4D-C1D-N1N-C2N
2	D	1001	NAP	O4D-C1D-N1N-C6N
2	D	1001	NAP	C2D-C1D-N1N-C2N
2	D	1001	NAP	C2D-C1D-N1N-C6N
3	A	1002	XOG	C6-C7-C8-O2
3	B	1002	XOG	C6-C7-C8-O2
3	C	1002	XOG	C6-C7-C8-O2
3	D	1002	XOG	C6-C7-C8-O2
3	B	1002	XOG	C4-C5-C6-C7
3	D	1002	XOG	C4-C5-C6-C7
3	A	1002	XOG	C4-C5-C6-C10
3	A	1002	XOG	C4-C5-C6-C7
3	B	1002	XOG	C4-C5-C6-C10
3	D	1002	XOG	C4-C5-C6-C10
2	C	1001	NAP	PN-O3-PA-O1A
2	D	1001	NAP	C5B-O5B-PA-O1A
2	D	1001	NAP	C5B-O5B-PA-O3
2	A	1001	NAP	C3D-C4D-C5D-O5D
2	D	1001	NAP	C2B-O2B-P2B-O3X
2	A	1001	NAP	O4D-C4D-C5D-O5D
2	D	1001	NAP	C3D-C4D-C5D-O5D
2	D	1001	NAP	PN-O3-PA-O1A
2	D	1001	NAP	O4D-C4D-C5D-O5D
3	C	1002	XOG	C4-C5-C6-C7
2	D	1001	NAP	O4B-C4B-C5B-O5B
2	B	1001	NAP	C2B-O2B-P2B-O3X
3	C	1002	XOG	C4-C5-C6-C10
2	C	1001	NAP	O4B-C4B-C5B-O5B
2	A	1001	NAP	PN-O3-PA-O2A
2	B	1001	NAP	PN-O3-PA-O2A
2	A	1001	NAP	O4B-C4B-C5B-O5B

There are no ring outliers.

11 monomers are involved in 27 short contacts:

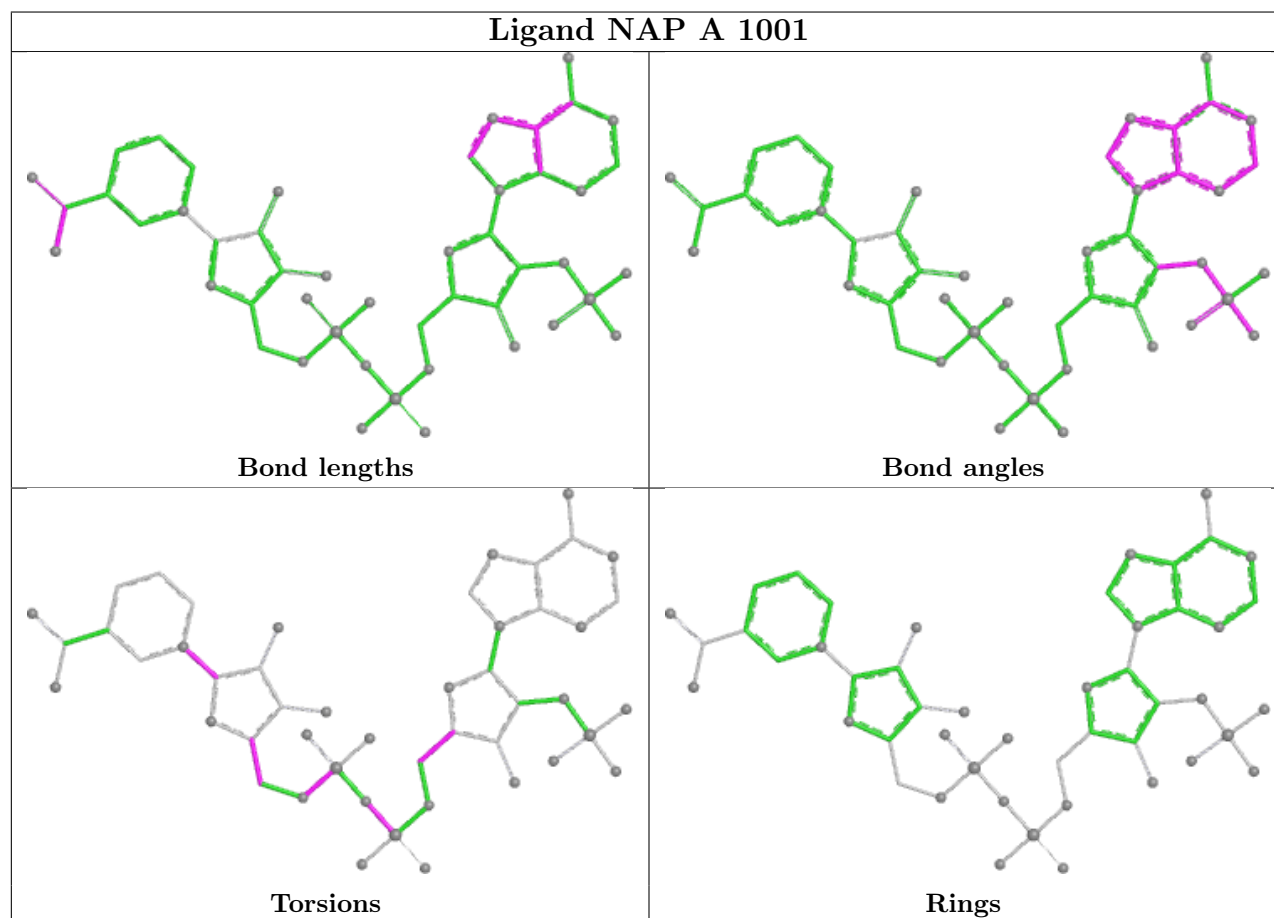
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	1002	XOG	3	0
2	A	1001	NAP	4	0
4	A	1003	PG4	2	0
2	D	1001	NAP	1	0
2	C	1001	NAP	3	0
3	D	1002	XOG	5	0

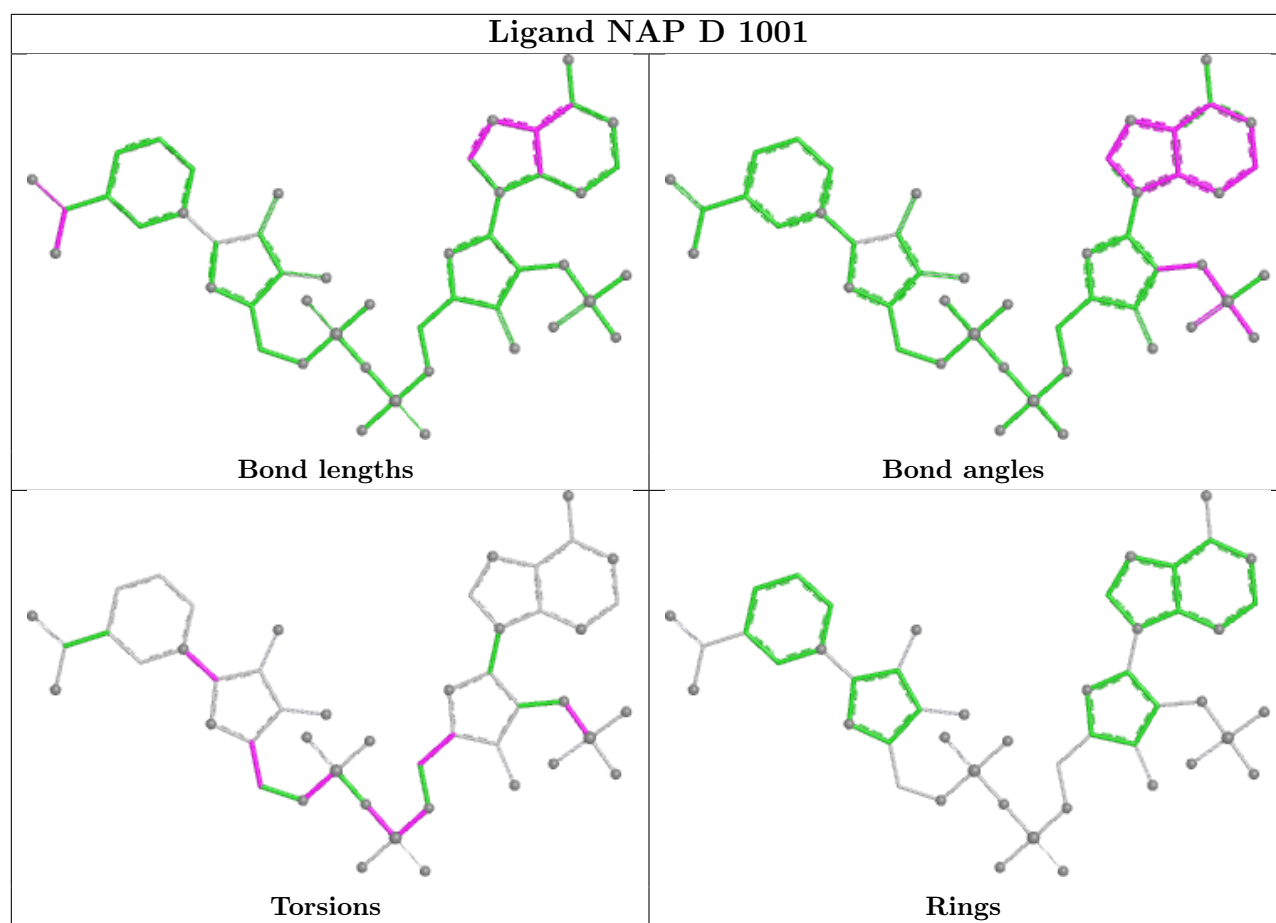
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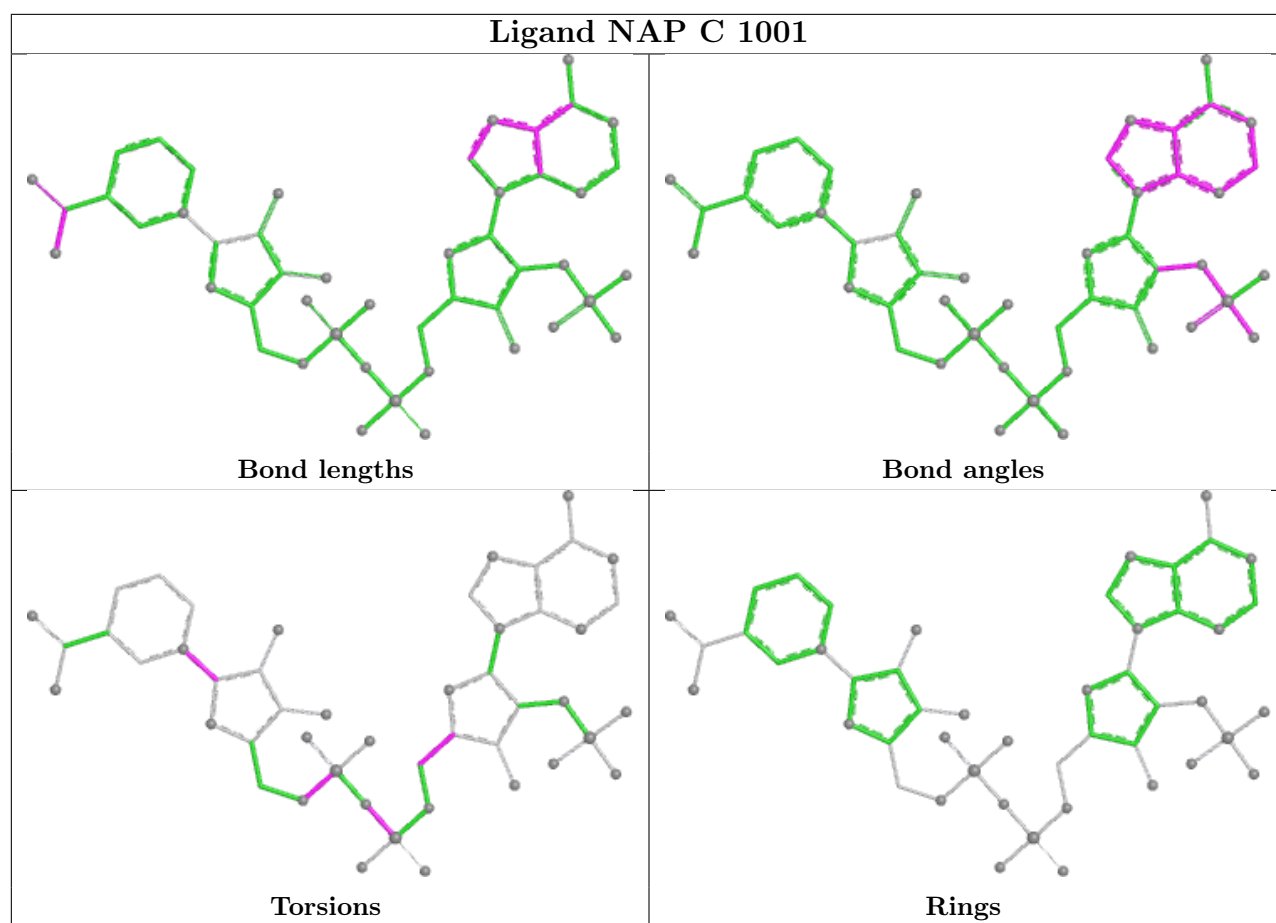
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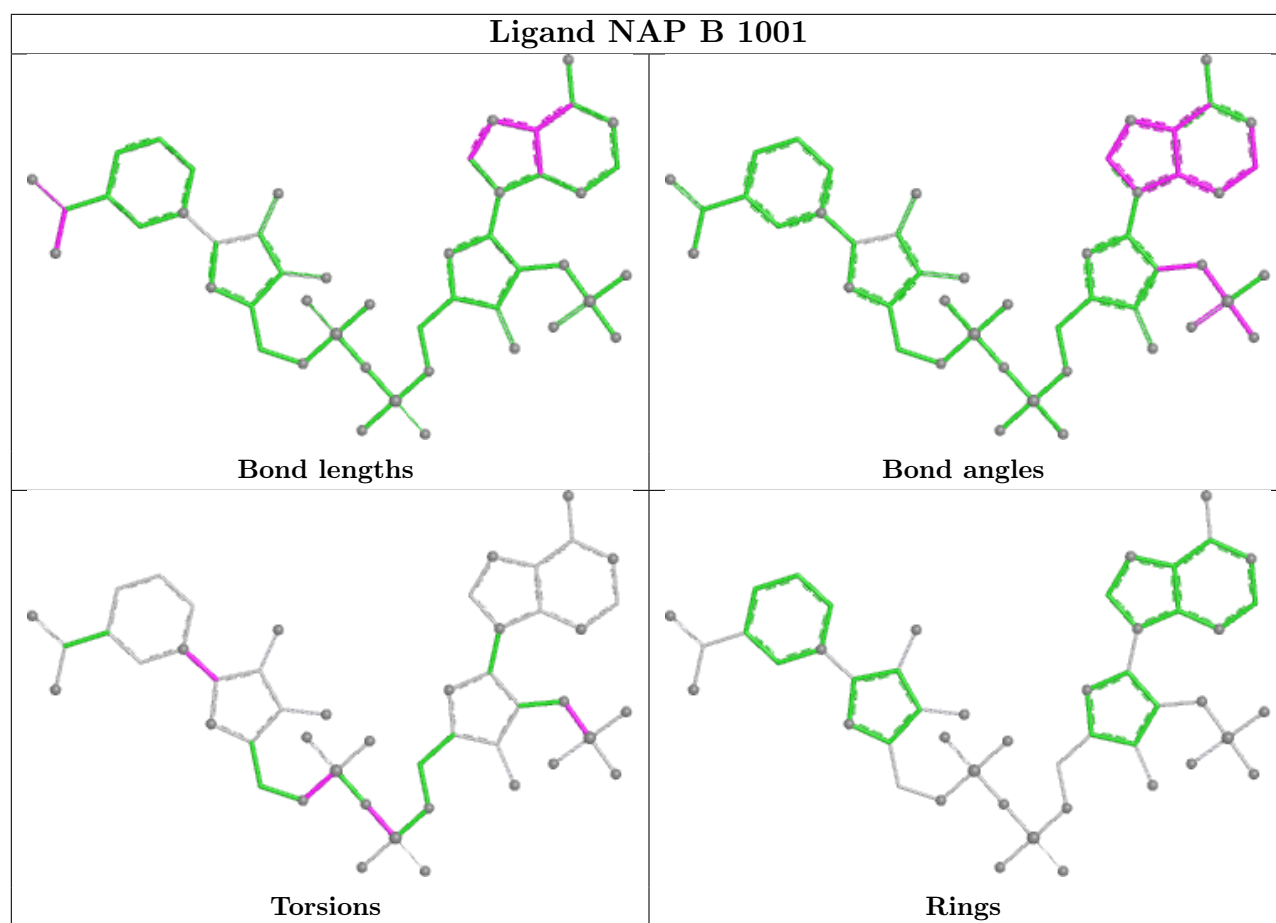
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	D	1003	PG4	1	0
3	A	1002	XOG	4	0
4	B	1003	PG4	1	0
3	B	1002	XOG	3	0
2	B	1001	NAP	2	0

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	365/365 (100%)	-1.60	0 100 100	35, 57, 94, 170	0
1	B	365/365 (100%)	-1.58	0 100 100	36, 57, 96, 176	0
1	C	365/365 (100%)	-1.55	0 100 100	38, 58, 96, 154	0
1	D	365/365 (100%)	-1.55	0 100 100	34, 58, 100, 177	0
All	All	1460/1460 (100%)	-1.57	0 100 100	34, 58, 98, 177	0

There are no RSRZ outliers to report.

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	XOG	A	1002	12/12	0.99	0.05	72,76,81,81	0
3	XOG	B	1002	12/12	0.99	0.05	75,77,79,81	0
3	XOG	C	1002	12/12	0.99	0.05	71,78,84,84	0
3	XOG	D	1002	12/12	0.99	0.06	78,80,89,90	0

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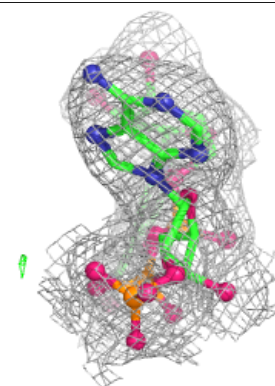
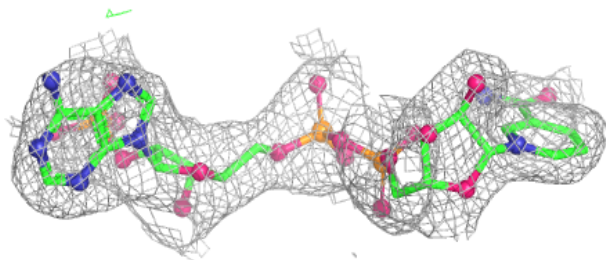
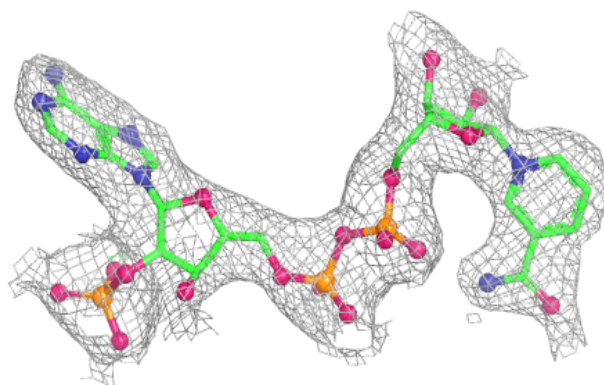
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	PG4	A	1003	13/13	0.99	0.05	77,80,83,83	0
4	PG4	B	1003	13/13	0.99	0.05	76,78,86,88	0
4	PG4	C	1003	13/13	0.99	0.04	68,74,81,83	0
4	PG4	D	1003	13/13	0.99	0.04	75,80,85,87	0
2	NAP	A	1001	48/48	1.00	0.02	27,54,61,83	0
2	NAP	B	1001	48/48	1.00	0.02	31,55,65,105	0
2	NAP	C	1001	48/48	1.00	0.02	32,57,62,84	0
2	NAP	D	1001	48/48	1.00	0.02	31,54,67,93	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.

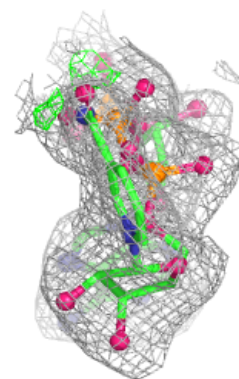
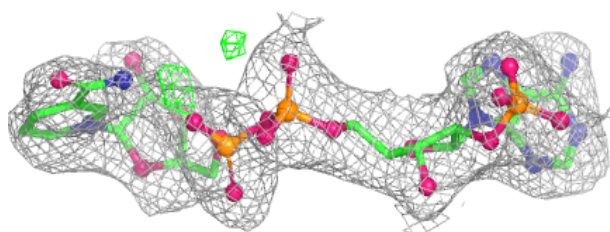
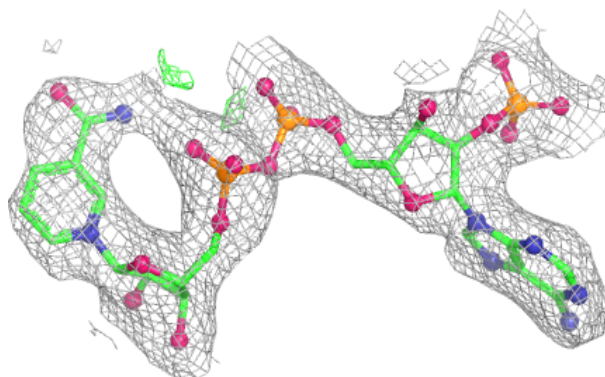
Electron density around NAP A 1001:

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

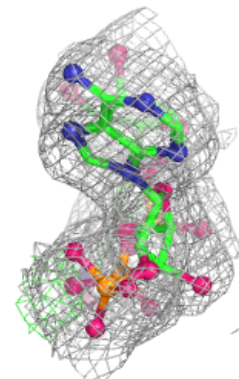
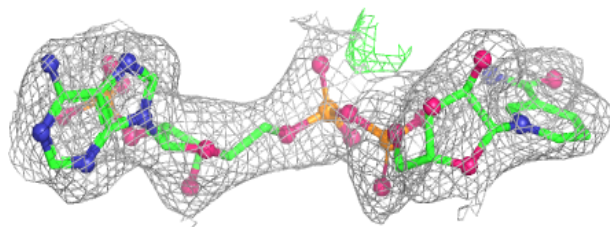
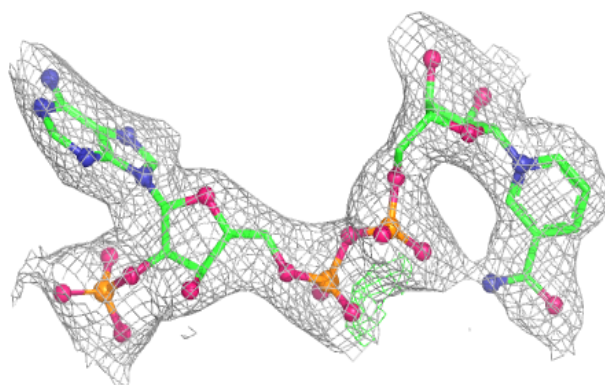


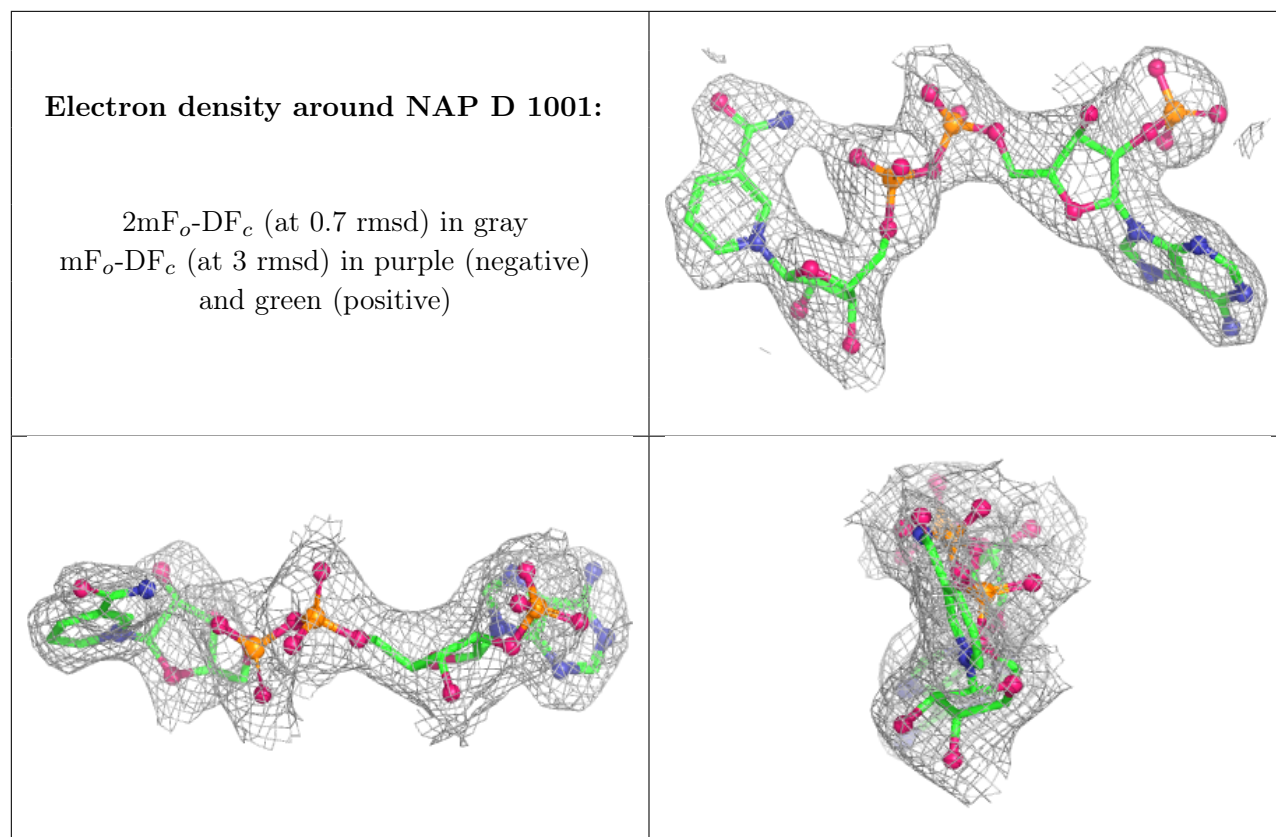
Electron density around NAP B 1001:

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

**Electron density around NAP C 1001:**

$2mF_o-DF_c$ (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.