



Full wwPDB X-ray Structure Validation Report ⓘ

May 17, 2020 – 03:33 am BST

PDB ID : 4CMC
Title : Crystal structure of pteridine reductase 1 (PTR1) from Trypanosoma brucei in ternary complex with cofactor and inhibitor
Authors : Barrack, K.L.; Hunter, W.N.
Deposited on : 2014-01-16
Resolution : 1.85 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

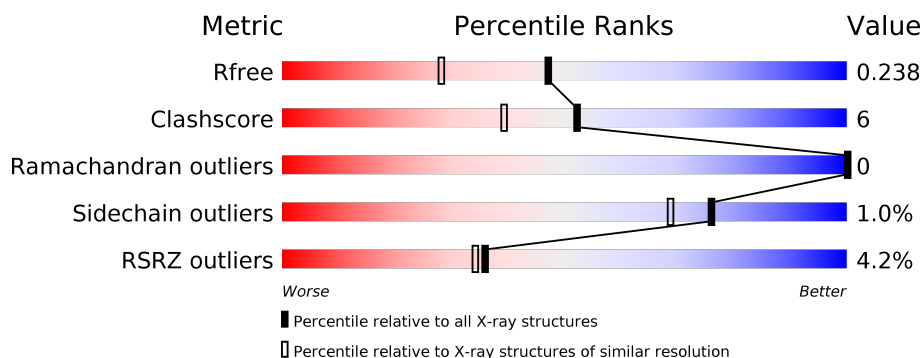
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.85 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}	130704	2469 (1.86-1.86)
Clashscore	141614	2625 (1.86-1.86)
Ramachandran outliers	138981	2592 (1.86-1.86)
Sidechain outliers	138945	2592 (1.86-1.86)
RSRZ outliers	127900	2436 (1.86-1.86)

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 on 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	288	<div> <div>6%</div> <div> <div></div> <div>76%</div> <div>11%</div> <div>13%</div> </div> </div>
1	B	288	<div> <div>2%</div> <div> <div></div> <div>75%</div> <div>11%</div> <div>14%</div> </div> </div>
1	C	288	<div> <div>4%</div> <div> <div></div> <div>80%</div> <div>7%</div> <div>14%</div> </div> </div>
1	D	288	<div> <div>2%</div> <div> <div></div> <div>77%</div> <div>10%</div> <div>13%</div> </div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 8592 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 PTERIDINE REDUCTASE 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	251	Total	C	N	O	S	0	3	0
			1886	1183	333	359	11			
1	B	249	Total	C	N	O	S	0	2	0
			1870	1175	329	355	11			
1	C	249	Total	C	N	O	S	0	2	0
			1868	1174	329	354	11			
1	D	251	Total	C	N	O	S	0	4	0
			1896	1190	335	360	11			

There are 80 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	expression tag	UNP O76290
A	-18	GLY	-	expression tag	UNP O76290
A	-17	SER	-	expression tag	UNP O76290
A	-16	SER	-	expression tag	UNP O76290
A	-15	HIS	-	expression tag	UNP O76290
A	-14	HIS	-	expression tag	UNP O76290
A	-13	HIS	-	expression tag	UNP O76290
A	-12	HIS	-	expression tag	UNP O76290
A	-11	HIS	-	expression tag	UNP O76290
A	-10	HIS	-	expression tag	UNP O76290
A	-9	SER	-	expression tag	UNP O76290
A	-8	SER	-	expression tag	UNP O76290
A	-7	GLY	-	expression tag	UNP O76290
A	-6	LEU	-	expression tag	UNP O76290
A	-5	VAL	-	expression tag	UNP O76290
A	-4	PRO	-	expression tag	UNP O76290
A	-3	ARG	-	expression tag	UNP O76290
A	-2	GLY	-	expression tag	UNP O76290
A	-1	SER	-	expression tag	UNP O76290
A	0	HIS	-	expression tag	UNP O76290
B	-19	MET	-	expression tag	UNP O76290

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-18	GLY	-	expression tag	UNP O76290
B	-17	SER	-	expression tag	UNP O76290
B	-16	SER	-	expression tag	UNP O76290
B	-15	HIS	-	expression tag	UNP O76290
B	-14	HIS	-	expression tag	UNP O76290
B	-13	HIS	-	expression tag	UNP O76290
B	-12	HIS	-	expression tag	UNP O76290
B	-11	HIS	-	expression tag	UNP O76290
B	-10	HIS	-	expression tag	UNP O76290
B	-9	SER	-	expression tag	UNP O76290
B	-8	SER	-	expression tag	UNP O76290
B	-7	GLY	-	expression tag	UNP O76290
B	-6	LEU	-	expression tag	UNP O76290
B	-5	VAL	-	expression tag	UNP O76290
B	-4	PRO	-	expression tag	UNP O76290
B	-3	ARG	-	expression tag	UNP O76290
B	-2	GLY	-	expression tag	UNP O76290
B	-1	SER	-	expression tag	UNP O76290
B	0	HIS	-	expression tag	UNP O76290
C	-19	MET	-	expression tag	UNP O76290
C	-18	GLY	-	expression tag	UNP O76290
C	-17	SER	-	expression tag	UNP O76290
C	-16	SER	-	expression tag	UNP O76290
C	-15	HIS	-	expression tag	UNP O76290
C	-14	HIS	-	expression tag	UNP O76290
C	-13	HIS	-	expression tag	UNP O76290
C	-12	HIS	-	expression tag	UNP O76290
C	-11	HIS	-	expression tag	UNP O76290
C	-10	HIS	-	expression tag	UNP O76290
C	-9	SER	-	expression tag	UNP O76290
C	-8	SER	-	expression tag	UNP O76290
C	-7	GLY	-	expression tag	UNP O76290
C	-6	LEU	-	expression tag	UNP O76290
C	-5	VAL	-	expression tag	UNP O76290
C	-4	PRO	-	expression tag	UNP O76290
C	-3	ARG	-	expression tag	UNP O76290
C	-2	GLY	-	expression tag	UNP O76290
C	-1	SER	-	expression tag	UNP O76290
C	0	HIS	-	expression tag	UNP O76290
D	-19	MET	-	expression tag	UNP O76290
D	-18	GLY	-	expression tag	UNP O76290
D	-17	SER	-	expression tag	UNP O76290

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-16	SER	-	expression tag	UNP O76290
D	-15	HIS	-	expression tag	UNP O76290
D	-14	HIS	-	expression tag	UNP O76290
D	-13	HIS	-	expression tag	UNP O76290
D	-12	HIS	-	expression tag	UNP O76290
D	-11	HIS	-	expression tag	UNP O76290
D	-10	HIS	-	expression tag	UNP O76290
D	-9	SER	-	expression tag	UNP O76290
D	-8	SER	-	expression tag	UNP O76290
D	-7	GLY	-	expression tag	UNP O76290
D	-6	LEU	-	expression tag	UNP O76290
D	-5	VAL	-	expression tag	UNP O76290
D	-4	PRO	-	expression tag	UNP O76290
D	-3	ARG	-	expression tag	UNP O76290
D	-2	GLY	-	expression tag	UNP O76290
D	-1	SER	-	expression tag	UNP O76290
D	0	HIS	-	expression tag	UNP O76290

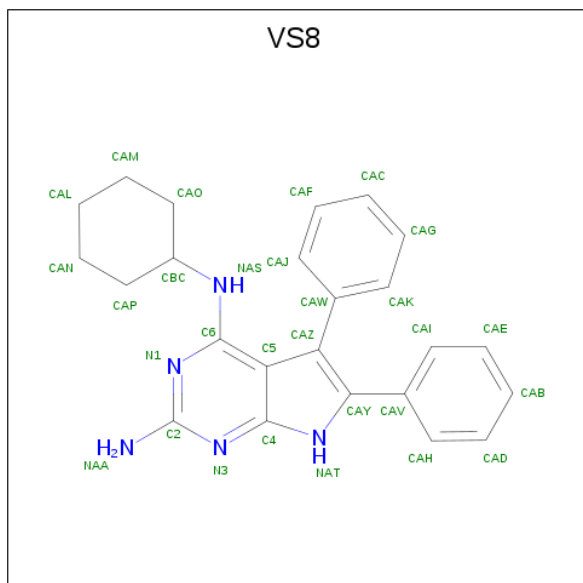
- # NAP
-
- The chemical structure of Naproxen (NAP) is shown, featuring a naphthalene ring system. The structure includes a carboxylic acid group (COOH) and a chiral center (C1) with a hydroxyl group (OH) and a naphthalene ring. The naphthalene ring is substituted with a carboxylic acid group (COOH) and a chiral center (C1) with a hydroxyl group (OH) and a naphthalene ring. The structure is labeled with various atoms and groups, including NH, COOH, OH, and COO⁻.

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

- Molecule 3 is N4-cyclohexyl-5,6-diphenyl-7H-pyrrolo[2,3-d]pyrimidine-2,4-diamine (three-letter code: VS8) (formula: C₂₄H₂₅N₅).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	N	0	0
			29	24	5		
3	B	1	Total	C	N	0	0
			29	24	5		
3	C	1	Total	C	N	0	0
			29	24	5		
3	D	1	Total	C	N	0	0
			29	24	5		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	217	Total	O	0	0
			217	217		
4	B	201	Total	O	0	0
			201	201		
4	C	189	Total	O	0	0
			189	189		

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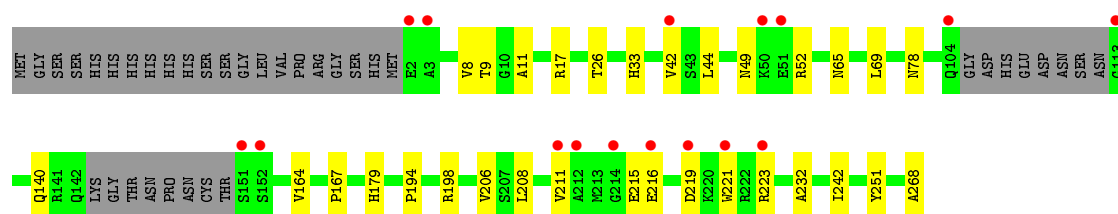
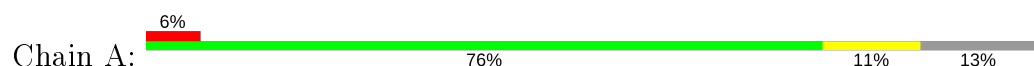
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	157	Total	O	0	0
			157	157		

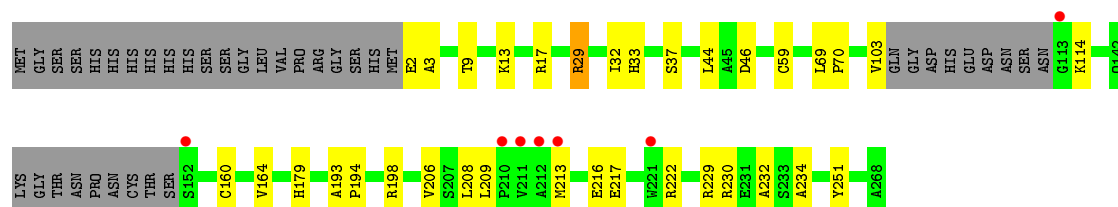
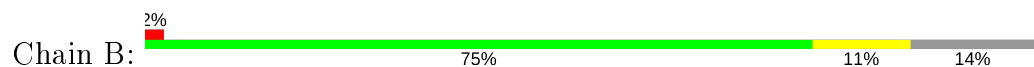
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

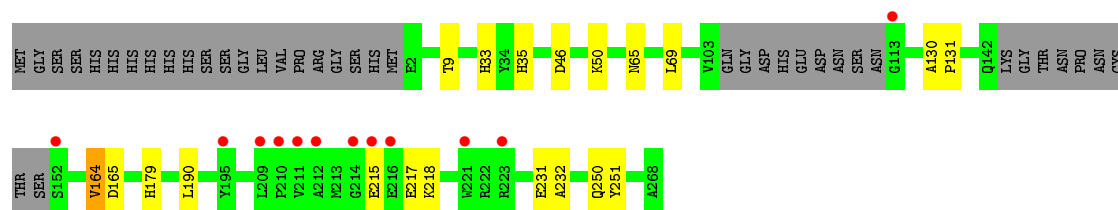
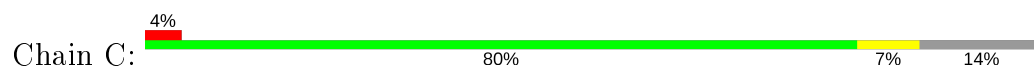
• Molecule 1: PTERIDINE REDUCTASE 1



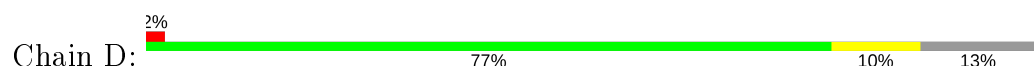
• Molecule 1: PTERIDINE REDUCTASE 1



• Molecule 1: PTERIDINE REDUCTASE 1



• Molecule 1: PTERIDINE REDUCTASE 1





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	74.19Å 89.47Å 84.39Å 90.00° 115.55° 90.00°	Depositor
Resolution (Å)	38.10 – 1.85 38.07 – 1.85	Depositor EDS
% Data completeness (in resolution range)	97.6 (38.10-1.85) 97.6 (38.07-1.85)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.04 (at 1.85Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.184 , 0.230 0.195 , 0.238	Depositor DCC
R_{free} test set	4177 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	16.6	Xtriage
Anisotropy	0.054	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 44.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.017 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8592	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 27.18 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 2.3017e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: VS8, 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	0.67	0/1913	0.76	1/2595 (0.0%)
1	B	0.68	0/1897	0.80	5/2574 (0.2%)
1	C	0.66	0/1895	0.79	1/2571 (0.0%)
1	D	0.66	0/1923	0.80	2/2609 (0.1%)
All	All	0.67	0/7628	0.79	9/10349 (0.1%)

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	159	LEU	CA-CB-CG	-8.15	96.55	115.30
1	B	198	ARG	NE-CZ-NH1	7.38	123.99	120.30
1	B	29	ARG	NE-CZ-NH2	-7.04	116.78	120.30
1	B	29	ARG	NE-CZ-NH1	6.01	123.31	120.30
1	B	198	ARG	NE-CZ-NH2	-5.97	117.31	120.30
1	B	46	ASP	CB-CG-OD1	5.62	123.35	118.30
1	D	82	ARG	NE-CZ-NH2	-5.21	117.70	120.30
1	A	198	ARG	NE-CZ-NH1	5.08	122.84	120.30
1	C	165	ASP	CB-CG-OD1	5.07	122.87	118.30

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	1886	0	1905	21	0
1	B	1870	0	1890	27	0
1	C	1868	0	1891	14	0
1	D	1896	0	1918	20	0
2	A	48	0	25	0	0
2	B	48	0	25	0	0
2	C	48	0	25	1	0
2	D	48	0	25	2	0
3	A	29	0	25	2	0
3	B	29	0	25	2	0
3	C	29	0	25	2	0
3	D	29	0	25	3	0
4	A	217	0	0	8	2
4	B	201	0	0	9	2
4	C	189	0	0	1	0
4	D	157	0	0	6	0
All	All	8592	0	7804	86	2

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 (86) 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:B:3:ALA:HB3	1:B:29:ARG:HD2	1.69	0.72
1:A:78:ASN:HB3	4:A:2093:HOH:O	1.90	0.71
1:A:140:GLN:HG3	4:A:2133:HOH:O	1.95	0.65
1:A:42:VAL:HG23	4:A:2041:HOH:O	1.97	0.65
1:A:219:ASP:OD2	1:A:223:ARG:NH1	2.30	0.64
1:B:213:MET:HG2	1:B:217[A]:GLU:HG3	1.83	0.61
1:C:215:GLU:HG3	1:C:218:LYS:CE	2.30	0.60
1:A:164:VAL:HG22	1:A:179:HIS:CD2	2.37	0.60
1:D:219:ASP:OD2	1:D:223:ARG:HD3	2.02	0.59
1:B:213:MET:HE3	4:B:2166:HOH:O	2.00	0.59
1:B:213:MET:CG	1:B:217[A]:GLU:HG3	2.34	0.58
1:C:215:GLU:HG3	1:C:218:LYS:HE2	1.86	0.57
4:B:2114:HOH:O	1:D:140:GLN:CG	2.51	0.57
1:B:222:ARG:HB3	1:B:229:ARG:O	2.04	0.57
1:A:167:PRO:HG2	1:C:190[B]:LEU:HD12	1.87	0.56
1:D:138:PHE:O	1:D:142:GLN:HG2	2.05	0.56
3:B:1270:VS8:CAK	3:B:1270:VS8:HAS	2.21	0.53
1:D:164:VAL:HG22	1:D:179:HIS:CD2	2.43	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:103:VAL:HG22	4:B:2113:HOH:O	2.08	0.53
1:B:13:LYS:HE2	1:B:37:SER:OG	2.09	0.53
1:A:140:GLN:CG	4:A:2133:HOH:O	2.55	0.52
1:D:46:ASP:HB3	1:D:50:LYS:HE3	1.91	0.52
4:B:2113:HOH:O	1:D:191:GLU:OE2	2.19	0.52
1:A:17:ARG:HG3	1:A:44:LEU:HD22	1.91	0.52
1:D:112:ASN:C	1:D:114:LYS:H	2.14	0.52
1:D:13:LYS:HE3	4:D:2008:HOH:O	2.09	0.51
3:D:1270:VS8:HAS	3:D:1270:VS8:CAJ	2.23	0.51
1:C:215:GLU:CG	1:C:218:LYS:HE3	2.41	0.51
1:A:232:ALA:HB2	1:B:251:TYR:CE2	2.46	0.51
1:D:26:THR:O	1:D:26:THR:HG22	2.12	0.50
1:D:114:LYS:HE2	4:D:2098:HOH:O	2.12	0.50
3:A:1270:VS8:HAS	3:A:1270:VS8:CAJ	2.26	0.49
1:B:206[B]:VAL:CG2	1:B:209:LEU:HD21	2.42	0.49
4:B:2114:HOH:O	1:D:140:GLN:HG3	2.13	0.49
1:C:9:THR:HA	1:C:33:HIS:HB3	1.95	0.48
1:A:42:VAL:CG2	4:A:2041:HOH:O	2.59	0.48
3:C:1270:VS8:HAS	3:C:1270:VS8:CAJ	2.26	0.48
1:D:115:THR:HG21	4:D:2100:HOH:O	2.14	0.47
1:D:65:ASN:HA	1:D:69:LEU:HD22	1.96	0.47
1:B:160:CYS:HB3	4:B:2138:HOH:O	2.15	0.46
1:A:215:GLU:HB3	4:A:2173:HOH:O	2.14	0.46
1:B:213:MET:HE2	1:B:217[A]:GLU:HG3	1.96	0.46
1:A:65:ASN:HA	1:A:69:LEU:HD22	1.97	0.46
1:B:213:MET:CE	1:B:217[B]:GLU:HG3	2.46	0.45
1:B:213:MET:HE2	1:B:217[A]:GLU:C	2.37	0.45
1:C:231:GLU:OE2	4:C:2144:HOH:O	2.21	0.45
1:B:213:MET:HE2	1:B:217[B]:GLU:CG	2.46	0.45
3:A:1270:VS8:CAH	3:A:1270:VS8:CAK	2.94	0.45
1:A:49:ASN:HA	1:A:52:ARG:O	2.16	0.45
1:A:9:THR:HA	1:A:33:HIS:HB3	1.97	0.45
1:C:215:GLU:HG3	1:C:218:LYS:HE3	1.96	0.45
1:A:268:ALA:HB3	4:A:2215:HOH:O	2.15	0.45
3:B:1270:VS8:CAB	4:B:2139:HOH:O	2.64	0.45
1:C:130:ALA:HB3	1:C:131:PRO:HD3	1.98	0.44
3:D:1270:VS8:CAB	4:D:2113:HOH:O	2.66	0.44
1:B:69:LEU:N	1:B:70:PRO:CD	2.81	0.44
1:C:65:ASN:HA	1:C:69:LEU:HD22	1.99	0.43
1:C:35:HIS:HB2	2:C:1269:NAP:C2A	2.48	0.43
1:A:251:TYR:CE2	1:B:232:ALA:HB2	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:1270:VS8:CAK	3:D:1270:VS8:CAI	2.96	0.43
1:D:214:GLY:HA3	4:D:2131:HOH:O	2.18	0.43
1:B:208:LEU:HD22	1:B:234:ALA:HB2	2.00	0.43
1:B:114:LYS:HE2	4:B:2117:HOH:O	2.18	0.43
3:C:1270:VS8:CAH	3:C:1270:VS8:CAK	2.97	0.42
1:D:17:ARG:HG3	1:D:44:LEU:HD22	2.02	0.42
1:B:9:THR:HA	1:B:33:HIS:HB3	2.01	0.42
1:C:251:TYR:CE2	1:D:232:ALA:HB2	2.55	0.42
1:C:164:VAL:HG22	1:C:179:HIS:CD2	2.54	0.42
1:D:15:ILE:HB	2:D:1269:NAP:H51N	2.01	0.42
1:B:206[B]:VAL:HG21	1:B:209:LEU:HD21	2.02	0.41
1:B:2:GLU:HA	4:B:2002:HOH:O	2.19	0.41
1:A:8:VAL:HG12	1:A:11:ALA:HB2	2.02	0.41
1:A:26:THR:O	1:A:26:THR:HG22	2.21	0.41
1:B:193:ALA:N	1:B:194:PRO:CD	2.83	0.41
1:B:33:HIS:HA	1:B:59:CYS:O	2.21	0.41
1:A:208:LEU:HD12	1:A:208:LEU:HA	1.85	0.41
1:B:32:ILE:HD12	1:B:44:LEU:HD23	2.03	0.41
1:B:17:ARG:HG3	1:B:44:LEU:HD22	2.02	0.41
1:C:232:ALA:HB2	1:D:251:TYR:CE2	2.56	0.40
1:C:46:ASP:O	1:C:50:LYS:HD3	2.20	0.40
1:A:194:PRO:HG3	4:A:2167:HOH:O	2.21	0.40
1:D:230:ARG:HD2	4:D:2143:HOH:O	2.21	0.40
1:D:35:HIS:HB2	2:D:1269:NAP:C2A	2.51	0.40
1:B:164:VAL:HG22	1:B:179:HIS:CD2	2.56	0.40
1:B:222:ARG:HG2	1:B:230:ARG:HA	2.03	0.40
1:A:206[B]:VAL:HG21	1:A:221:TRP:CE3	2.56	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:2011:HOH:O	4:B:2167:HOH:O[2_555]	2.17	0.03
4:A:2177:HOH:O	4:B:2178:HOH:O[2_555]	2.18	0.02

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	248/288 (86%)	240 (97%)	8 (3%)	0	100	100
1	B	245/288 (85%)	236 (96%)	9 (4%)	0	100	100
1	C	245/288 (85%)	234 (96%)	11 (4%)	0	100	100
1	D	249/288 (86%)	237 (95%)	12 (5%)	0	100	100
All	All	987/1152 (86%)	947 (96%)	40 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	201/231 (87%)	198 (98%)	3 (2%)	65	53
1	B	199/231 (86%)	198 (100%)	1 (0%)	88	86
1	C	199/231 (86%)	196 (98%)	3 (2%)	65	53
1	D	202/231 (87%)	201 (100%)	1 (0%)	88	86
All	All	801/924 (87%)	793 (99%)	8 (1%)	76	69

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	211	VAL
1	A	216	GLU

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Mol	Chain	Res	Type
1	A	242	ILE
1	B	216	GLU
1	C	164	VAL
1	C	217	GLU
1	C	250	GLN
1	D	152	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	25	GLN

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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

8 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	VS8	C	1270	-	32,33,33	2.74	5 (15%)	36,46,46	1.97	8 (22%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	VS8	A	1270	-	32,33,33	3.09	8 (25%)	36,46,46	1.78	7 (19%)
2	NAP	D	1269	-	45,52,52	1.31	3 (6%)	56,80,80	1.37	9 (16%)
2	NAP	B	1269	-	45,52,52	1.06	2 (4%)	56,80,80	1.35	8 (14%)
3	VS8	D	1270	-	32,33,33	2.71	7 (21%)	36,46,46	1.72	6 (16%)
2	NAP	C	1269	-	45,52,52	0.96	1 (2%)	56,80,80	1.31	6 (10%)
3	VS8	B	1270	-	32,33,33	2.91	7 (21%)	36,46,46	1.99	10 (27%)
2	NAP	A	1269	-	45,52,52	1.08	1 (2%)	56,80,80	1.33	9 (16%)

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	VS8	C	1270	-	-	0/12/20/20	0/5/5/5
3	VS8	A	1270	-	-	0/12/20/20	0/5/5/5
2	NAP	D	1269	-	-	0/31/67/67	0/5/5/5
2	NAP	B	1269	-	-	1/31/67/67	0/5/5/5
3	VS8	D	1270	-	-	0/12/20/20	0/5/5/5
2	NAP	C	1269	-	-	0/31/67/67	0/5/5/5
3	VS8	B	1270	-	-	0/12/20/20	0/5/5/5
2	NAP	A	1269	-	-	0/31/67/67	0/5/5/5

All (34) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1270	VS8	CAZ-CAW	-11.62	1.35	1.49
3	B	1270	VS8	CAV-CAY	-10.31	1.37	1.49
3	C	1270	VS8	CAZ-CAW	-10.01	1.37	1.49
3	D	1270	VS8	CAZ-CAW	-9.71	1.38	1.49
3	B	1270	VS8	CAZ-CAW	-8.99	1.38	1.49
3	D	1270	VS8	CAV-CAY	-8.74	1.39	1.49
3	A	1270	VS8	CAV-CAY	-8.50	1.39	1.49
3	C	1270	VS8	CAV-CAY	-8.28	1.39	1.49
3	C	1270	VS8	C6-C5	-5.84	1.36	1.44
3	B	1270	VS8	C6-C5	-5.75	1.37	1.44
3	D	1270	VS8	C6-C5	-4.83	1.38	1.44
3	A	1270	VS8	C6-C5	-4.62	1.38	1.44
3	A	1270	VS8	CAZ-CAY	-4.52	1.36	1.41
2	D	1269	NAP	C7N-N7N	-4.41	1.24	1.33
2	A	1269	NAP	P2B-O2B	4.31	1.67	1.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	1269	NAP	P2B-O2B	3.75	1.66	1.59
3	A	1270	VS8	CAZ-C5	-3.53	1.37	1.42
2	D	1269	NAP	O7N-C7N	3.51	1.30	1.24
3	B	1270	VS8	CAZ-CAY	-3.33	1.37	1.41
3	B	1270	VS8	CAZ-C5	-3.14	1.38	1.42
2	D	1269	NAP	P2B-O2B	3.00	1.65	1.59
2	B	1269	NAP	P2B-O2B	2.99	1.65	1.59
3	D	1270	VS8	CAZ-C5	-2.90	1.38	1.42
3	A	1270	VS8	CAO-CBC	2.79	1.58	1.52
3	A	1270	VS8	C6-N1	2.68	1.36	1.32
3	C	1270	VS8	CAZ-C5	-2.64	1.38	1.42
3	C	1270	VS8	CAZ-CAY	-2.60	1.38	1.41
2	B	1269	NAP	C5A-C4A	2.44	1.47	1.40
3	D	1270	VS8	CAZ-CAY	-2.44	1.38	1.41
3	D	1270	VS8	CAO-CBC	2.36	1.57	1.52
3	B	1270	VS8	CAO-CBC	2.35	1.57	1.52
3	D	1270	VS8	C6-N1	2.21	1.35	1.32
3	A	1270	VS8	CBC-NAS	2.20	1.51	1.46
3	B	1270	VS8	C6-N1	2.16	1.35	1.32

All (63) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1270	VS8	N3-C2-N1	-6.79	118.17	127.22
3	C	1270	VS8	N3-C2-N1	-6.34	118.77	127.22
3	D	1270	VS8	N3-C2-N1	-5.97	119.26	127.22
3	A	1270	VS8	N3-C2-N1	-5.01	120.53	127.22
3	A	1270	VS8	CAM-CAO-CBC	-4.66	102.34	111.11
2	B	1269	NAP	C3N-C7N-N7N	4.21	122.81	117.75
2	A	1269	NAP	C1B-N9A-C4A	-4.16	119.34	126.64
3	C	1270	VS8	CAH-CAV-CAY	3.80	126.62	120.61
3	B	1270	VS8	C2-N3-C4	3.78	119.67	115.36
3	D	1270	VS8	C2-N3-C4	3.77	119.66	115.36
3	C	1270	VS8	C2-N1-C6	3.71	122.85	114.54
3	B	1270	VS8	C2-N1-C6	3.69	122.80	114.54
2	C	1269	NAP	N3A-C2A-N1A	-3.60	123.06	128.68
2	D	1269	NAP	N3A-C2A-N1A	-3.46	123.28	128.68
2	B	1269	NAP	O7N-C7N-C3N	-3.35	115.62	119.63
2	A	1269	NAP	N3A-C2A-N1A	-3.35	123.45	128.68
2	C	1269	NAP	O2B-P2B-O1X	-3.30	96.67	109.39
3	B	1270	VS8	CAH-CAV-CAY	3.28	125.80	120.61
3	D	1270	VS8	C2-N1-C6	3.19	121.69	114.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	1270	VS8	CAE-CAI-CAV	3.18	124.55	120.56
2	C	1269	NAP	C1B-N9A-C4A	-3.15	121.10	126.64
2	C	1269	NAP	O3X-P2B-O2X	3.13	119.62	107.64
3	C	1270	VS8	C2-N3-C4	3.03	118.81	115.36
2	D	1269	NAP	C3N-C7N-N7N	3.02	121.38	117.75
3	D	1270	VS8	NAA-C2-N3	2.98	122.65	117.79
3	A	1270	VS8	C2-N1-C6	2.96	121.17	114.54
3	A	1270	VS8	C2-N3-C4	2.95	118.73	115.36
2	D	1269	NAP	C2A-N1A-C6A	2.94	123.78	118.75
2	A	1269	NAP	C3N-C7N-N7N	2.91	121.25	117.75
2	B	1269	NAP	C1B-N9A-C4A	-2.89	121.57	126.64
3	B	1270	VS8	CAM-CAO-CBC	-2.88	105.69	111.11
2	B	1269	NAP	N3A-C2A-N1A	-2.88	124.18	128.68
2	C	1269	NAP	O2N-PN-O1N	2.75	125.82	112.24
2	D	1269	NAP	C6N-N1N-C2N	-2.69	119.52	121.97
3	A	1270	VS8	CAH-CAV-CAY	2.61	124.74	120.61
2	D	1269	NAP	O3X-P2B-O2X	2.61	117.59	107.64
2	D	1269	NAP	C4A-C5A-N7A	-2.59	106.70	109.40
3	B	1270	VS8	NAA-C2-N3	2.59	122.00	117.79
2	D	1269	NAP	C5A-C6A-N6A	2.52	124.19	120.35
3	C	1270	VS8	C5-C6-NAS	2.47	122.98	120.66
3	B	1270	VS8	CAI-CAV-CAY	-2.46	116.70	120.61
3	A	1270	VS8	CAW-CAZ-CAY	-2.42	122.08	126.88
2	A	1269	NAP	O2X-P2B-O2B	-2.34	95.50	105.99
2	B	1269	NAP	O2N-PN-O1N	2.31	123.65	112.24
3	D	1270	VS8	CAI-CAV-CAY	2.29	124.23	120.61
2	B	1269	NAP	C2A-N1A-C6A	2.28	122.65	118.75
2	A	1269	NAP	O4D-C1D-C2D	-2.24	103.65	106.93
2	A	1269	NAP	O3X-P2B-O2X	2.22	116.11	107.64
2	C	1269	NAP	O2A-PA-O1A	2.18	123.01	112.24
3	B	1270	VS8	CAG-CAK-CAW	2.18	123.30	120.56
2	A	1269	NAP	O7N-C7N-C3N	-2.17	117.04	119.63
2	D	1269	NAP	O2N-PN-O1N	2.16	122.91	112.24
3	C	1270	VS8	CAM-CAO-CBC	-2.14	107.08	111.11
3	C	1270	VS8	NAA-C2-N3	2.14	121.27	117.79
3	B	1270	VS8	CAE-CAI-CAV	2.12	123.22	120.56
2	A	1269	NAP	C2A-N1A-C6A	2.10	122.34	118.75
3	D	1270	VS8	CAN-CAP-CBC	-2.05	107.25	111.11
2	D	1269	NAP	C1B-N9A-C4A	-2.04	123.05	126.64
3	B	1270	VS8	CAP-CBC-CAO	-2.04	107.28	110.82
3	A	1270	VS8	NAA-C2-N1	2.01	120.38	117.25
2	B	1269	NAP	N6A-C6A-N1A	2.01	122.74	118.57

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1269	NAP	O2A-PA-O1A	2.01	122.16	112.24
2	A	1269	NAP	O3X-P2B-O1X	2.00	118.52	110.68

There are no chirality outliers.

All (1) torsion outliers are listed below:

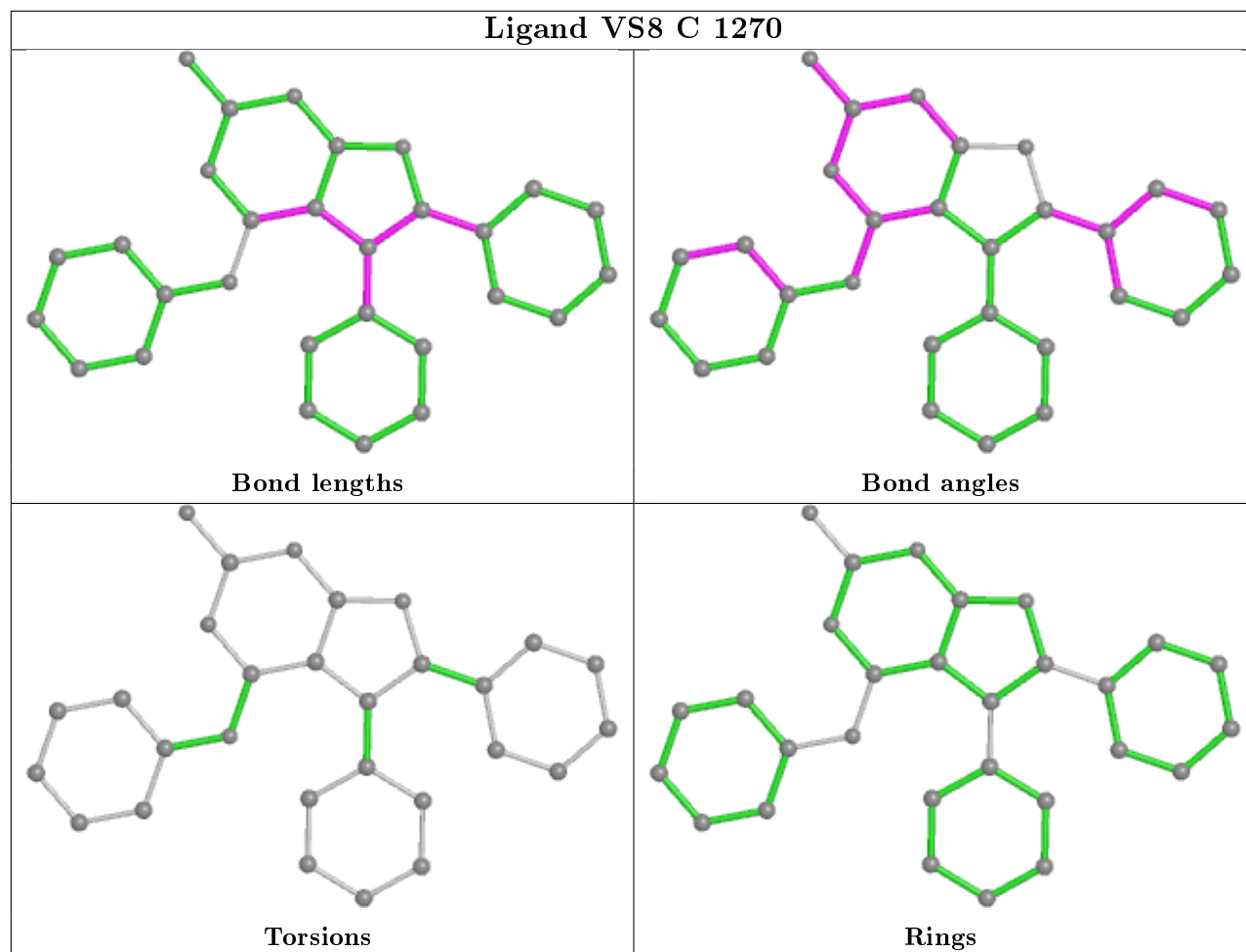
Mol	Chain	Res	Type	Atoms
2	B	1269	NAP	C3B-C2B-O2B-P2B

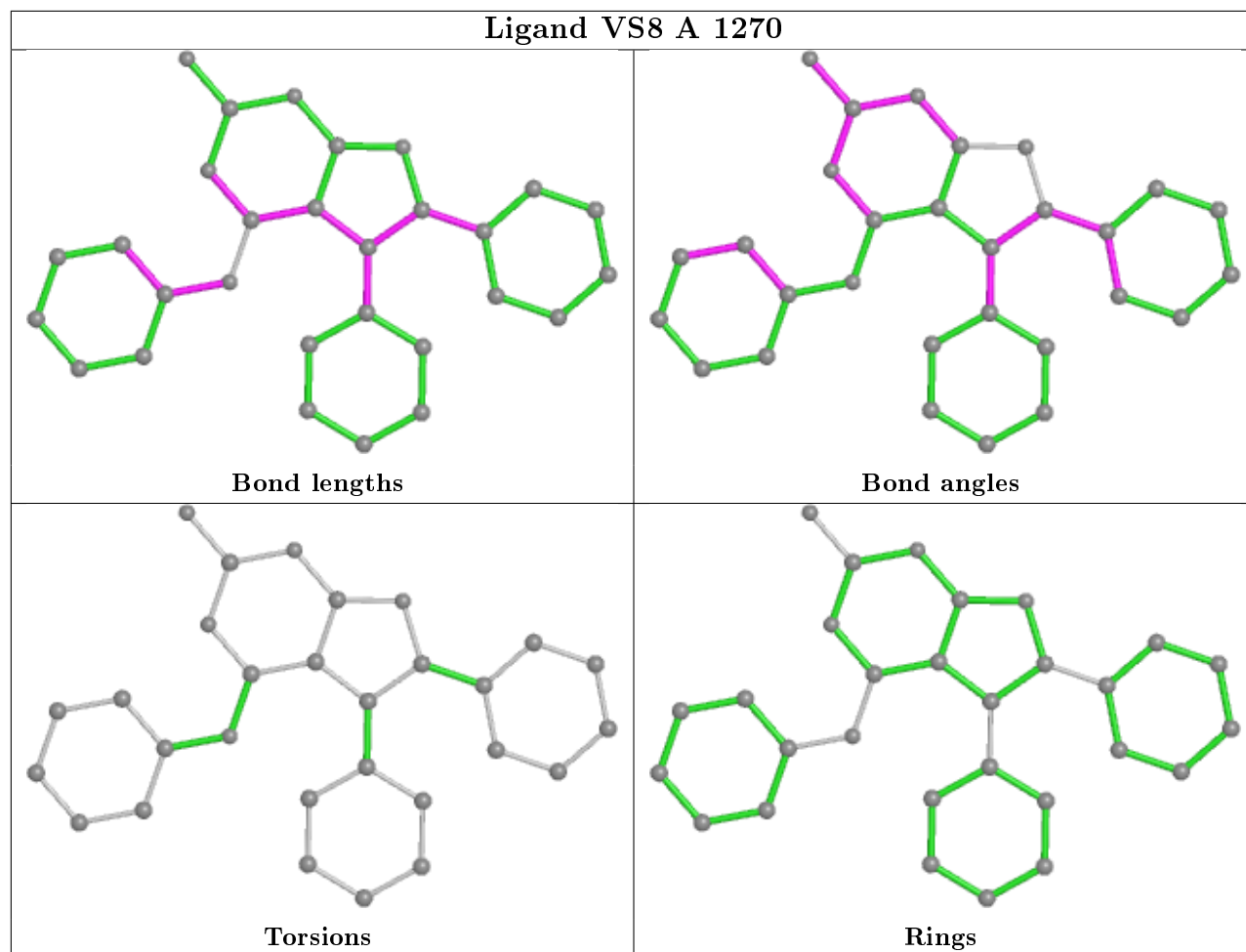
There are no ring outliers.

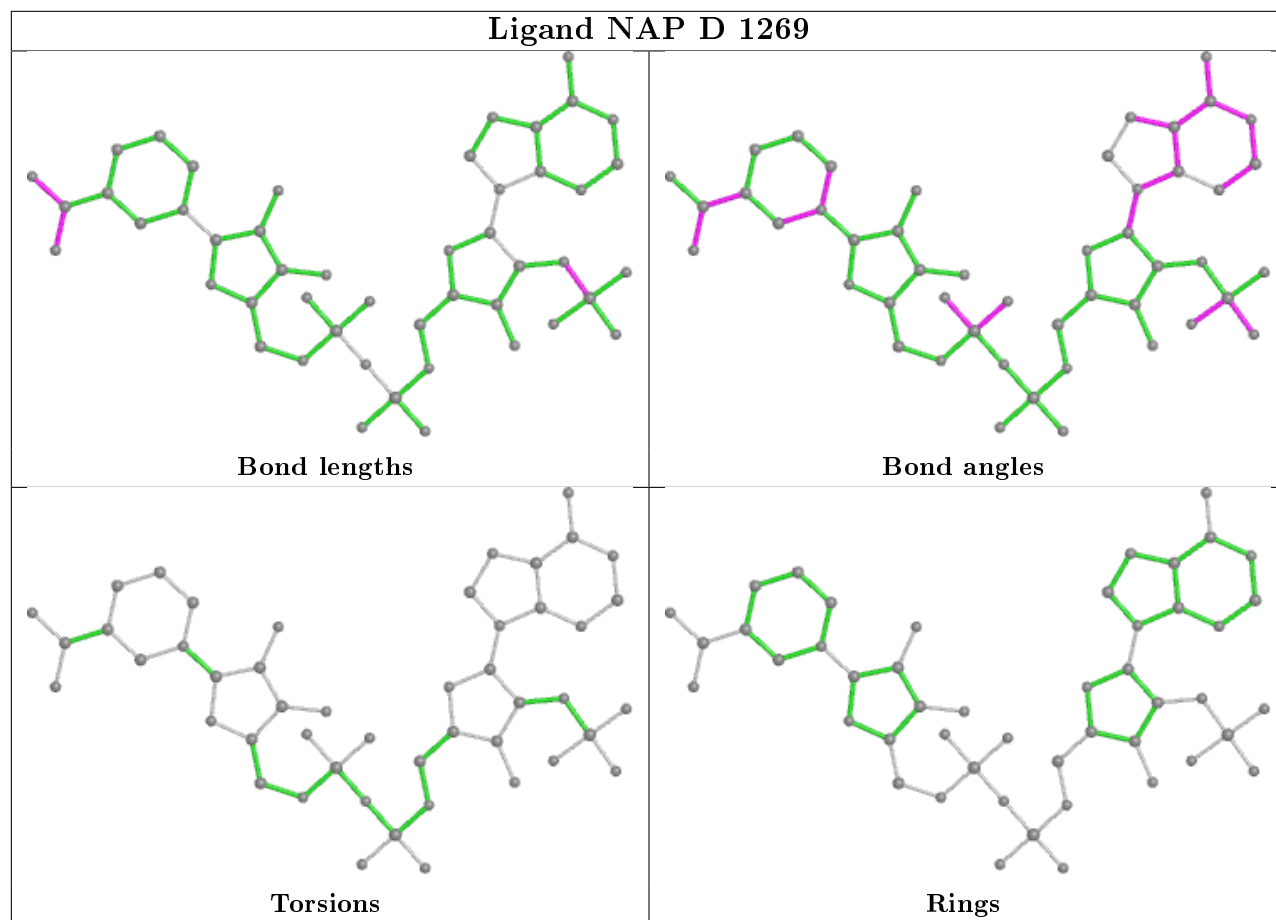
6 monomers are involved in 12 short contacts:

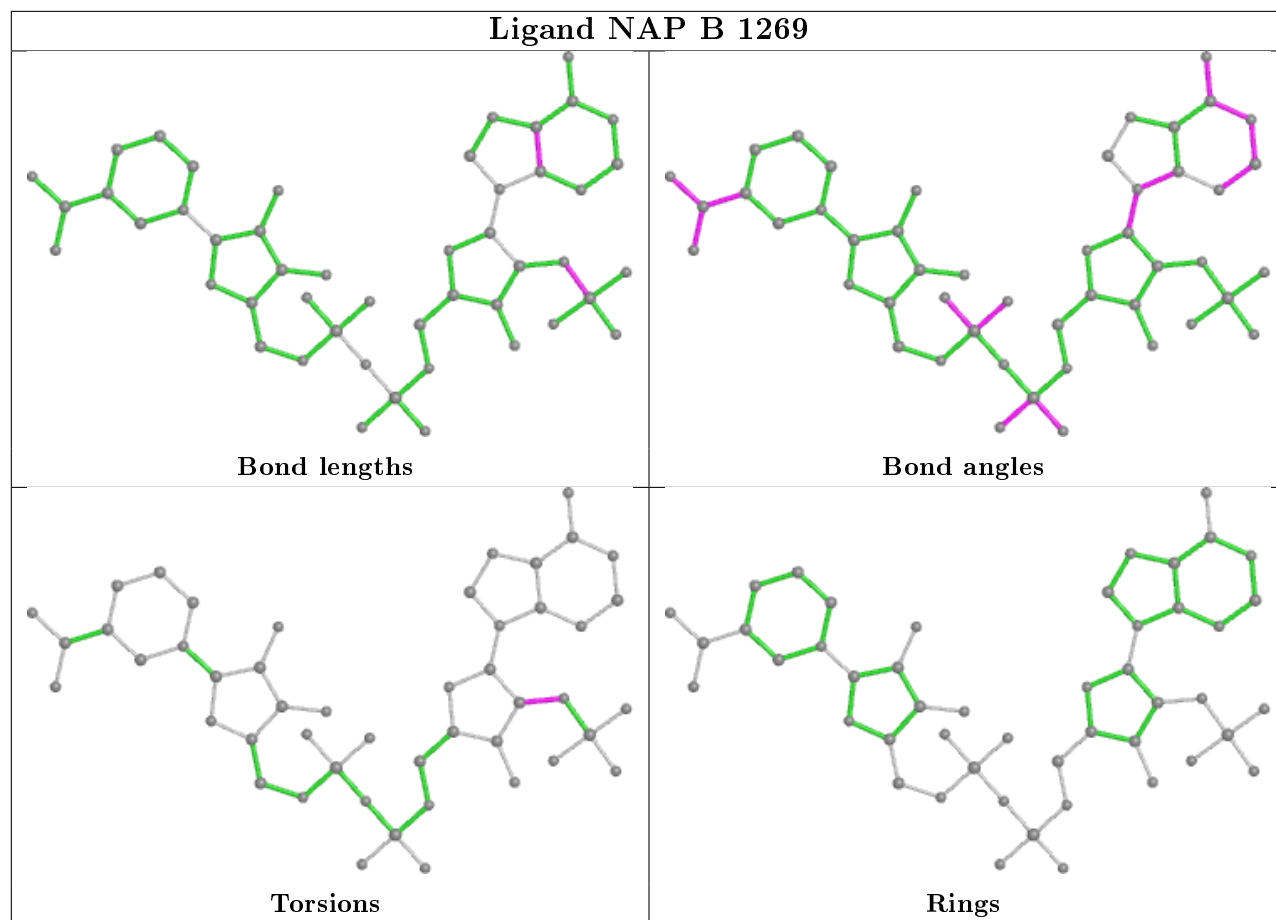
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	1270	VS8	2	0
3	A	1270	VS8	2	0
2	D	1269	NAP	2	0
3	D	1270	VS8	3	0
2	C	1269	NAP	1	0
3	B	1270	VS8	2	0

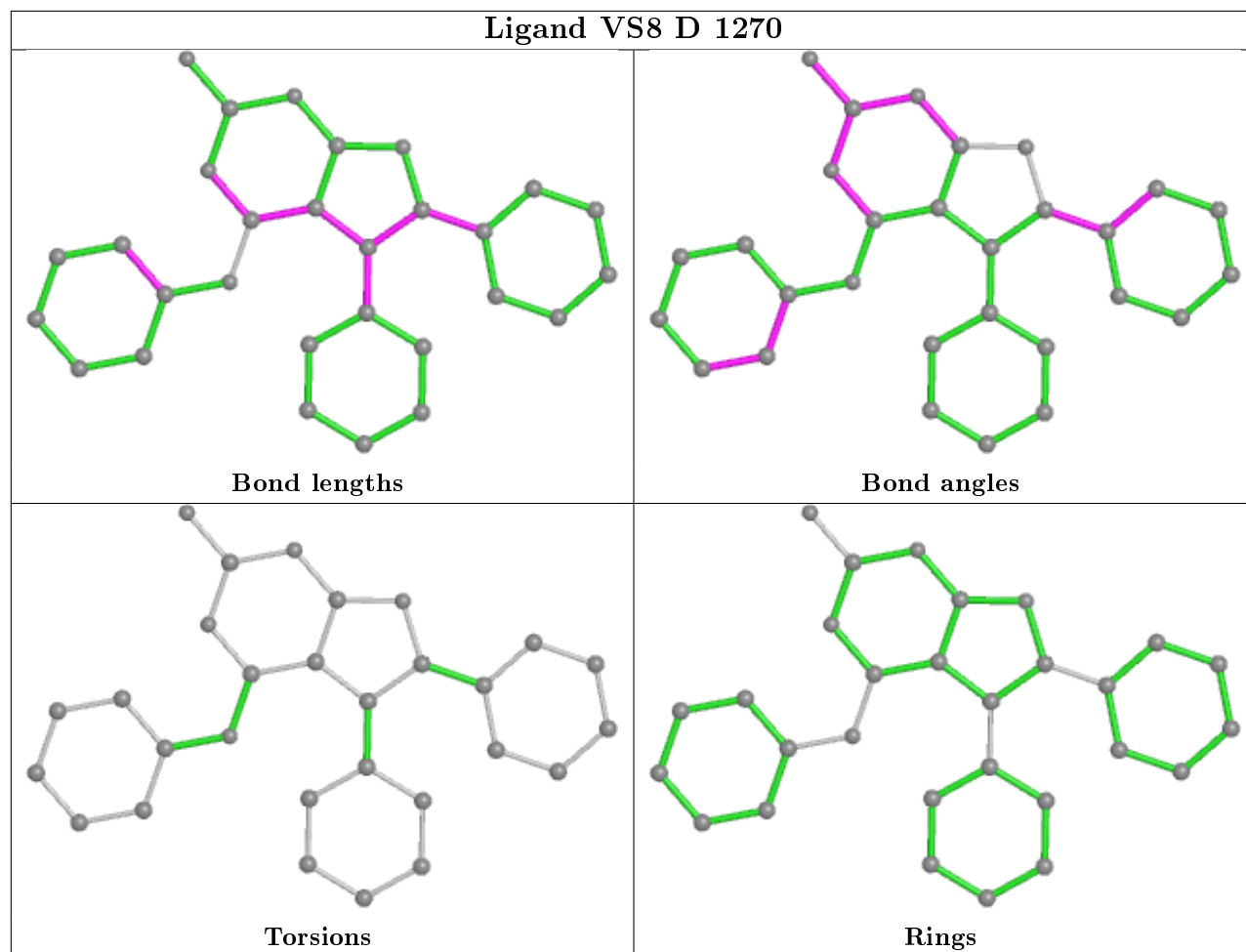
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

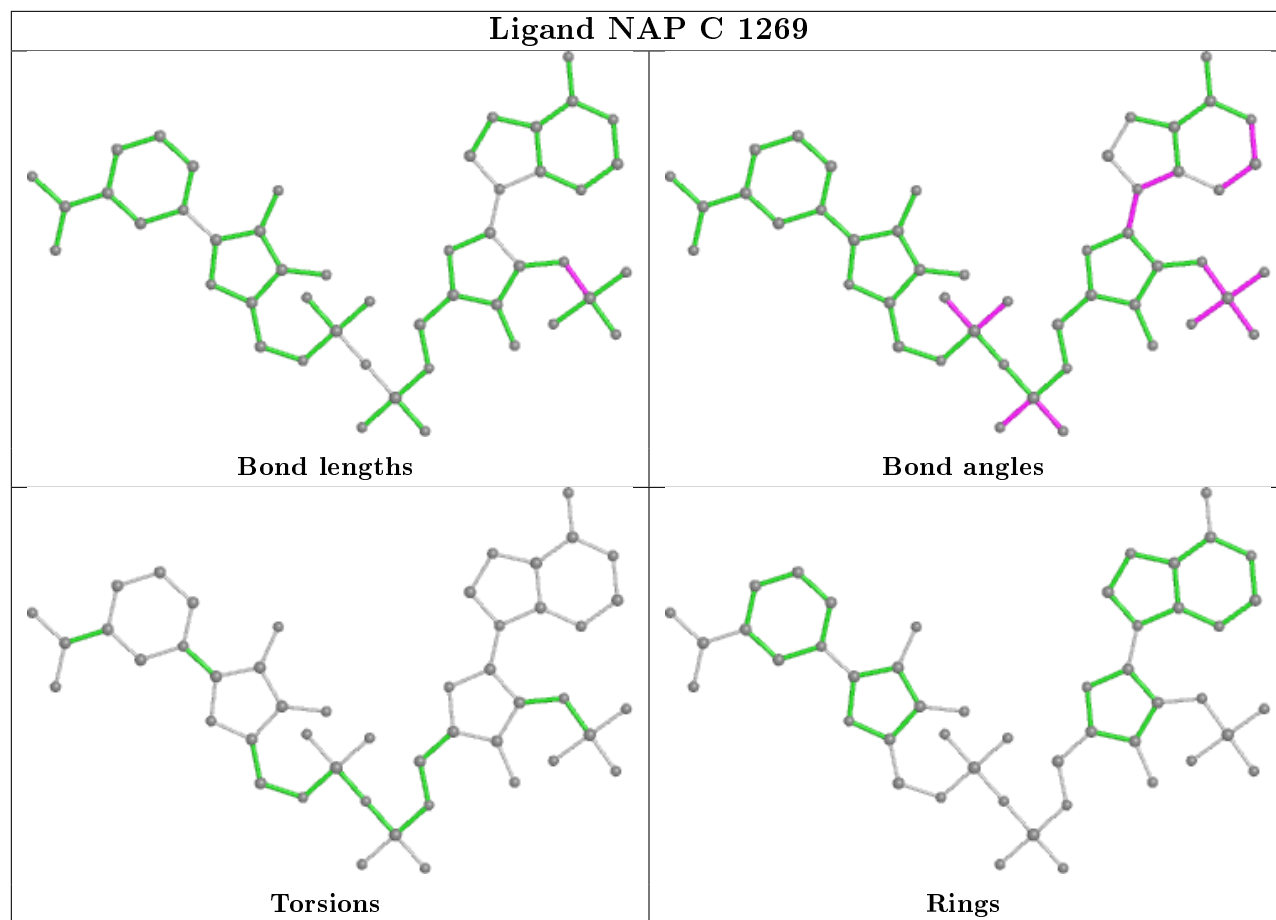


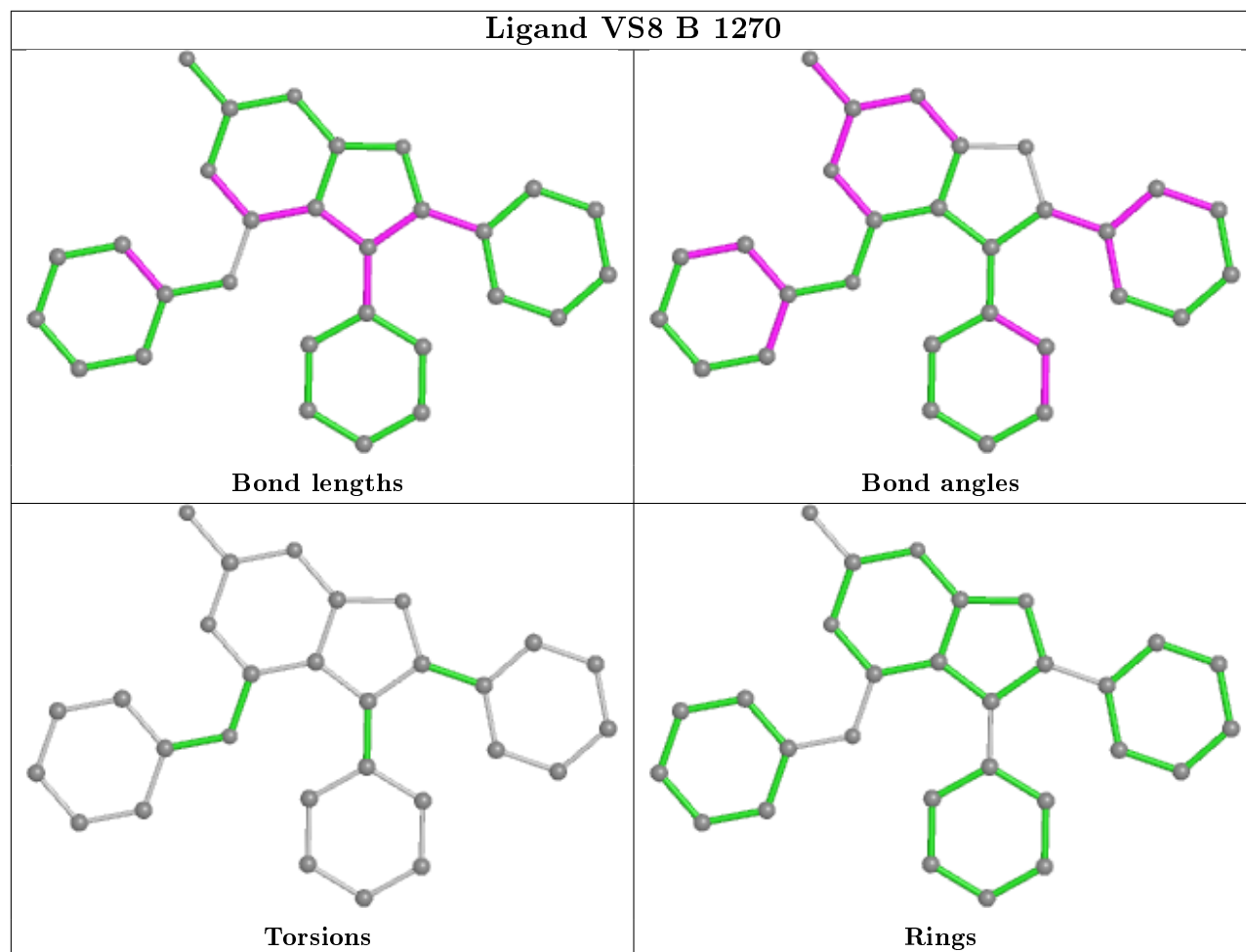


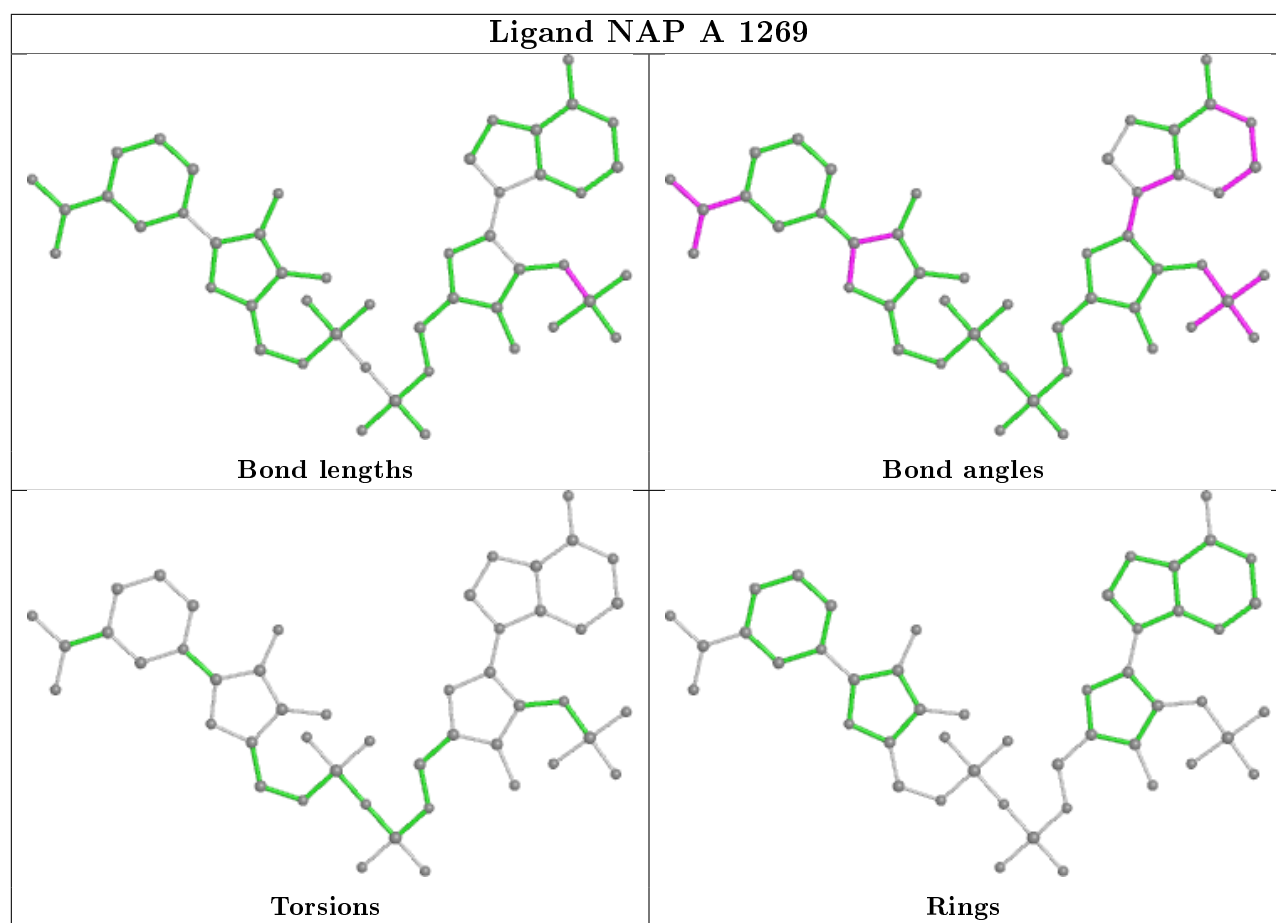












5.7 Other polymers [i](#)

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 ⓘ

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	251/288 (87%)	0.17	16 (6%) 19 18	9, 18, 45, 65	0
1	B	249/288 (86%)	-0.10	7 (2%) 53 52	9, 17, 35, 53	0
1	C	249/288 (86%)	-0.07	12 (4%) 30 29	8, 16, 38, 63	0
1	D	251/288 (87%)	0.05	7 (2%) 53 52	9, 17, 37, 69	0
All	All	1000/1152 (86%)	0.01	42 (4%) 36 34	8, 17, 38, 69	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	211	VAL	8.3
1	C	212	ALA	7.4
1	D	112	ASN	5.4
1	D	104	GLN	5.4
1	C	113	GLY	5.0
1	B	211	VAL	5.0
1	A	211	VAL	4.9
1	B	212	ALA	4.6
1	A	104	GLN	4.3
1	B	113	GLY	3.8
1	D	113	GLY	3.7
1	C	216	GLU	3.7
1	D	2	GLU	3.6
1	B	152	SER	3.6
1	C	209	LEU	3.6
1	B	221	TRP	3.4
1	A	2	GLU	3.3
1	A	151	SER	3.3
1	C	152	SER	3.2
1	D	152	SER	3.1
1	C	223	ARG	3.0

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Mol	Chain	Res	Type	RSRZ
1	D	211	VAL	2.9
1	A	223	ARG	2.9
1	A	212	ALA	2.8
1	A	113	GLY	2.8
1	A	216	GLU	2.8
1	C	221	TRP	2.7
1	A	51	GLU	2.6
1	A	50	LYS	2.6
1	C	214	GLY	2.6
1	D	221	TRP	2.5
1	A	42	VAL	2.5
1	C	210	PRO	2.4
1	A	214	GLY	2.4
1	C	215	GLU	2.4
1	A	152	SER	2.4
1	A	3	ALA	2.3
1	B	213	MET	2.2
1	A	221	TRP	2.2
1	A	219	ASP	2.2
1	C	195	TYR	2.2
1	B	210	PRO	2.2

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no carbohydrates 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
3	VS8	A	1270	29/29	0.87	0.15	19,31,36,36	0
3	VS8	C	1270	29/29	0.88	0.17	16,26,36,36	0

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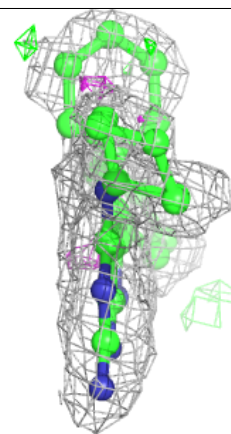
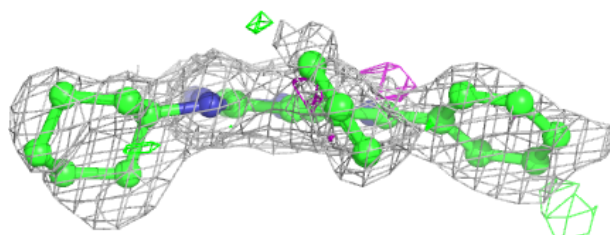
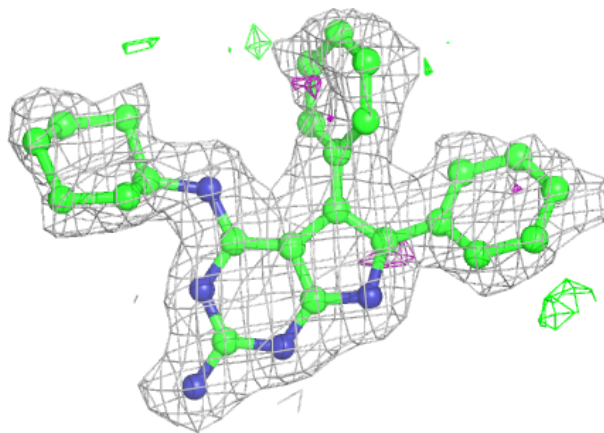
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	VS8	D	1270	29/29	0.89	0.15	15,25,33,35	0
3	VS8	B	1270	29/29	0.89	0.17	16,26,38,38	0
2	NAP	D	1269	48/48	0.97	0.07	11,14,18,19	0
2	NAP	B	1269	48/48	0.97	0.07	11,14,16,17	0
2	NAP	C	1269	48/48	0.98	0.07	9,12,14,16	0
2	NAP	A	1269	48/48	0.98	0.07	12,14,17,20	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.

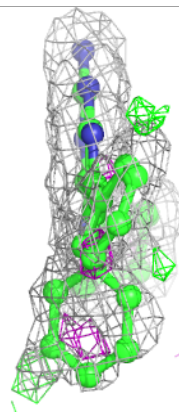
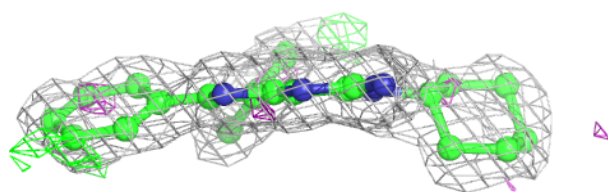
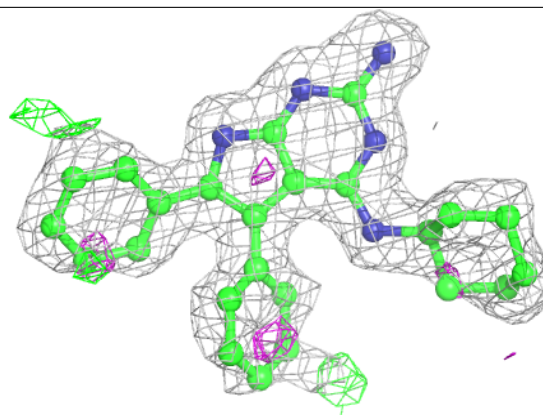
Electron density around VS8 A 1270:

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



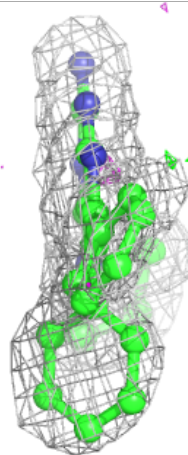
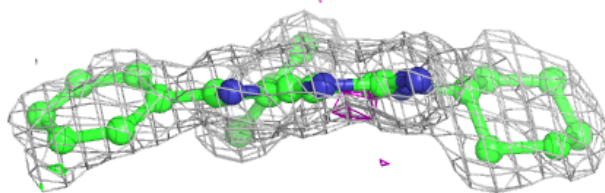
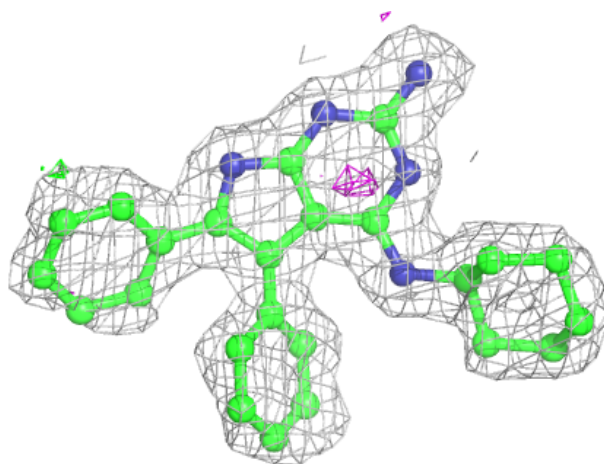
Electron density around VS8 C 1270:

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



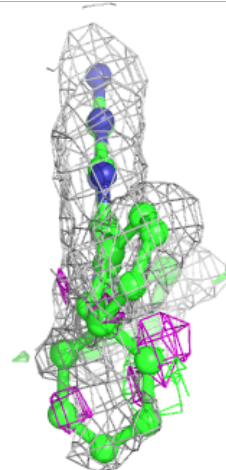
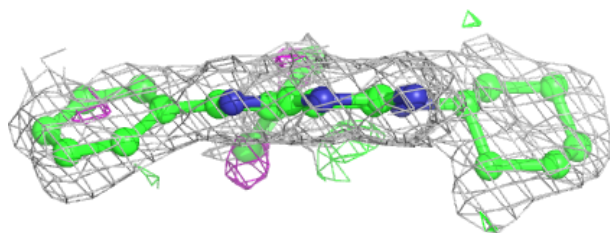
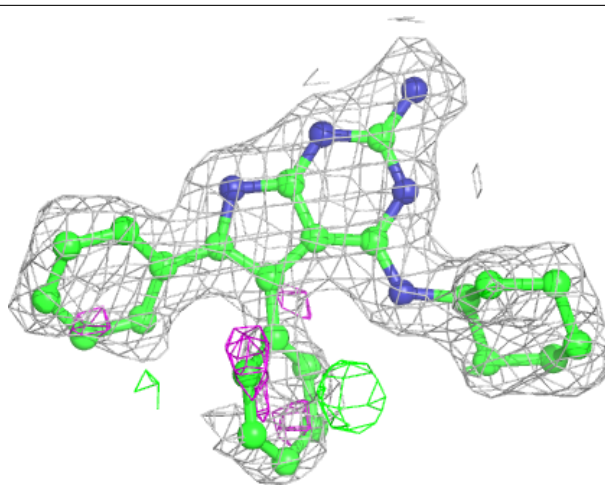
Electron density around VS8 D 1270:

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



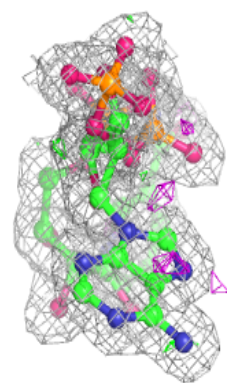
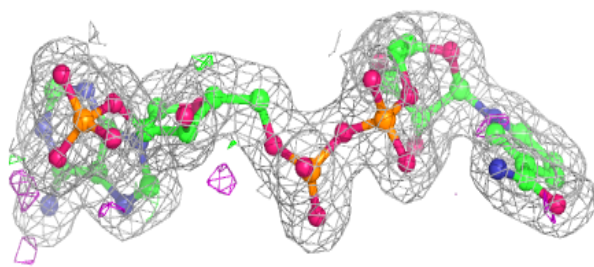
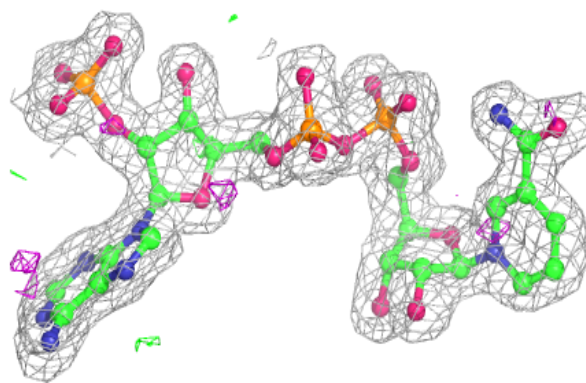
Electron density around VS8 B 1270:

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

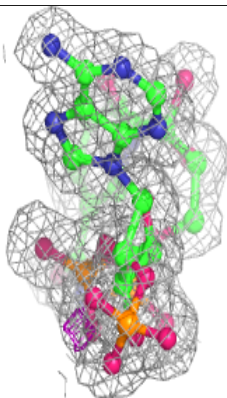
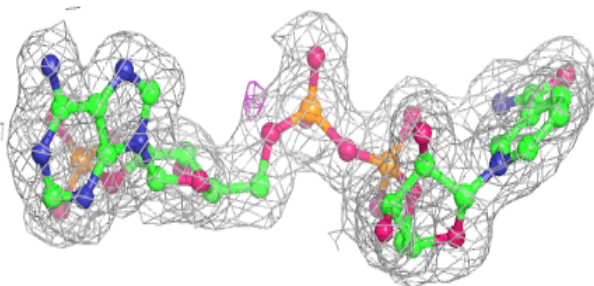
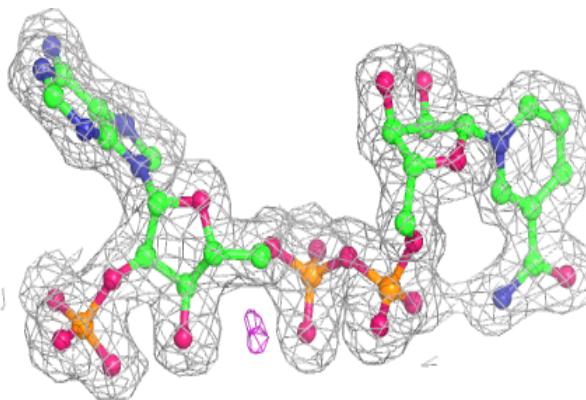


Electron density around NAP D 1269:

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

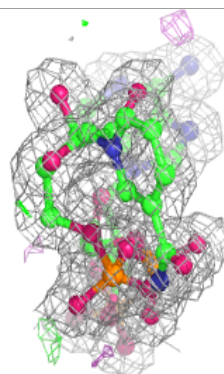
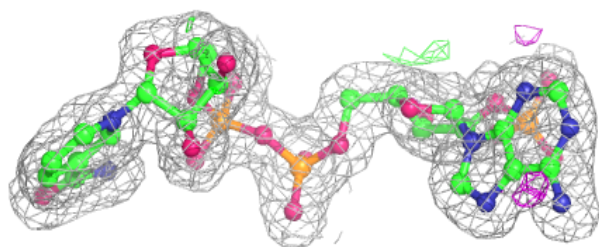
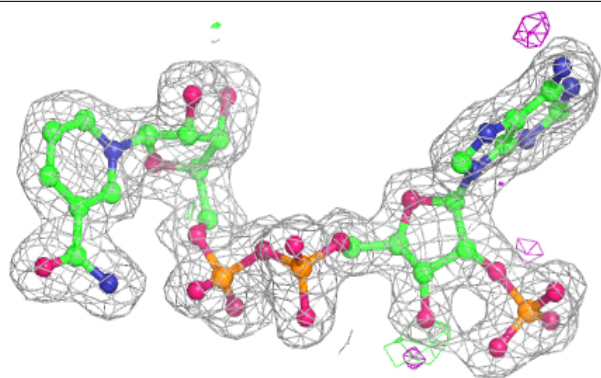
**Electron density around NAP B 1269:**

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

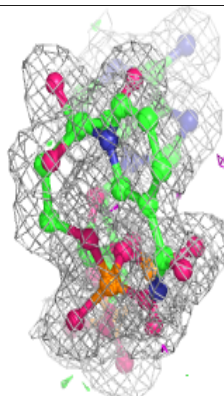
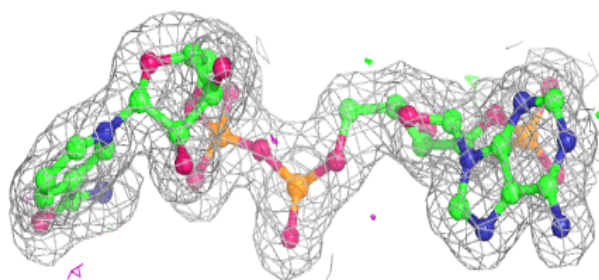
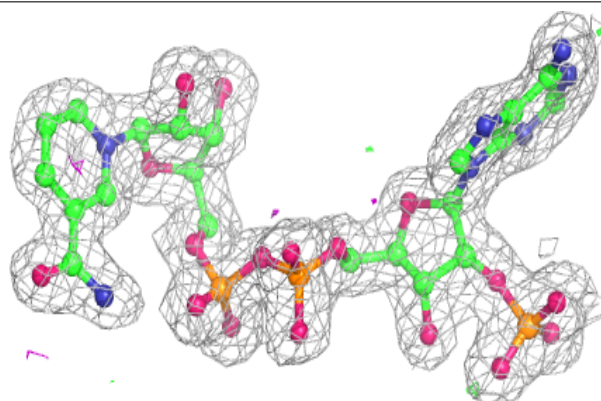


Electron density around NAP C 1269:

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

**Electron density around NAP A 1269:**

$2mF_o-DF_c$ (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.