



Full wwPDB X-ray Structure Validation Report ⓘ

May 20, 2026 – 08:14 PM EDT

PDB ID : 9Y92 / pdb_00009y92
Title : Nanorana parkeri saxiphilin:STX-C13-OBz (co-crystal)
Authors : Zakrzewska, S.; Minor, D.L.
Deposited on : 2025-09-12
Resolution : 1.90 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

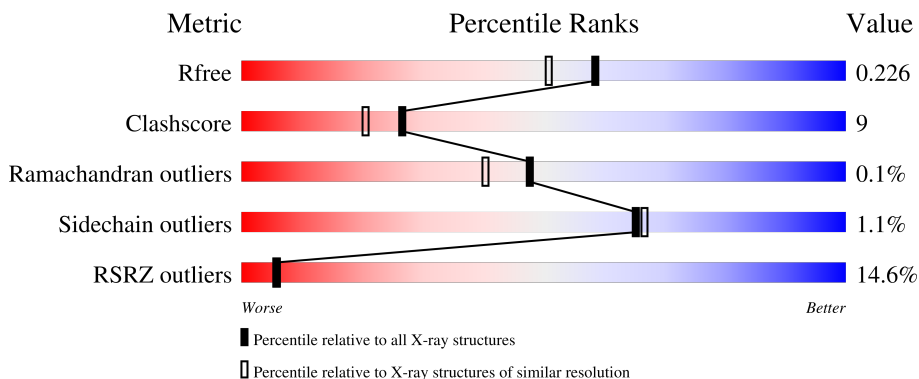
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	854	<div> <div>14%</div> <div>79%</div> <div>17%</div> <div>.</div> </div>

2 Entry composition [i](#)

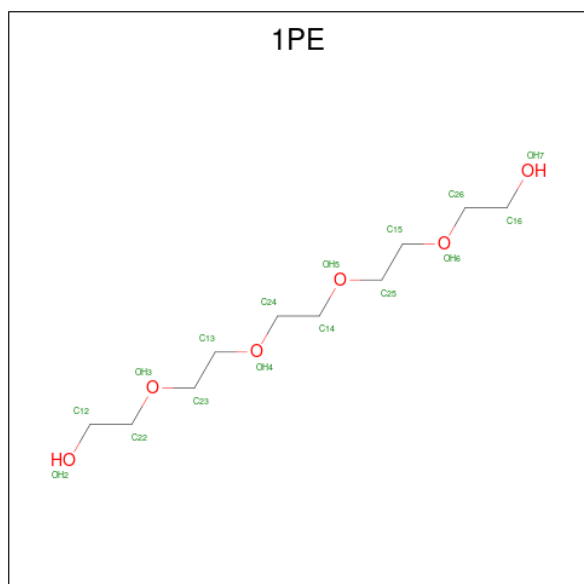
There are 4 unique types of molecules in this entry. The entry contains 6841 atoms, of which 22 are hydrogens and 0 are deuteriums.

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 Saxiphilin.

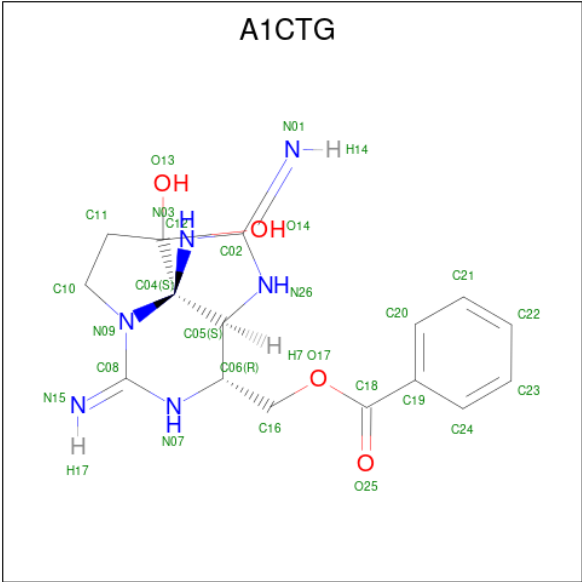
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	821	6377	3995	1103	1220	59	0	2	0

- Molecule 2 is PENTAETHYLENE GLYCOL (CCD ID: 1PE) (formula: $C_{10}H_{22}O_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	H	O		
2	A	1	38	10	22	6	0	0

- Molecule 3 is [(2Z,3aS,4R,6Z,7R,10aS)-10,10-dihydroxy-2,6-diiminooctahydro-1H,8H-pyrrolo[1,2-c]purin-4-yl]methyl benzoate (CCD ID: A1CTG) (formula: $C_{16}H_{20}N_6O_4$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			26	16	6	4		

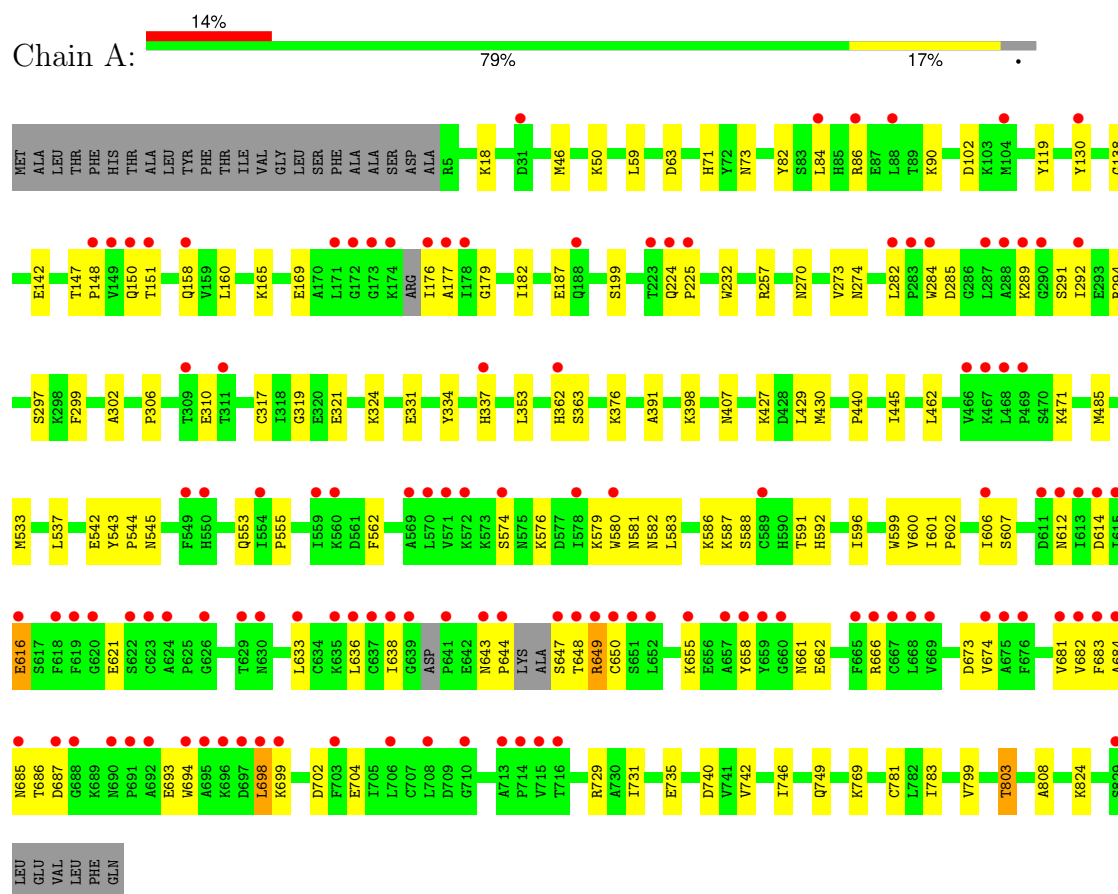
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	400	Total	O	0	0
			400	400		

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: Saxiphilin



4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α , β , γ	228.64Å 228.64Å 67.18Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	43.21 – 1.90 43.21 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.9 (43.21-1.90) 99.9 (43.21-1.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.41 (at 1.89Å)	Xtriage
Refinement program	PHENIX (1.21.2_5419: ???)	Depositor
R, R_{free}	0.198 , 0.226 0.198 , 0.226	Depositor DCC
R_{free} test set	4989 reflections (4.83%)	wwPDB-VP
Wilson B-factor (Å ²)	46.7	Xtriage
Anisotropy	0.317	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 47.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.010 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6841	wwPDB-VP
Average B, all atoms (Å ²)	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.60% 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: A1CTG, 1PE

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	0.32	0/6528	0.46	0/8826

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	6377	0	6179	117	0
2	A	16	22	22	1	0
3	A	26	0	0	0	0
4	A	400	0	0	9	0
All	All	6819	22	6201	117	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 9.

All (117) 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:84:LEU:HD22	1:A:824:LYS:HE3	1.34	1.07
1:A:102:ASP:H	1:A:803:THR:HG21	1.23	0.99
1:A:616:GLU:HA	1:A:633:LEU:HD21	1.43	0.99
1:A:73:ASN:HA	1:A:398:LYS:HE3	1.55	0.87
1:A:102:ASP:N	1:A:803:THR:HG21	1.92	0.84
1:A:587:LYS:HG2	1:A:673:ASP:OD2	1.78	0.82
1:A:799:VAL:O	1:A:803:THR:HG23	1.81	0.80
1:A:661:ASN:HB3	1:A:681:VAL:HG22	1.64	0.79
1:A:86:ARG:HG2	1:A:86:ARG:HH11	1.49	0.78
1:A:662:GLU:HG3	1:A:685:ASN:OD1	1.88	0.73
1:A:50:LYS:HE2	1:A:71:HIS:O	1.89	0.72
1:A:225:PRO:HG3	1:A:362[A]:HIS:ND1	2.06	0.70
1:A:46:MET:CE	1:A:63:ASP:HB3	2.21	0.70
1:A:179:GLY:HA2	1:A:199:SER:OG	1.92	0.69
1:A:684:ALA:C	1:A:685:ASN:HD22	2.01	0.68
1:A:224:GLN:HG3	1:A:225:PRO:O	1.93	0.68
1:A:147:THR:HB	1:A:148:PRO:HD2	1.77	0.67
1:A:187:GLU:HA	1:A:187:GLU:OE1	1.93	0.67
1:A:638:ILE:HG13	1:A:658:TYR:CE2	2.30	0.66
1:A:147:THR:H	1:A:150:GLN:NE2	1.93	0.66
1:A:699:LYS:HB2	1:A:702:ASP:OD2	1.95	0.65
1:A:485:MET:HE3	1:A:769:LYS:HD3	1.79	0.65
1:A:542:GLU:OE1	1:A:729:ARG:HD3	1.96	0.64
1:A:294:ARG:HG2	1:A:310:GLU:OE2	1.97	0.64
1:A:362[A]:HIS:HD2	4:A:1344:HOH:O	1.79	0.64
1:A:86:ARG:HD3	1:A:232:TRP:CE2	2.34	0.63
1:A:661:ASN:HB3	1:A:681:VAL:CG2	2.29	0.62
1:A:285:ASP:OD1	1:A:289:LYS:HD2	2.02	0.60
1:A:602:PRO:O	1:A:606:ILE:HG12	2.02	0.59
1:A:165:LYS:HE2	1:A:169:GLU:OE2	2.03	0.58
1:A:621:GLU:HB3	1:A:636:LEU:CD1	2.34	0.57
1:A:587:LYS:N	1:A:587:LYS:HD3	2.20	0.57
1:A:225:PRO:HG3	1:A:362[B]:HIS:ND1	2.18	0.57
1:A:648:THR:O	1:A:649:ARG:HB2	2.03	0.57
1:A:86:ARG:HD2	4:A:1332:HOH:O	2.05	0.57
1:A:749:GLN:OE1	1:A:781:CYS:HB2	2.04	0.56
1:A:151:THR:HG22	4:A:1341:HOH:O	2.04	0.56
1:A:644:PRO:HD3	1:A:647:SER:HB3	1.87	0.56
1:A:73:ASN:CA	1:A:398:LYS:HE3	2.32	0.56
1:A:596:ILE:HA	1:A:600:VAL:HB	1.89	0.55
1:A:130:TYR:CE2	1:A:147:THR:HG22	2.42	0.55
1:A:337[B]:HIS:ND1	1:A:427:LYS:HD2	2.21	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:376:LYS:HE2	4:A:1187:HOH:O	2.05	0.55
1:A:553:GLN:O	1:A:555:PRO:HD3	2.07	0.55
1:A:362[B]:HIS:HD2	4:A:1344:HOH:O	1.89	0.54
1:A:284:TRP:HA	1:A:289:LYS:HD3	1.88	0.54
1:A:580:TRP:CE3	1:A:583:LEU:HD11	2.42	0.54
1:A:693:GLU:HG2	1:A:694:TRP:N	2.22	0.54
1:A:647:SER:HB2	1:A:655:LYS:HE3	1.89	0.53
1:A:638:ILE:O	1:A:649:ARG:HG2	2.08	0.53
1:A:407:ASN:ND2	1:A:440:PRO:HB3	2.24	0.52
1:A:579:LYS:H	1:A:582:ASN:HD21	1.56	0.52
1:A:46:MET:HE1	1:A:63:ASP:HB3	1.91	0.52
1:A:46:MET:HE2	1:A:59:LEU:HD13	1.92	0.51
1:A:621:GLU:C	1:A:636:LEU:HD13	2.35	0.51
1:A:306:PRO:HG2	1:A:334:TYR:HA	1.93	0.50
1:A:86:ARG:HG2	1:A:86:ARG:NH1	2.21	0.50
1:A:607:SER:HB2	1:A:612:ASN:HA	1.93	0.49
1:A:177:ALA:HB3	1:A:182:ILE:HD11	1.94	0.48
1:A:257:ARG:HA	1:A:302:ALA:O	2.13	0.48
1:A:294:ARG:O	1:A:297:SER:HB3	2.12	0.48
1:A:666:ARG:HG2	1:A:694:TRP:CZ2	2.49	0.48
1:A:682:VAL:HG13	1:A:683:PHE:N	2.28	0.48
1:A:666:ARG:HG2	1:A:694:TRP:CH2	2.47	0.48
1:A:587:LYS:HA	1:A:621:GLU:O	2.12	0.48
1:A:599:TRP:C	1:A:602:PRO:HD2	2.38	0.48
1:A:274:ASN:HB3	2:A:901:1PE:H142	1.95	0.48
1:A:84:LEU:CD2	1:A:824:LYS:HE3	2.24	0.48
1:A:543:TYR:CZ	1:A:545:ASN:HB3	2.49	0.47
1:A:803:THR:HA	1:A:808:ALA:N	2.29	0.47
1:A:158:GLN:HG2	4:A:1082:HOH:O	2.14	0.47
1:A:160:LEU:HA	4:A:1123:HOH:O	2.13	0.47
1:A:285:ASP:OD1	1:A:285:ASP:N	2.47	0.47
1:A:321:GLU:OE2	1:A:321:GLU:HA	2.15	0.47
1:A:587:LYS:HB3	1:A:621:GLU:O	2.14	0.47
1:A:661:ASN:O	1:A:681:VAL:HG11	2.14	0.46
1:A:142:GLU:OE2	1:A:142:GLU:HA	2.14	0.46
1:A:543:TYR:HB2	1:A:783:ILE:HD12	1.98	0.46
1:A:537:LEU:CD1	1:A:735:GLU:HG3	2.47	0.46
1:A:86:ARG:HD3	1:A:232:TRP:NE1	2.31	0.45
1:A:176:ILE:HD12	1:A:176:ILE:O	2.16	0.45
1:A:574:SER:O	1:A:576:LYS:HG3	2.16	0.45
1:A:130:TYR:CZ	1:A:147:THR:HG22	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:317:CYS:HB3	1:A:331:GLU:OE1	2.17	0.44
1:A:586:LYS:C	1:A:587:LYS:HD3	2.43	0.44
1:A:614:ASP:HB3	1:A:616:GLU:HG3	2.00	0.44
1:A:638:ILE:HD11	1:A:658:TYR:OH	2.18	0.44
1:A:429:LEU:O	1:A:430:MET:HB2	2.17	0.44
1:A:46:MET:HE2	1:A:59:LEU:CD1	2.47	0.43
1:A:581:ASN:C	1:A:581:ASN:OD1	2.60	0.43
1:A:544:PRO:HD2	1:A:562:PHE:O	2.18	0.43
1:A:18:LYS:HE2	1:A:18:LYS:HB2	1.77	0.43
1:A:742:VAL:O	1:A:746:ILE:HG12	2.18	0.43
1:A:273:VAL:HG11	1:A:292:ILE:HG12	1.99	0.43
1:A:291:SER:OG	1:A:294:ARG:HG3	2.19	0.43
1:A:471:LYS:HD3	1:A:740:ASP:OD2	2.18	0.43
1:A:579:LYS:H	1:A:582:ASN:ND2	2.16	0.42
1:A:225:PRO:HG3	1:A:362[A]:HIS:HD1	1.82	0.42
1:A:537:LEU:HD12	1:A:735:GLU:HG3	2.00	0.42
1:A:803:THR:HG22	1:A:808:ALA:HB2	2.01	0.42
1:A:643:ASN:HD22	1:A:643:ASN:H	1.67	0.42
1:A:591:THR:O	1:A:592:HIS:HB3	2.19	0.42
1:A:119:TYR:CZ	1:A:138:GLY:HA3	2.55	0.42
1:A:319:GLY:O	1:A:324:LYS:HB2	2.19	0.42
1:A:601:ILE:HB	1:A:602:PRO:HD3	2.02	0.41
1:A:682:VAL:CG1	1:A:683:PHE:N	2.84	0.41
1:A:90:LYS:NZ	4:A:1023:HOH:O	2.53	0.41
1:A:731:ILE:HD12	1:A:731:ILE:N	2.34	0.41
1:A:270:ASN:ND2	1:A:462:LEU:HD12	2.35	0.41
1:A:682:VAL:O	1:A:686:THR:HG23	2.20	0.41
1:A:224:GLN:OE1	1:A:225:PRO:HG2	2.20	0.41
1:A:533:MET:HE2	4:A:1143:HOH:O	2.20	0.40
1:A:82:TYR:CE1	1:A:391:ALA:HB2	2.56	0.40
1:A:588:SER:HA	1:A:674:VAL:O	2.22	0.40
1:A:647:SER:N	1:A:655:LYS:HB3	2.36	0.40
1:A:687:ASP:N	1:A:698:LEU:O	2.51	0.40
1:A:282:LEU:CD1	1:A:299:PHE:HB2	2.52	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	815/854 (95%)	779 (96%)	35 (4%)	1 (0%)	48	40

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	649	ARG

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	707/731 (97%)	699 (99%)	8 (1%)	65	67

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

Mol	Chain	Res	Type
1	A	353	LEU
1	A	363	SER
1	A	445	ILE
1	A	616	GLU
1	A	650	CYS
1	A	698	LEU
1	A	704	GLU
1	A	803	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	145	ASN
1	A	150	GLN
1	A	158	GLN
1	A	188	GLN
1	A	190	ASN
1	A	230	HIS
1	A	249	GLN
1	A	379	ASN
1	A	407	ASN
1	A	437	HIS
1	A	608	ASN

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 ⓘ

2 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	1PE	A	901	-	15,15,15	0.52	0	14,14,14	0.30	0
3	A1CTG	A	902	-	23,29,29	0.53	0	22,45,45	1.10	3 (13%)

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	1PE	A	901	-	-	6/13/13/13	-
3	A1CTG	A	902	-	-	0/9/57/57	0/4/4/4

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	902	A1CTG	N03-C04-N09	-2.87	107.50	112.45
3	A	902	A1CTG	C05-C06-N07	2.07	111.14	108.86
3	A	902	A1CTG	N03-C02-N26	2.04	111.29	109.21

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	901	1PE	OH6-C15-C25-OH5
2	A	901	1PE	C12-C22-OH3-C23
2	A	901	1PE	C14-C24-OH4-C13
2	A	901	1PE	OH4-C13-C23-OH3
2	A	901	1PE	C25-C15-OH6-C26
2	A	901	1PE	C16-C26-OH6-C15

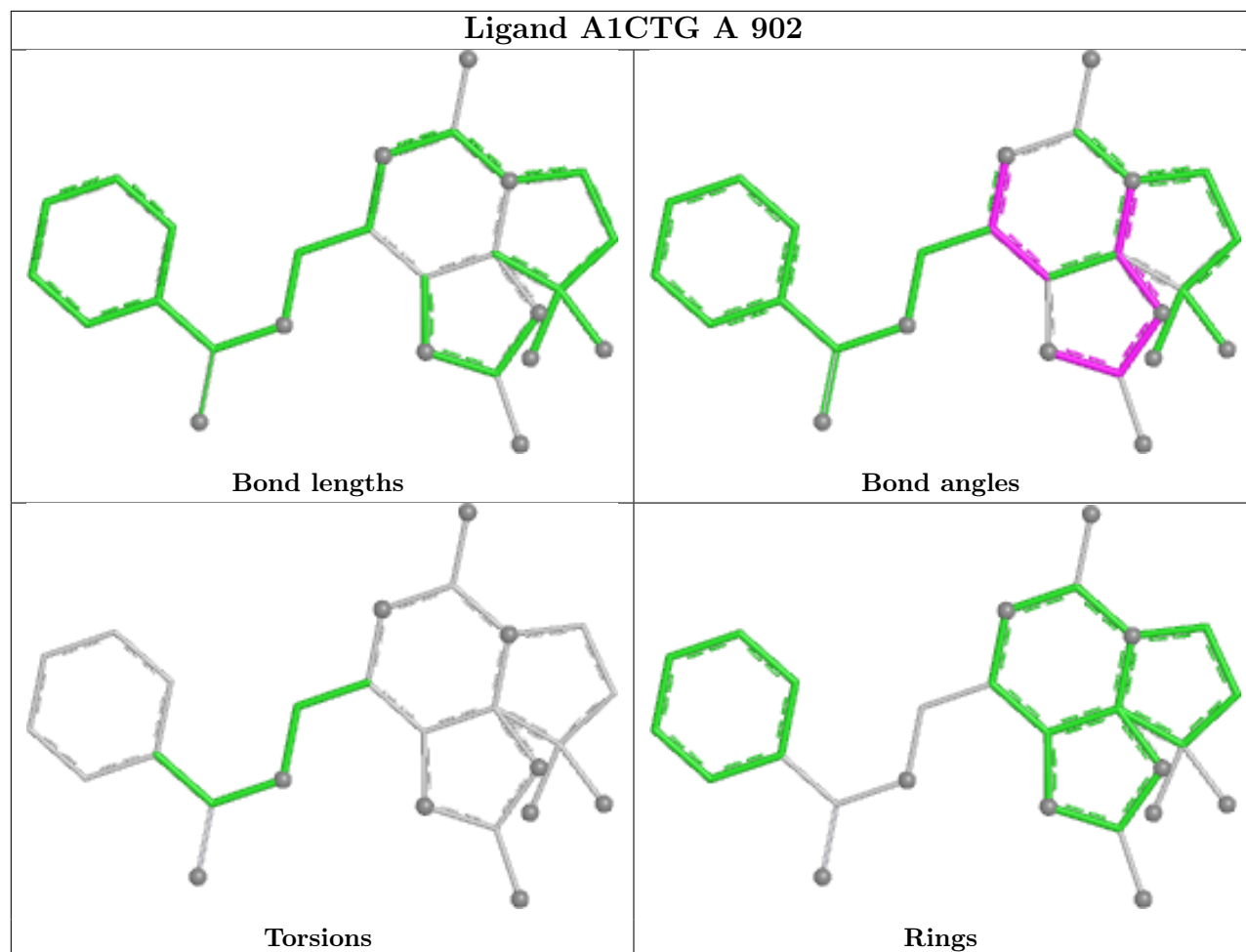
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	901	1PE	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and

any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	821/854 (96%)	0.90	120 (14%) 6 6	27, 52, 103, 137	2 (0%)

All (120) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	176	ILE	9.3
1	A	613	ILE	7.4
1	A	641	PRO	7.0
1	A	638	ILE	6.0
1	A	284	TRP	5.2
1	A	466	VAL	5.1
1	A	554	ILE	4.5
1	A	648	THR	4.5
1	A	639	GLY	4.4
1	A	644	PRO	4.3
1	A	698	LEU	4.2
1	A	287	LEU	4.0
1	A	177	ALA	4.0
1	A	657	ALA	3.9
1	A	658	TYR	3.8
1	A	571	VAL	3.8
1	A	665	PHE	3.8
1	A	468	LEU	3.8
1	A	283	PRO	3.8
1	A	703	PHE	3.8
1	A	178	ILE	3.7
1	A	695	ALA	3.7
1	A	649	ARG	3.6
1	A	225	PRO	3.5
1	A	647	SER	3.4
1	A	224	GLN	3.4
1	A	669	VAL	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	688	GLY	3.3
1	A	637	CYS	3.3
1	A	149	VAL	3.3
1	A	694	TRP	3.3
1	A	88	LEU	3.2
1	A	570	LEU	3.2
1	A	714	PRO	3.2
1	A	86	ARG	3.2
1	A	624	ALA	3.2
1	A	629	THR	3.1
1	A	148	PRO	3.1
1	A	635	LYS	3.1
1	A	151	THR	3.1
1	A	223	THR	3.1
1	A	569	ALA	3.1
1	A	84	LEU	3.1
1	A	150	GLN	3.0
1	A	682	VAL	3.0
1	A	589	CYS	3.0
1	A	174	LYS	2.9
1	A	684	ALA	2.9
1	A	469	PRO	2.9
1	A	676	PHE	2.9
1	A	559	ILE	2.9
1	A	173	GLY	2.9
1	A	692	ALA	2.8
1	A	651	SER	2.8
1	A	104	MET	2.8
1	A	171	LEU	2.8
1	A	620	GLY	2.8
1	A	158	GLN	2.7
1	A	578	ILE	2.7
1	A	130	TYR	2.7
1	A	699	LYS	2.7
1	A	713	ALA	2.6
1	A	633	LEU	2.6
1	A	675	ALA	2.6
1	A	612	ASN	2.6
1	A	683	PHE	2.6
1	A	660	GLY	2.6
1	A	618	PHE	2.5
1	A	362[A]	HIS	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	580	TRP	2.5
1	A	615	ILE	2.5
1	A	630	ASN	2.5
1	A	550	HIS	2.5
1	A	655	LYS	2.5
1	A	626	GLY	2.4
1	A	288	ALA	2.4
1	A	611	ASP	2.4
1	A	636	LEU	2.4
1	A	697	ASP	2.4
1	A	666	ARG	2.4
1	A	681	VAL	2.4
1	A	572	LYS	2.4
1	A	292	ILE	2.4
1	A	290	GLY	2.4
1	A	623	CYS	2.4
1	A	614	ASP	2.3
1	A	668	LEU	2.3
1	A	659	TYR	2.3
1	A	549	PHE	2.3
1	A	289	LYS	2.3
1	A	715	VAL	2.3
1	A	560	LYS	2.3
1	A	650	CYS	2.3
1	A	282	LEU	2.3
1	A	311	THR	2.3
1	A	696	LYS	2.3
1	A	716	THR	2.3
1	A	616	GLU	2.2
1	A	31	ASP	2.2
1	A	467	LYS	2.2
1	A	667	CYS	2.2
1	A	652	LEU	2.2
1	A	691	PRO	2.2
1	A	829	SER	2.2
1	A	619	PHE	2.2
1	A	172	GLY	2.2
1	A	685	ASN	2.2
1	A	606	ILE	2.1
1	A	188	GLN	2.1
1	A	622	SER	2.1
1	A	337[A]	HIS	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	706	LEU	2.1
1	A	309	THR	2.0
1	A	643	ASN	2.0
1	A	708	LEU	2.0
1	A	687	ASP	2.0
1	A	574	SER	2.0
1	A	690	ASN	2.0
1	A	710	GLY	2.0
1	A	674	VAL	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](#)

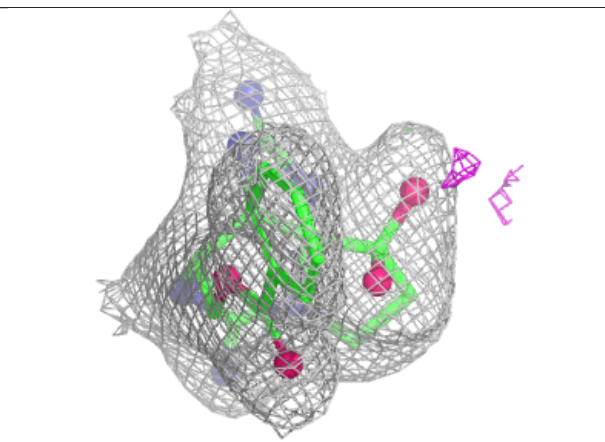
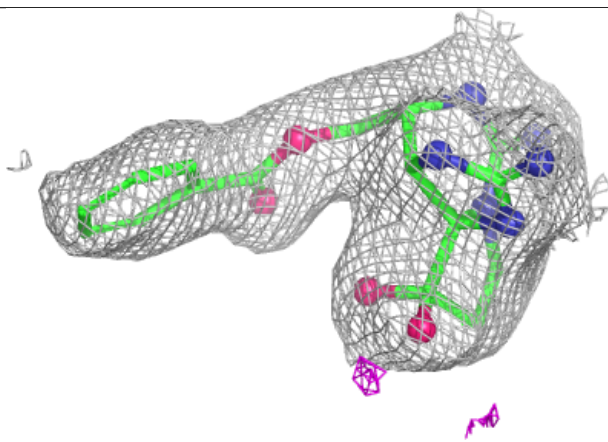
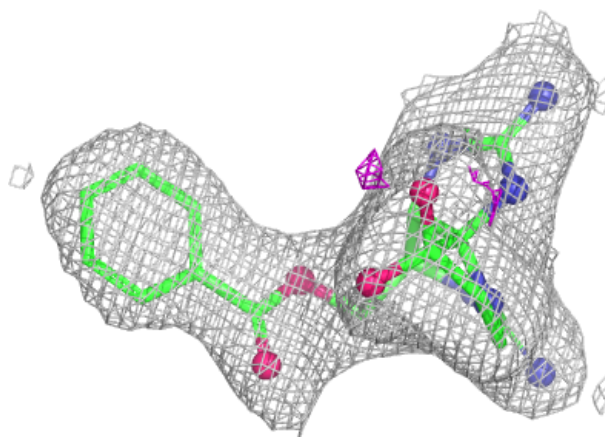
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	1PE	A	901	16/16	0.80	0.16	61,76,92,95	0
3	A1CTG	A	902	26/26	0.94	0.10	45,51,69,71	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 A1CTG A 902:

$2mF_o-DF_c$ (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.