



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 17, 2026 – 11:28 PM UTC

PDB ID : 6KDT / pdb_00006kdt
Title : Crystal structure of human DNMT3B (Q772R)-DNMT3L complex
Authors : Lin, C.-C.; Chen, Y.-P.; Yang, W.-Z.; Shen, C.-K.; Yuan, H.S.
Deposited on : 2019-07-02
Resolution : 2.87 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

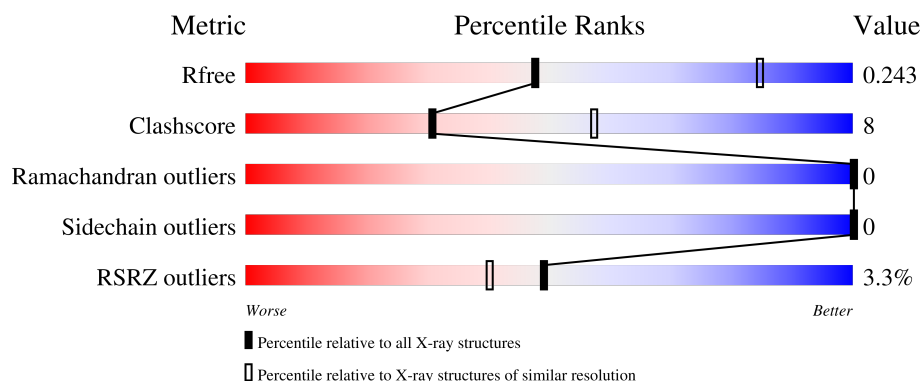
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.87 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	3557 (2.90-2.86)
Clashscore	190562	3801 (2.90-2.86)
Ramachandran outliers	187476	3699 (2.90-2.86)
Sidechain outliers	187428	3702 (2.90-2.86)
RSRZ outliers	180081	3558 (2.90-2.86)

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 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	286	0% 80% 15% 5%
1	D	286	2% 79% 17% 5%
2	B	204	5% 69% 28% .
2	C	204	5% 83% 15% .

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 7895 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 DNA (cytosine-5)-methyltransferase 3B.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	272	Total	C	N	O	S	0	0	0
			2200	1412	391	384	13			
1	D	273	Total	C	N	O	S	0	0	0
			2204	1414	392	385	13			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	568	GLY	-	expression tag	UNP Q9UBC3
A	569	HIS	-	expression tag	UNP Q9UBC3
A	570	MET	-	expression tag	UNP Q9UBC3
A	772	ARG	GLN	engineered mutation	UNP Q9UBC3
D	568	GLY	-	expression tag	UNP Q9UBC3
D	569	HIS	-	expression tag	UNP Q9UBC3
D	570	MET	-	expression tag	UNP Q9UBC3
D	772	ARG	GLN	engineered mutation	UNP Q9UBC3

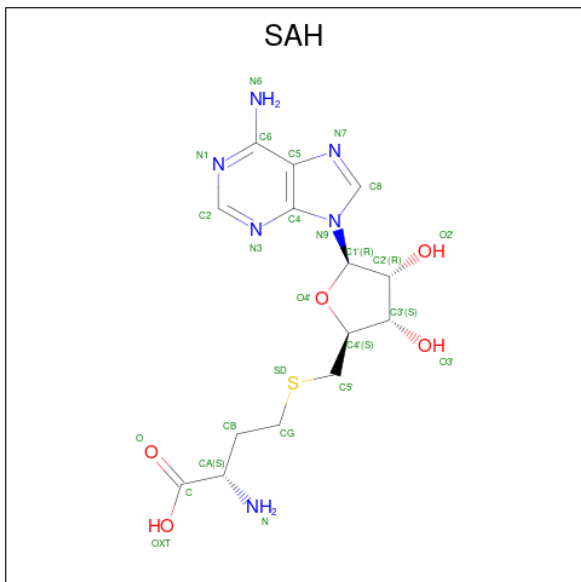
- Molecule 2 is a protein called DNA (cytosine-5)-methyltransferase 3-like.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	198	Total	C	N	O	S	0	0	0
			1631	1061	279	287	4			
2	C	200	Total	C	N	O	S	0	0	0
			1642	1069	281	287	5			

There are 4 discrepancies between the modelled and reference sequences:

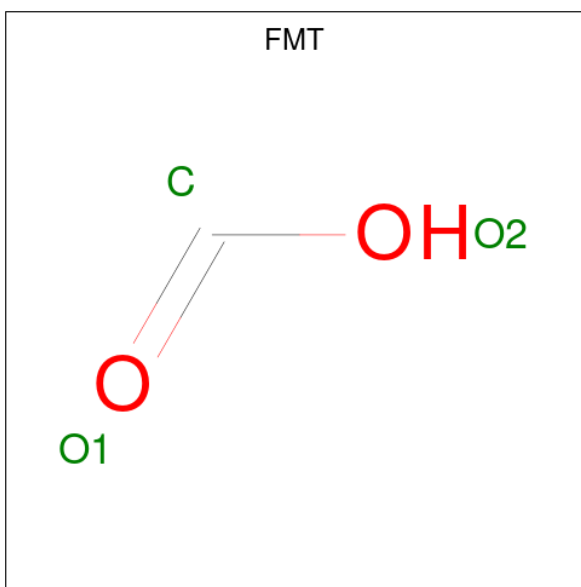
Chain	Residue	Modelled	Actual	Comment	Reference
B	176	GLY	-	expression tag	UNP Q9UJW3
B	177	HIS	-	expression tag	UNP Q9UJW3
C	176	GLY	-	expression tag	UNP Q9UJW3
C	177	HIS	-	expression tag	UNP Q9UJW3

- Molecule 3 is S-ADENOSYL-L-HOMOCYSTEINE (CCD ID: SAH) (formula: $\text{C}_{14}\text{H}_{20}\text{N}_6\text{O}_5\text{S}$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total 26	C 14	N 6	O 5	S 1	0	0
3	D	1	Total 26	C 14	N 6	O 5	S 1	0	0

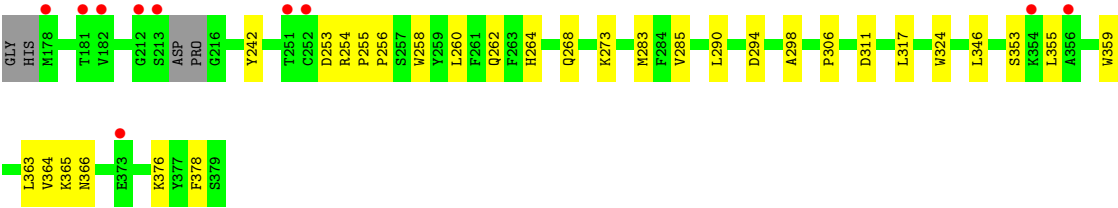
- Molecule 4 is FORMIC ACID (CCD ID: FMT) (formula: CH_2O_2).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			3	1	2		
4	D	1	Total	C	O	0	0
			3	1	2		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	56	Total	O	0	0
			56	56		
5	B	15	Total	O	0	0
			15	15		
5	C	28	Total	O	0	0
			28	28		
5	D	61	Total	O	0	0
			61	61		



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	60.35Å 240.78Å 232.13Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	27.37 – 2.87 27.37 – 2.87	Depositor EDS
% Data completeness (in resolution range)	93.3 (27.37-2.87) 93.4 (27.37-2.87)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.47 (at 2.85Å)	Xtriage
Refinement program	PHENIX 1.14-3260-000	Depositor
R, R_{free}	0.194 , 0.242 0.194 , 0.243	Depositor DCC
R_{free} test set	1786 reflections (4.52%)	wwPDB-VP
Wilson B-factor (Å ²)	53.2	Xtriage
Anisotropy	0.144	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 31.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7895	wwPDB-VP
Average B, all atoms (Å ²)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.26% 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: SAH, FMT

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	0.41	0/2252	0.57	0/3037
1	D	0.42	0/2256	0.58	0/3042
2	B	0.43	0/1683	0.60	0/2290
2	C	0.43	0/1693	0.56	0/2302
All	All	0.42	0/7884	0.58	0/10671

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 [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	2200	0	2212	29	0
1	D	2204	0	2215	41	0
2	B	1631	0	1612	41	0
2	C	1642	0	1631	21	0
3	A	26	0	19	0	0
3	D	26	0	19	2	0
4	A	3	0	1	0	0
4	D	3	0	1	0	0
5	A	56	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	B	15	0	0	0	0
5	C	28	0	0	0	0
5	D	61	0	0	3	0
All	All	7895	0	7710	128	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 8.

All (128) 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)
2:B:210:GLU:HG3	2:B:213:SER:HB3	1.11	1.10
2:B:210:GLU:HG3	2:B:213:SER:CB	1.87	1.02
2:B:273:LYS:HD2	2:B:273:LYS:H	1.47	0.79
2:B:210:GLU:CG	2:B:213:SER:HB3	2.05	0.79
1:D:592:LYS:HA	1:D:592:LYS:HE2	1.72	0.71
1:D:703:LYS:NZ	1:D:705:GLY:HA3	2.07	0.69
1:D:655:SER:HA	1:D:701:ALA:O	1.93	0.67
2:B:186:ARG:NE	2:B:186:ARG:HA	2.09	0.66
2:B:322:ARG:NH1	2:B:340:GLU:OE1	2.21	0.66
1:D:703:LYS:HE2	1:D:705:GLY:H	1.60	0.65
2:B:204:THR:HG22	2:B:209:LEU:HB2	1.80	0.63
1:D:703:LYS:HD3	1:D:705:GLY:H	1.63	0.63
2:B:191:ARG:NH1	2:B:237:PRO:HG2	2.14	0.63
2:B:214:ASP:OD1	2:B:215:PRO:HD2	1.99	0.62
1:D:703:LYS:HD3	1:D:705:GLY:N	2.15	0.62
2:B:249:GLY:HA3	2:B:354:LYS:HD3	1.82	0.61
1:D:703:LYS:CE	1:D:705:GLY:H	2.14	0.61
2:C:264:HIS:O	2:C:268:GLN:HG2	2.01	0.61
1:A:764:ARG:NH2	1:A:798:ASP:OD2	2.33	0.60
1:D:662:LYS:HG3	1:D:666:GLU:HB2	1.83	0.60
1:D:594:LEU:O	1:D:847:LYS:HD3	2.02	0.60
1:D:592:LYS:HE3	1:D:598:VAL:HG23	1.84	0.60
2:C:317:LEU:HG	2:C:346:LEU:HD22	1.83	0.59
2:B:184:VAL:HG12	2:B:187:ARG:HE	1.68	0.59
1:A:819:SER:HA	1:D:801:TRP:CE2	2.38	0.58
2:B:214:ASP:CG	2:B:215:PRO:HD2	2.28	0.58
2:B:305:GLU:HG3	2:B:306:PRO:HD2	1.86	0.58
1:D:644:LEU:HA	1:D:693:PHE:O	2.03	0.57
1:D:634:LYS:HD3	1:D:638:GLU:OE1	2.04	0.57
2:C:311:ASP:HB2	2:C:363:LEU:HD21	1.86	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:210:GLU:O	2:B:213:SER:HB3	2.05	0.56
2:B:253:ASP:OD2	2:B:253:ASP:N	2.38	0.56
1:D:840:ARG:HD3	5:D:1046:HOH:O	2.03	0.56
2:B:283:MET:HE3	2:B:327:ILE:HD12	1.87	0.56
2:C:254:ARG:HB3	2:C:255:PRO:CD	2.35	0.56
2:B:372:ARG:HG2	2:B:372:ARG:HH11	1.71	0.56
1:D:703:LYS:HZ3	1:D:705:GLY:HA3	1.69	0.56
1:D:656:ASN:OD1	1:D:661:ARG:NH1	2.39	0.56
2:C:256:PRO:HB2	2:C:290:LEU:HD23	1.89	0.55
2:B:187:ARG:NH1	2:B:373:GLU:O	2.40	0.55
1:D:703:LYS:CD	1:D:705:GLY:H	2.20	0.54
1:D:823:ARG:O	1:D:827:GLN:HG2	2.07	0.54
1:D:764:ARG:NH1	1:D:804:GLU:OE1	2.41	0.54
1:A:644:LEU:HA	1:A:693:PHE:O	2.07	0.54
1:A:617:LYS:HE3	1:A:815:TYR:HA	1.90	0.53
1:D:762:TYR:CE1	1:D:763:ASN:HB2	2.43	0.53
2:B:273:LYS:H	2:B:273:LYS:CD	2.21	0.53
2:C:306:PRO:HB3	2:C:324:TRP:CE2	2.44	0.53
2:C:260:LEU:HD21	2:C:298:ALA:HA	1.91	0.52
1:A:772:ARG:NH2	5:A:1001:HOH:O	2.43	0.52
2:C:253:ASP:C	2:C:254:ARG:HG3	2.34	0.52
1:D:662:LYS:CG	1:D:666:GLU:HB2	2.39	0.52
2:C:290:LEU:HD22	2:C:294:ASP:HB3	1.92	0.51
1:A:730:HIS:CE1	1:A:772:ARG:HE	2.28	0.51
2:B:210:GLU:O	2:B:213:SER:CB	2.58	0.51
2:B:347:ALA:O	2:B:351:GLN:HG2	2.09	0.51
2:B:209:LEU:HD21	2:B:218:LEU:HB2	1.92	0.51
1:A:654:LEU:HD11	1:A:669:GLY:HA2	1.93	0.50
2:C:285:VAL:HG11	2:C:359:TRP:HH2	1.76	0.50
1:A:801:TRP:CE2	1:D:819:SER:HA	2.47	0.50
2:B:292:LYS:HA	2:B:295:LEU:HD12	1.93	0.50
2:B:200:LYS:O	2:B:204:THR:HG23	2.11	0.49
1:A:580:LEU:HD22	1:A:628:VAL:HG23	1.93	0.49
2:C:258:TRP:CE2	2:C:262:GLN:HG3	2.46	0.49
2:C:365:LYS:HG3	2:C:366:ASN:N	2.27	0.49
1:A:694:TRP:CH2	1:A:737:GLY:HA2	2.46	0.49
1:A:590:VAL:HG21	1:A:839:ILE:HB	1.95	0.49
1:A:836:VAL:N	1:A:837:PRO:HD2	2.28	0.48
3:D:901:SAH:N	5:D:1001:HOH:O	2.35	0.47
1:A:756:LEU:HD21	1:A:760:LEU:HD21	1.97	0.47
1:A:724:ILE:HG13	1:A:770:LYS:HD2	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:739:LEU:O	1:D:742:MET:HG3	2.14	0.47
1:D:755:GLU:HG2	1:D:769:LYS:HE2	1.96	0.47
2:B:306:PRO:HB3	2:B:324:TRP:CE2	2.50	0.46
1:A:828:LYS:O	1:A:832:ARG:HG3	2.15	0.46
1:D:693:PHE:CD1	1:D:740:PRO:HD3	2.50	0.46
1:D:626:ASN:HB2	5:D:1010:HOH:O	2.15	0.46
1:A:586:THR:HG23	1:A:836:VAL:HG12	1.98	0.46
1:A:626:ASN:HB3	1:A:627:ASP:H	1.54	0.46
2:C:253:ASP:C	2:C:254:ARG:CG	2.89	0.46
2:C:376:LYS:HG2	2:C:378:PHE:CE1	2.51	0.46
1:A:820:ASN:OD1	1:D:801:TRP:HA	2.16	0.45
1:A:844:ALA:O	1:A:847:LYS:HG3	2.17	0.45
2:B:339:SER:HB3	2:B:342:GLU:CD	2.42	0.45
1:D:654:LEU:HD13	1:D:698:ASN:OD1	2.16	0.45
1:D:823:ARG:NH1	1:D:827:GLN:OE1	2.49	0.45
1:A:661:ARG:CZ	1:A:703:LYS:HE3	2.46	0.45
2:C:254:ARG:CB	2:C:255:PRO:CD	2.95	0.44
2:C:258:TRP:HD1	1:D:665:TYR:CZ	2.35	0.44
1:A:730:HIS:ND1	1:A:772:ARG:HG3	2.31	0.44
2:B:246:PRO:HG2	2:B:252:CYS:SG	2.58	0.44
2:B:346:LEU:HD12	2:B:350:LYS:HB2	1.99	0.44
2:B:190:VAL:HG22	2:B:375:PHE:CG	2.52	0.44
1:A:759:CYS:HB3	1:A:807:ARG:HB2	2.00	0.44
1:D:718:PRO:HB3	1:D:736:TRP:CD2	2.53	0.44
2:B:256:PRO:HB3	2:B:289:VAL:HG12	2.00	0.43
2:B:210:GLU:HB3	2:B:378:PHE:CG	2.53	0.43
2:B:247:PRO:HG3	2:B:359:TRP:CD1	2.53	0.43
1:D:724:ILE:HG13	1:D:770:LYS:HE3	2.01	0.43
1:A:655:SER:HA	1:A:701:ALA:O	2.19	0.43
1:D:617:LYS:HE3	1:D:815:TYR:HA	2.01	0.42
2:B:361:THR:HA	2:B:364:VAL:HG12	2.01	0.42
1:A:702:MET:HB2	1:A:702:MET:HE3	1.77	0.42
1:D:724:ILE:HG13	1:D:770:LYS:CE	2.49	0.42
1:D:617:LYS:HD3	1:D:617:LYS:HA	1.77	0.42
1:D:588:TYR:CE1	1:D:592:LYS:HG3	2.55	0.42
2:B:196:PHE:O	2:B:222:VAL:HG12	2.20	0.42
2:C:273:LYS:HA	2:C:273:LYS:HD2	1.67	0.42
2:B:258:TRP:CE2	2:B:262:GLN:HG3	2.54	0.42
2:C:242:TYR:CE1	2:C:283:MET:HE2	2.55	0.42
1:A:645:VAL:O	1:A:694:TRP:HA	2.20	0.41
2:B:216:GLY:O	2:B:217:GLN:C	2.60	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:304:MET:HE3	2:B:304:MET:HB2	1.79	0.41
2:B:367:CYS:O	2:B:370:PRO:HD2	2.20	0.41
1:D:787:GLN:HG2	1:D:788:LEU:H	1.85	0.41
1:A:661:ARG:H	1:A:661:ARG:HG3	1.66	0.41
1:A:699:VAL:HG12	1:A:701:ALA:H	1.84	0.41
2:C:285:VAL:HG11	2:C:364:VAL:HG11	2.01	0.41
1:A:702:MET:HE1	1:A:707:LYS:HA	2.03	0.41
1:A:801:TRP:CD2	1:D:819:SER:HA	2.56	0.41
2:B:339:SER:OG	2:B:340:GLU:N	2.53	0.41
2:B:311:ASP:HB2	2:B:363:LEU:HD21	2.02	0.40
1:D:595:GLY:O	1:D:852:CYS:HB3	2.21	0.40
1:D:765:ILE:O	1:D:791:VAL:HG23	2.21	0.40
2:C:242:TYR:HE1	2:C:283:MET:HE2	1.87	0.40
2:C:353:SER:HB2	2:C:355:LEU:HD13	2.03	0.40
2:B:185:TRP:CE3	2:B:185:TRP:HA	2.55	0.40
1:D:581:PHE:HB3	3:D:901:SAH:O4'	2.22	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	268/286 (94%)	259 (97%)	9 (3%)	0	100	100
1	D	269/286 (94%)	257 (96%)	12 (4%)	0	100	100
2	B	194/204 (95%)	186 (96%)	8 (4%)	0	100	100
2	C	196/204 (96%)	189 (96%)	7 (4%)	0	100	100
All	All	927/980 (95%)	891 (96%)	36 (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	237/249 (95%)	237 (100%)	0	100	100
1	D	237/249 (95%)	237 (100%)	0	100	100
2	B	182/185 (98%)	182 (100%)	0	100	100
2	C	182/185 (98%)	182 (100%)	0	100	100
All	All	838/868 (96%)	838 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	349	ASN
2	B	366	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](#)

4 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	SAH	A	901	-	27,28,28	1.15	4 (14%)	36,40,40	2.21	9 (25%)
4	FMT	D	902	-	2,2,2	0.65	0	1,1,1	0.21	0
3	SAH	D	901	-	27,28,28	1.13	3 (11%)	36,40,40	2.07	12 (33%)
4	FMT	A	902	-	2,2,2	0.61	0	1,1,1	0.11	0

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	SAH	A	901	-	-	6/15/31/31	0/3/3/3
3	SAH	D	901	-	-	4/15/31/31	0/3/3/3

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	901	SAH	C2-N3	2.94	1.39	1.33
3	D	901	SAH	C2-N1	2.88	1.39	1.33
3	D	901	SAH	C2-N3	2.69	1.38	1.33
3	D	901	SAH	C8-N7	2.44	1.36	1.31
3	A	901	SAH	C2-N1	2.32	1.38	1.33
3	A	901	SAH	C8-N7	2.21	1.35	1.31
3	A	901	SAH	OXT-C	-2.08	1.24	1.30

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	901	SAH	N3-C2-N1	-5.75	119.88	128.58
3	A	901	SAH	C5-C4-N3	-5.60	119.01	126.72
3	D	901	SAH	C5-C4-N3	-5.40	119.28	126.72
3	D	901	SAH	N3-C2-N1	-4.85	121.25	128.58
3	A	901	SAH	C2-N3-C4	4.15	121.97	111.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	901	SAH	C5-N7-C8	3.88	109.55	103.45
3	A	901	SAH	N9-C8-N7	-3.85	108.48	113.94
3	A	901	SAH	OXT-C-O	-3.79	115.48	124.08
3	A	901	SAH	N3-C4-N9	3.56	133.23	127.17
3	D	901	SAH	C2-N3-C4	3.52	120.42	111.83
3	D	901	SAH	C5-N7-C8	3.32	108.66	103.45
3	D	901	SAH	OXT-C-O	-3.29	116.61	124.08
3	D	901	SAH	N3-C4-N9	3.26	132.71	127.17
3	D	901	SAH	N9-C8-N7	-3.26	109.31	113.94
3	A	901	SAH	C4-C5-N7	-2.95	107.21	110.58
3	D	901	SAH	C4-C5-N7	-2.67	107.53	110.58
3	D	901	SAH	C6-C5-C4	2.59	120.71	117.18
3	A	901	SAH	C5'-SD-CG	-2.25	95.59	102.26
3	D	901	SAH	C5'-SD-CG	-2.20	95.74	102.26
3	D	901	SAH	O4'-C1'-N9	2.04	112.00	108.09
3	D	901	SAH	C5-C4-N9	2.02	108.01	105.81

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	901	SAH	N-CA-CB-CG
3	D	901	SAH	N-CA-CB-CG
3	D	901	SAH	C2'-C1'-N9-C8
3	A	901	SAH	O-C-CA-CB
3	A	901	SAH	C-CA-CB-CG
3	D	901	SAH	C-CA-CB-CG
3	A	901	SAH	CB-CG-SD-C5'
3	A	901	SAH	OXT-C-CA-CB
3	A	901	SAH	C2'-C1'-N9-C8
3	D	901	SAH	O4'-C1'-N9-C8

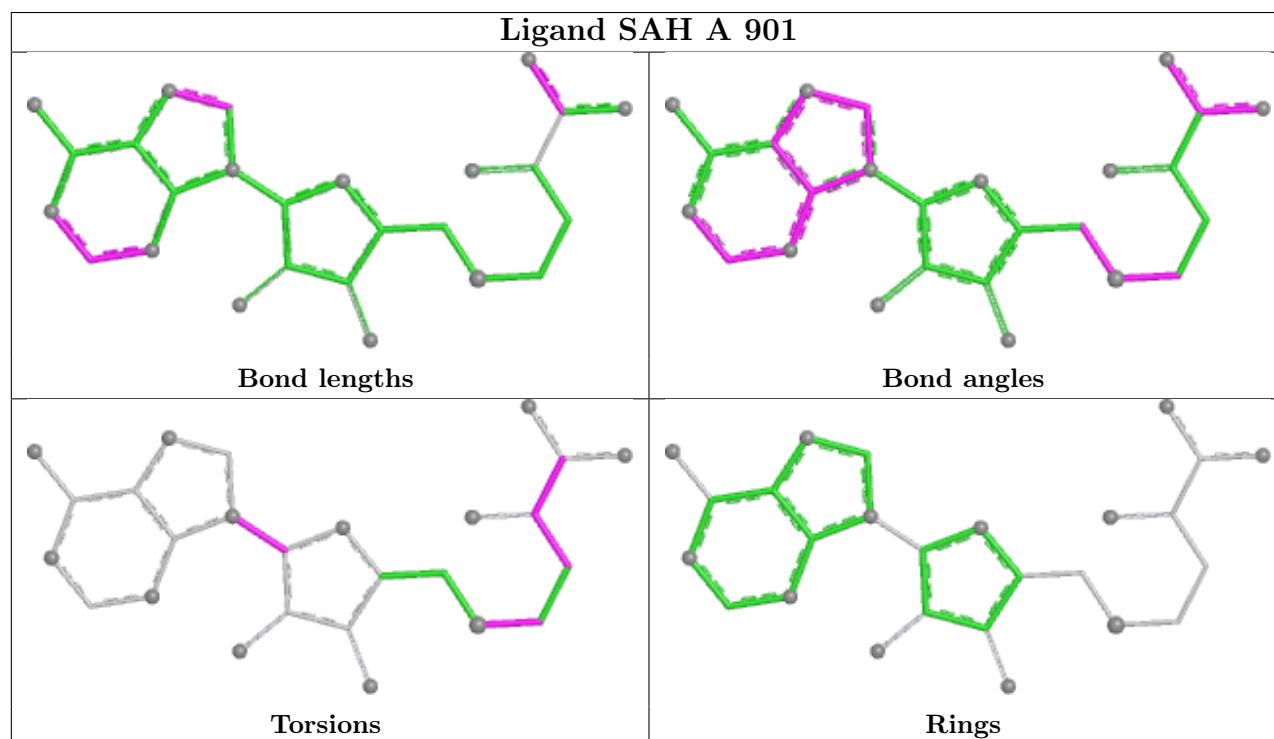
There are no ring outliers.

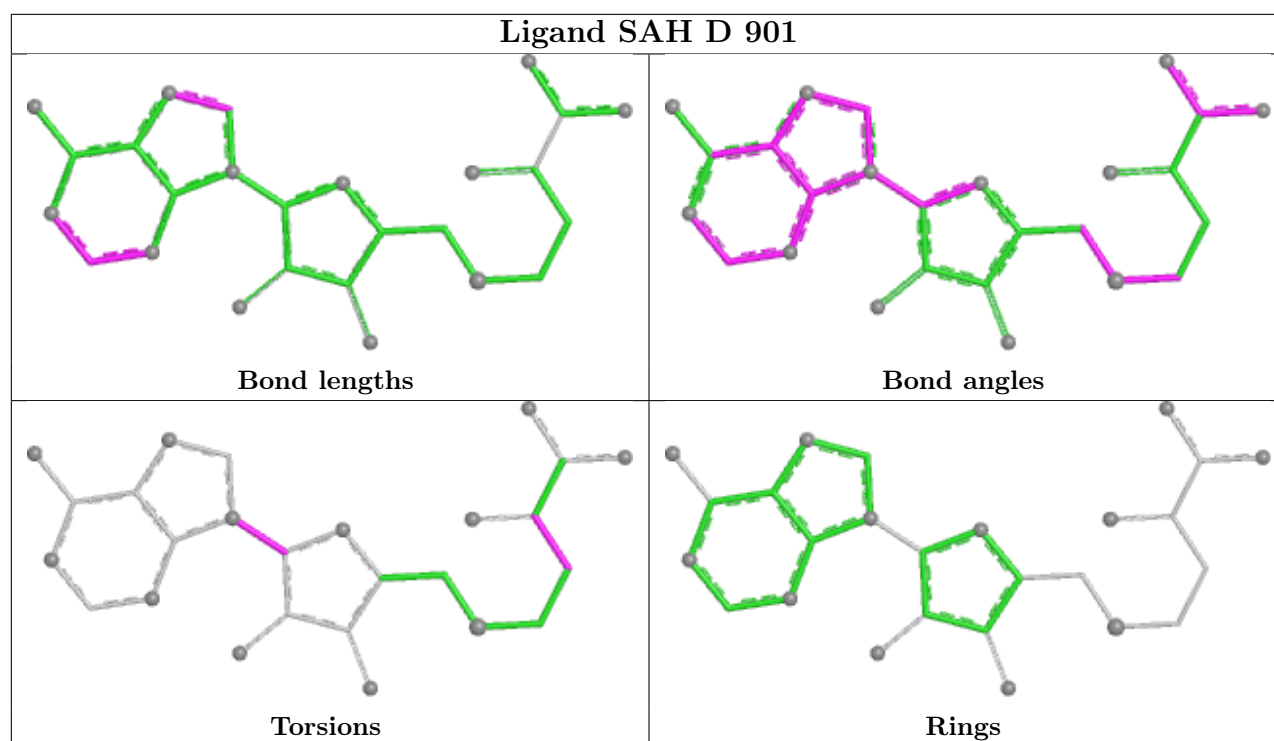
1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	901	SAH	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	272/286 (95%)	-0.37	3 (1%) 78 72	28, 45, 97, 143	0
1	D	273/286 (95%)	-0.32	7 (2%) 57 48	26, 42, 98, 133	0
2	B	198/204 (97%)	0.22	11 (5%) 30 23	34, 58, 133, 184	0
2	C	200/204 (98%)	-0.01	10 (5%) 34 27	28, 53, 128, 153	0
All	All	943/980 (96%)	-0.15	31 (3%) 49 40	26, 48, 115, 184	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	215	PRO	6.0
1	D	788	LEU	3.7
2	B	249	GLY	3.6
1	A	773	THR	3.6
1	D	773	THR	3.6
2	B	185	TRP	3.4
2	C	252	CYS	3.1
1	D	657	VAL	3.1
2	B	350	LYS	2.9
2	C	213	SER	2.9
2	C	356	ALA	2.7
1	D	569	HIS	2.7
2	B	250	HIS	2.6
2	B	351	GLN	2.6
2	C	182	VAL	2.6
2	B	216	GLY	2.5
1	A	788	LEU	2.4
2	B	179	PHE	2.4
1	D	656	ASN	2.4
2	C	373	GLU	2.4
1	A	655	SER	2.3

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Mol	Chain	Res	Type	RSRZ
2	C	181	THR	2.3
2	C	251	THR	2.3
2	C	178	MET	2.3
1	D	703	LYS	2.2
1	D	652	ASN	2.2
2	B	217	GLN	2.2
2	B	181	THR	2.2
2	B	273	LYS	2.1
2	C	354	LYS	2.1
2	C	212	GLY	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](#)

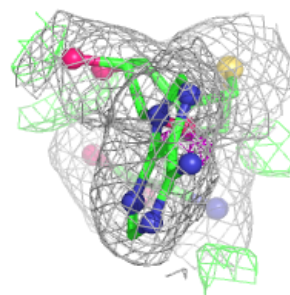
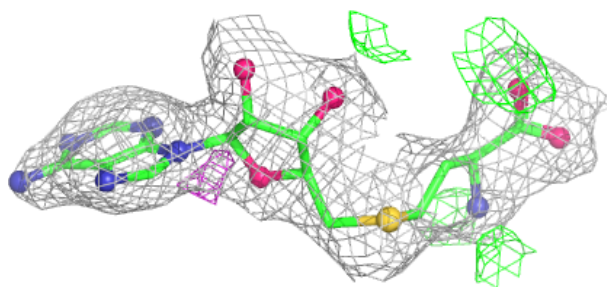
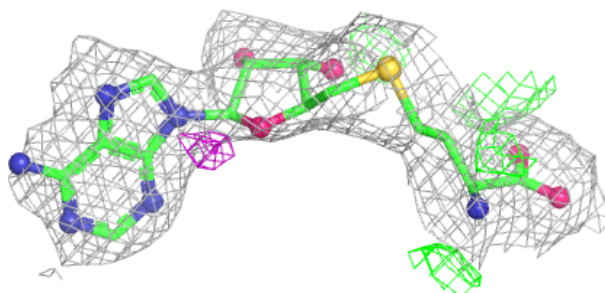
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	FMT	A	902	3/3	0.75	0.23	55,55,55,57	0
4	FMT	D	902	3/3	0.86	0.14	52,52,53,58	0
3	SAH	A	901	26/26	0.90	0.11	39,59,73,76	0
3	SAH	D	901	26/26	0.92	0.09	38,52,68,69	0

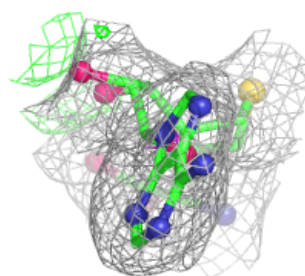
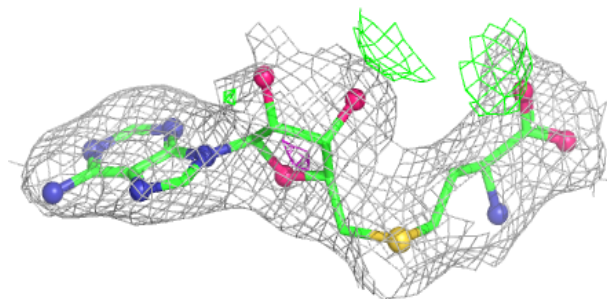
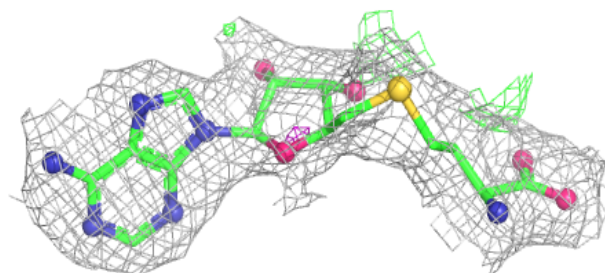
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around SAH A 901:

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

**Electron density around SAH D 901:**

$2mF_o-DF_c$ (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.