



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 03:18 AM UTC

PDB ID : 4QEP / pdb_00004qep
Title : crystal structure of KRYPTONITE in complex with mCHG DNA and SAH
Authors : Du, J.; Li, S.; Patel, D.J.
Deposited on : 2014-05-17
Resolution : 3.10 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

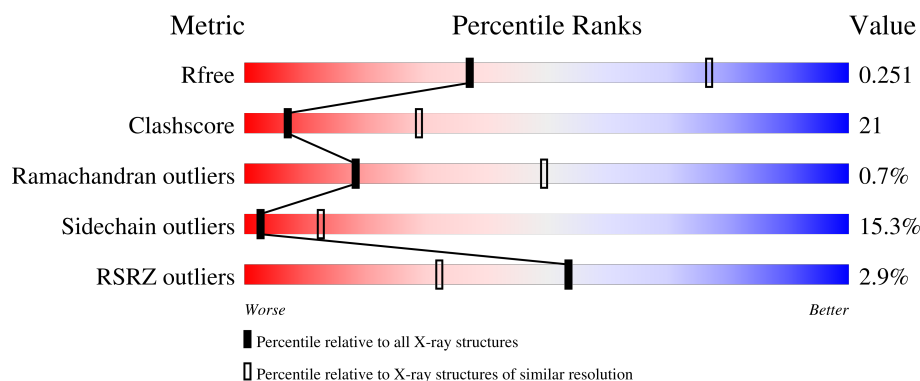
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.10 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	1456 (3.10-3.10)
Clashscore	190562	1539 (3.10-3.10)
Ramachandran outliers	187476	1467 (3.10-3.10)
Sidechain outliers	187428	1467 (3.10-3.10)
RSRZ outliers	180081	1456 (3.10-3.10)

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	533	2% 48% 32% 7% 12%
2	C	15	7% 20% 40% 13% 27%
3	D	15	7% 20% 40% 27% 13%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	SAH	A	801	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 4210 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 Histone-lysine N-methyltransferase, H3 lysine-9 specific SUVH4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	467	Total	C	N	O	S	0	0	0
			3690	2332	647	681	30			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	92	SER	-	expression tag	UNP Q8GZB6

- Molecule 2 is a DNA chain called DNA (5'-D(*GP*GP*TP*AP*CP*TP*(5CM)P*AP*GP*CP*AP*GP*TP*AP*T)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	11	Total	C	N	O	P	0	0	0
			224	108	43	63	10			

- Molecule 3 is a DNA chain called DNA (5'-D(*AP*CP*TP*GP*CP*TP*GP*AP*GP*TP*AP*CP*CP*AP*T)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	13	Total	C	N	O	P	0	0	0
			266	127	47	79	13			

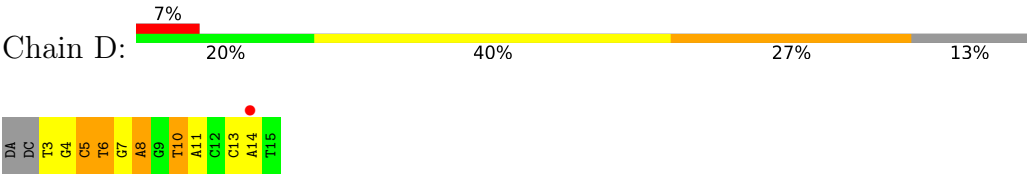
- Molecule 4 is S-ADENOSYL-L-HOMOCYSTEINE (CCD ID: SAH) (formula: C₁₄H₂₀N₆O₅S).



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

- Molecule 5 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	4	Total	Zn	0	0
			4	4		



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	54.35Å 95.57Å 121.75Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.63 – 3.10 49.63 – 3.10	Depositor EDS
% Data completeness (in resolution range)	89.8 (49.63-3.10) 89.8 (49.63-3.10)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	0.15	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.29 (at 3.12Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.2_1309)	Depositor
R, R_{free}	0.210 , 0.251 0.212 , 0.251	Depositor DCC
R_{free} test set	517 reflections (4.27%)	wwPDB-VP
Wilson B-factor (Å ²)	58.5	Xtriage
Anisotropy	0.631	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 50.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	4210	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.93% 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, 5CM, ZN

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	0.95	6/3767 (0.2%)	1.01	17/5082 (0.3%)
2	C	0.89	2/228 (0.9%)	1.22	4/348 (1.1%)
3	D	0.89	4/297 (1.3%)	1.03	2/456 (0.4%)
All	All	0.94	12/4292 (0.3%)	1.03	23/5886 (0.4%)

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	5	DC	O3'-P	-8.75	1.48	1.61
2	C	4	DA	O3'-P	-6.39	1.51	1.61
2	C	5	DC	O3'-P	-6.13	1.51	1.61
1	A	619	CYS	CA-C	-6.06	1.45	1.52
3	D	7	DG	O3'-P	-5.78	1.52	1.61
1	A	143	TYR	C-O	-5.69	1.19	1.24
1	A	269	THR	CA-CB	-5.63	1.45	1.53
1	A	299	ARG	C-O	-5.58	1.17	1.24
1	A	220	THR	CA-CB	-5.42	1.45	1.53
3	D	6	DT	O3'-P	-5.37	1.53	1.61
3	D	8	DA	O3'-P	-5.31	1.53	1.61
1	A	241	ARG	C-O	-5.00	1.17	1.24

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	5	DC	O3'-P-O5'	-8.80	90.80	104.00
2	C	2	DG	O3'-P-O5'	-8.20	91.70	104.00
1	A	345	ILE	CA-C-N	7.56	127.32	119.76
1	A	345	ILE	C-N-CA	7.56	127.32	119.76
1	A	615	GLY	N-CA-C	-6.41	105.66	114.64
1	A	177	ASN	N-CA-C	6.40	119.64	110.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	600	VAL	CB-CA-C	-6.25	101.51	110.83
1	A	339	GLY	N-CA-C	-6.23	106.55	115.64
2	C	2	DG	C2'-C3'-O3'	-6.22	102.16	111.50
1	A	614	CYS	CA-C-N	6.21	132.57	122.69
1	A	614	CYS	C-N-CA	6.21	132.57	122.69
3	D	6	DT	C4'-C3'-O3'	-6.00	100.99	110.00
1	A	143	TYR	CA-C-N	5.88	127.19	119.84
1	A	143	TYR	C-N-CA	5.88	127.19	119.84
1	A	397	CYS	CA-CB-SG	5.83	127.81	114.40
1	A	623	LEU	N-CA-C	-5.80	104.33	111.40
1	A	350	ARG	N-CA-C	-5.53	106.38	114.39
1	A	432	GLY	CA-C-N	5.40	125.02	119.56
1	A	432	GLY	C-N-CA	5.40	125.02	119.56
1	A	457	GLY	N-CA-C	-5.38	100.42	113.18
2	C	10	DC	P-O5'-C5'	-5.28	112.08	120.00
1	A	620	ARG	N-CA-C	-5.22	106.91	113.28
3	D	10	DT	C2'-C3'-O3'	5.02	119.03	111.50

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	3690	0	3641	148	0
2	C	224	0	127	5	0
3	D	266	0	148	14	0
4	A	26	0	19	9	0
5	A	4	0	0	0	0
All	All	4210	0	3935	168	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 21.

All (168) 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:554:CYS:N	1:A:614:CYS:SG	2.28	1.05
1:A:459:ALA:HB2	1:A:551:ASN:ND2	1.77	0.99
1:A:334:GLU:N	1:A:334:GLU:OE1	1.96	0.96
1:A:100:GLU:HB3	1:A:103:LEU:HD13	1.55	0.89
1:A:553:SER:HA	1:A:614:CYS:SG	2.15	0.87
3:D:4:DG:H8	3:D:4:DG:O5'	1.58	0.85
4:A:801:SAH:O4'	4:A:801:SAH:HB2	1.78	0.84
4:A:801:SAH:H8	4:A:801:SAH:H5'2	1.60	0.84
1:A:600:VAL:HG21	1:A:624:TYR:CE1	2.12	0.83
1:A:459:ALA:CB	1:A:551:ASN:HD22	1.94	0.81
1:A:600:VAL:HG21	1:A:624:TYR:CD1	2.15	0.80
1:A:228:THR:OG1	1:A:229:GLY:N	2.12	0.79
1:A:459:ALA:CB	1:A:551:ASN:ND2	2.44	0.79
1:A:600:VAL:O	1:A:600:VAL:HG23	1.85	0.75
1:A:100:GLU:OE1	1:A:103:LEU:CD1	2.35	0.74
1:A:456:LYS:NZ	1:A:588:GLU:OE2	2.21	0.74
1:A:419:ARG:O	1:A:561:GLN:NE2	2.21	0.74
1:A:100:GLU:OE1	1:A:103:LEU:HD13	1.90	0.71
1:A:204:SER:OG	1:A:271:ARG:NH1	2.24	0.70
2:C:9:DG:H2''	2:C:10:DC:H5''	1.72	0.69
1:A:100:GLU:CD	1:A:101:PRO:HD2	2.16	0.69
1:A:456:LYS:NZ	1:A:588:GLU:CD	2.51	0.69
1:A:459:ALA:HB2	1:A:551:ASN:HD21	1.57	0.69
4:A:801:SAH:H8	4:A:801:SAH:C5'	2.24	0.68
1:A:458:TRP:HB2	4:A:801:SAH:OXT	1.94	0.68
3:D:4:DG:H2''	3:D:5:DC:O5'	1.93	0.68
1:A:100:GLU:CG	1:A:101:PRO:HD2	2.24	0.68
1:A:124:LEU:HD23	1:A:130:LYS:HE2	1.76	0.68
1:A:535:GLU:OE1	1:A:571:ARG:NH1	2.27	0.68
1:A:597:LEU:HD21	1:A:622:ARG:HD3	1.76	0.67
1:A:397:CYS:SG	1:A:425:CYS:HA	2.36	0.66
1:A:194:LYS:HG2	1:A:252:TYR:OH	1.97	0.65
1:A:271:ARG:NH2	1:A:273:TYR:OH	2.30	0.65
1:A:553:SER:CA	1:A:614:CYS:SG	2.85	0.64
1:A:455:LYS:O	1:A:455:LYS:HG3	1.97	0.64
1:A:619:CYS:SG	1:A:621:LYS:N	2.68	0.64
3:D:4:DG:O5'	3:D:4:DG:C8	2.47	0.63
1:A:257:ARG:HB3	1:A:275:TYR:HE1	1.63	0.63
1:A:456:LYS:NZ	1:A:588:GLU:OE1	2.32	0.62
1:A:125:SER:O	1:A:130:LYS:NZ	2.30	0.62
1:A:103:LEU:HD12	1:A:103:LEU:N	2.15	0.62
1:A:450:VAL:HG22	1:A:460:VAL:HG12	1.81	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:600:VAL:CG2	1:A:624:TYR:CD1	2.83	0.61
1:A:597:LEU:O	1:A:598:ASP:HB2	2.00	0.61
1:A:540:ALA:O	1:A:548:ARG:NH2	2.35	0.60
1:A:329:GLU:HG3	1:A:331:LEU:H	1.66	0.60
1:A:104:LYS:O	1:A:107:LYS:HG2	2.02	0.60
1:A:282:GLU:OE1	1:A:299:ARG:NH1	2.35	0.59
1:A:499:CYS:SG	1:A:562:CYS:HB2	2.42	0.59
1:A:412:ASP:O	1:A:445:ARG:NH2	2.28	0.59
1:A:100:GLU:HG2	1:A:101:PRO:HD2	1.84	0.59
1:A:184:MET:O	1:A:187:GLU:HG2	2.03	0.58
1:A:221:GLY:HA3	1:A:243:ASN:OD1	2.02	0.58
1:A:459:ALA:HB1	1:A:551:ASN:HD22	1.69	0.58
1:A:617:LEU:C	1:A:618:ASN:HD22	2.13	0.57
1:A:99:LEU:O	1:A:104:LYS:NZ	2.26	0.57
1:A:583:ILE:HG23	1:A:587:GLN:HE21	1.68	0.57
1:A:452:ARG:HH21	1:A:457:GLY:HA2	1.68	0.56
1:A:194:LYS:CG	1:A:252:TYR:OH	2.53	0.56
1:A:458:TRP:O	4:A:801:SAH:OXT	2.22	0.56
1:A:462:SER:O	1:A:586:MET:N	2.36	0.55
1:A:156:VAL:HA	1:A:275:TYR:HD2	1.72	0.55
1:A:607:VAL:HG13	1:A:608:LYS:N	2.22	0.54
3:D:5:DC:H4'	3:D:6:DT:OP1	2.07	0.54
1:A:156:VAL:HA	1:A:275:TYR:CD2	2.43	0.53
1:A:552:HIS:HB3	4:A:801:SAH:HN61	1.72	0.53
1:A:553:SER:C	1:A:614:CYS:SG	2.90	0.53
1:A:288:LYS:NZ	1:A:294:THR:OG1	2.41	0.53
2:C:11:DA:N1	3:D:4:DG:C2	2.77	0.52
1:A:224:GLY:HA2	1:A:235:LYS:HG2	1.91	0.52
1:A:395:CYS:O	1:A:399:LYS:HG3	2.10	0.52
1:A:113:ASN:ND2	1:A:468:ALA:HB3	2.26	0.51
1:A:226:ASN:O	1:A:231:LYS:HA	2.10	0.51
1:A:416:ILE:HG22	1:A:417:GLU:HG2	1.92	0.51
1:A:453:SER:OG	1:A:457:GLY:O	2.20	0.51
1:A:118:LEU:HD23	1:A:416:ILE:HD13	1.93	0.51
1:A:154:ILE:HD11	1:A:169:VAL:HG21	1.93	0.50
1:A:455:LYS:O	1:A:455:LYS:CG	2.59	0.50
1:A:100:GLU:OE2	1:A:101:PRO:HD2	2.11	0.50
1:A:458:TRP:N	4:A:801:SAH:OXT	2.45	0.49
1:A:374:VAL:HG22	1:A:481:ARG:HG3	1.94	0.49
1:A:600:VAL:O	1:A:600:VAL:CG2	2.55	0.49
1:A:132:VAL:O	1:A:136:ILE:HG13	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:408:VAL:HG12	1:A:416:ILE:HG12	1.95	0.49
2:C:10:DC:O2	3:D:4:DG:N2	2.24	0.49
1:A:214:ALA:HA	1:A:260:ARG:NH2	2.27	0.49
1:A:217:VAL:HB	1:A:300:LEU:HD12	1.95	0.48
1:A:240:GLU:O	1:A:241:ARG:HG3	2.14	0.48
1:A:280:LYS:O	1:A:300:LEU:HA	2.14	0.48
1:A:390:THR:OG1	1:A:391:ASP:N	2.45	0.48
1:A:165:GLU:O	1:A:169:VAL:HG12	2.13	0.48
1:A:174:HIS:CB	1:A:177:ASN:HB2	2.44	0.48
1:A:302:ARG:CZ	1:A:309:LEU:HD12	2.43	0.48
1:A:415:LEU:HD23	1:A:559:PHE:HZ	1.79	0.48
3:D:5:DC:C4'	3:D:6:DT:OP1	2.61	0.48
3:D:3:DT:H2'	3:D:3:DT:O2	2.12	0.48
1:A:353:ASP:O	1:A:355:PRO:HD3	2.14	0.47
1:A:499:CYS:SG	1:A:570:ILE:HD11	2.54	0.47
1:A:198:ALA:N	1:A:245:ALA:HB1	2.30	0.47
1:A:345:ILE:HG12	1:A:448:LEU:HB2	1.97	0.47
1:A:453:SER:OG	1:A:456:LYS:HB2	2.14	0.46
1:A:557:ASN:O	1:A:558:LEU:HD23	2.16	0.46
1:A:345:ILE:HA	1:A:346:PRO:HD2	1.78	0.46
1:A:329:GLU:HG3	1:A:330:GLY:N	2.30	0.46
1:A:601:HIS:ND1	1:A:607:VAL:HA	2.31	0.45
1:A:439:THR:HG22	1:A:471:PRO:HG3	1.98	0.45
1:A:335:ASP:OD1	1:A:345:ILE:N	2.40	0.45
3:D:3:DT:OP1	3:D:3:DT:C7	2.64	0.45
1:A:266:SER:O	1:A:267:SER:C	2.60	0.45
1:A:332:VAL:HG12	1:A:348:THR:HA	1.97	0.45
4:A:801:SAH:C8	4:A:801:SAH:C3'	2.94	0.45
1:A:545:ASN:O	1:A:548:ARG:HG2	2.16	0.45
1:A:103:LEU:CD1	1:A:103:LEU:N	2.79	0.45
1:A:332:VAL:CG1	1:A:347:ALA:O	2.66	0.44
1:A:614:CYS:HB3	1:A:616:ALA:H	1.83	0.44
1:A:116:TYR:O	1:A:120:VAL:HG13	2.17	0.44
1:A:280:LYS:HE3	1:A:280:LYS:HB2	1.76	0.44
1:A:443:ARG:HE	1:A:443:ARG:HB3	1.60	0.44
1:A:619:CYS:SG	1:A:620:ARG:N	2.91	0.44
1:A:208:GLU:H	1:A:208:GLU:HG3	1.36	0.44
1:A:583:ILE:HG21	1:A:589:LEU:HD21	1.98	0.44
1:A:100:GLU:OE2	1:A:101:PRO:HG2	2.18	0.44
1:A:178:GLY:O	1:A:201:ILE:HA	2.18	0.44
1:A:612:CYS:HB3	1:A:619:CYS:SG	2.58	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:597:LEU:HD23	1:A:624:TYR:O	2.17	0.43
1:A:349:ASN:ND2	1:A:452:ARG:HD3	2.33	0.43
1:A:197:LEU:HD21	1:A:254:VAL:HG21	2.00	0.43
1:A:336:ILE:HD13	1:A:450:VAL:HG21	2.00	0.43
4:A:801:SAH:C8	4:A:801:SAH:H3'	2.49	0.43
1:A:247:LYS:HG3	1:A:284:PHE:CZ	2.54	0.42
1:A:401:ASN:CG	1:A:406:PRO:HG3	2.44	0.42
1:A:623:LEU:HA	1:A:623:LEU:HD12	1.86	0.42
1:A:264:CYS:HB2	1:A:269:THR:O	2.20	0.42
1:A:347:ALA:HA	1:A:450:VAL:O	2.19	0.42
1:A:606:LYS:HG3	1:A:607:VAL:HB	2.01	0.42
3:D:13:DC:H2''	3:D:14:DA:N7	2.34	0.42
1:A:129:LEU:HB3	3:D:8:DA:H5'	2.02	0.42
1:A:279:TYR:CE1	1:A:302:ARG:HG3	2.55	0.42
1:A:349:ASN:HD22	1:A:452:ARG:HD3	1.85	0.42
1:A:375:ILE:HD13	1:A:536:PHE:HZ	1.85	0.42
1:A:227:LEU:HD12	1:A:227:LEU:H	1.85	0.41
1:A:104:LYS:HA	1:A:107:LYS:HE3	2.02	0.41
1:A:149:GLY:O	1:A:257:ARG:HD3	2.20	0.41
1:A:243:ASN:HB3	1:A:298:TYR:OH	2.21	0.41
1:A:104:LYS:O	1:A:107:LYS:CG	2.69	0.41
1:A:332:VAL:HG13	1:A:347:ALA:O	2.21	0.41
3:D:10:DT:H2'	3:D:11:DA:C8	2.55	0.41
1:A:115:GLN:H	1:A:115:GLN:HG2	1.66	0.41
1:A:610:LEU:HD23	1:A:610:LEU:HA	1.68	0.41
1:A:146:LYS:HE3	1:A:181:TYR:O	2.20	0.41
1:A:613:TYR:O	1:A:614:CYS:C	2.64	0.41
1:A:618:ASN:O	1:A:619:CYS:C	2.63	0.41
2:C:7:5CM:O5'	2:C:7:5CM:H2'	2.21	0.41
1:A:459:ALA:HA	1:A:549:PHE:O	2.21	0.40
1:A:466:ILE:HB	1:A:583:ILE:HB	2.03	0.40
1:A:196:PRO:O	1:A:245:ALA:HA	2.21	0.40
1:A:558:LEU:HD11	1:A:589:LEU:HB3	2.04	0.40
1:A:151:LEU:HA	1:A:151:LEU:HD12	1.80	0.40
1:A:262:HIS:HB2	1:A:272:VAL:HG12	2.01	0.40
1:A:334:GLU:N	1:A:334:GLU:CD	2.75	0.40
1:A:424:GLU:OE2	1:A:440:SER:OG	2.39	0.40
1:A:201:ILE:HD11	1:A:258:VAL:HG22	2.03	0.40
1:A:240:GLU:C	1:A:241:ARG:HG3	2.45	0.40
1:A:244:LEU:HD12	1:A:244:LEU:HA	1.90	0.40
2:C:1:DG:C2	3:D:14:DA:C2	3.10	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:126:ARG:HB3	1:A:129:LEU:HD22	2.04	0.40
1:A:127:PRO:HA	1:A:130:LYS:HB2	2.02	0.40
3:D:6:DT:O5'	3:D:6:DT:H2'	2.20	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	457/533 (86%)	435 (95%)	19 (4%)	3 (1%)	18 49

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	607	VAL
1	A	267	SER
1	A	354	SER

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	406/461 (88%)	344 (85%)	62 (15%)	3 12

All (62) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	99	LEU
1	A	106	THR
1	A	114	LYS
1	A	120	VAL
1	A	121	GLN
1	A	124	LEU
1	A	128	ASP
1	A	129	LEU
1	A	145	ARG
1	A	151	LEU
1	A	169	VAL
1	A	191	SER
1	A	203	MET
1	A	208	GLU
1	A	211	LEU
1	A	215	ASP
1	A	216	THR
1	A	220	THR
1	A	222	GLN
1	A	228	THR
1	A	230	ASN
1	A	259	THR
1	A	267	SER
1	A	270	LYS
1	A	272	VAL
1	A	282	GLU
1	A	283	LYS
1	A	288	LYS
1	A	300	LEU
1	A	304	GLU
1	A	308	GLU
1	A	309	LEU
1	A	331	LEU
1	A	332	VAL
1	A	334	GLU
1	A	356	VAL
1	A	378	LYS
1	A	381	THR
1	A	385	CYS
1	A	391	ASP
1	A	392	SER
1	A	397	CYS
1	A	408	VAL

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Mol	Chain	Res	Type
1	A	422	VAL
1	A	434	LYS
1	A	436	VAL
1	A	442	LYS
1	A	452	ARG
1	A	460	VAL
1	A	550	ILE
1	A	555	GLU
1	A	560	VAL
1	A	566	SER
1	A	568	GLN
1	A	587	GLN
1	A	597	LEU
1	A	607	VAL
1	A	608	LYS
1	A	609	GLN
1	A	610	LEU
1	A	614	CYS
1	A	621	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	113	ASN
1	A	206	GLN
1	A	225	HIS
1	A	230	ASN
1	A	233	GLN
1	A	551	ASN
1	A	568	GLN
1	A	587	GLN
1	A	618	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

1 non-standard protein/DNA/RNA residue is modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
2	5CM	C	7	2	18,21,22	2.40	6 (33%)	24,30,33	4.49	10 (41%)

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	5CM	C	7	2	-	2/7/21/22	0/2/2/2

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	7	5CM	C5-C4	-6.37	1.39	1.44
2	C	7	5CM	C5A-C5	-4.41	1.40	1.50
2	C	7	5CM	C4-N3	3.80	1.40	1.34
2	C	7	5CM	C6-C5	2.97	1.39	1.34
2	C	7	5CM	C2-N1	-2.53	1.34	1.40
2	C	7	5CM	O4'-C4'	2.15	1.49	1.45

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	7	5CM	O4'-C1'-C2'	-16.27	75.85	106.25
2	C	7	5CM	O4'-C4'-C5'	-9.76	78.07	109.33
2	C	7	5CM	O4'-C4'-C3'	-6.92	89.87	105.65
2	C	7	5CM	C5'-C4'-C3'	3.69	135.28	114.64
2	C	7	5CM	C3'-C2'-C1'	-3.59	93.80	102.60
2	C	7	5CM	C5-C4-N4	-3.06	117.08	121.39
2	C	7	5CM	N4-C4-N3	2.59	123.20	118.51
2	C	7	5CM	C5-C6-N1	-2.51	120.58	123.31
2	C	7	5CM	C2'-C3'-C4'	-2.10	98.55	102.80
2	C	7	5CM	O2-C2-N3	-2.00	119.17	122.33

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	7	5CM	C3'-C4'-C5'-O5'
2	C	7	5CM	C2'-C1'-N1-C2

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	7	5CM	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 5 ligands modelled in this entry, 4 are monoatomic - leaving 1 for Mogul analysis.

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
4	SAH	A	801	-	27,28,28	1.66	5 (18%)	36,40,40	2.17	10 (27%)

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
4	SAH	A	801	-	-	7/15/31/31	0/3/3/3

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	801	SAH	C5-N7	-3.63	1.32	1.39
4	A	801	SAH	C4-N9	-3.38	1.30	1.37
4	A	801	SAH	C5-C4	3.32	1.45	1.39
4	A	801	SAH	C8-N9	-3.28	1.32	1.37
4	A	801	SAH	OXT-C	-2.24	1.23	1.30

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	801	SAH	C2'-C1'-N9	-7.43	94.83	113.30
4	A	801	SAH	C5-C4-N3	-4.89	119.99	126.72
4	A	801	SAH	N3-C2-N1	-3.56	123.20	128.58
4	A	801	SAH	N3-C4-N9	3.49	133.10	127.17
4	A	801	SAH	C2-N3-C4	3.42	120.18	111.83
4	A	801	SAH	C4-C5-N7	-3.30	106.81	110.58
4	A	801	SAH	C2-N1-C6	2.67	123.11	118.73
4	A	801	SAH	C6-C5-N7	2.53	136.97	132.09
4	A	801	SAH	O4'-C1'-N9	2.12	112.17	108.09
4	A	801	SAH	C4-N9-C8	2.12	107.96	105.74

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	801	SAH	O-C-CA-N
4	A	801	SAH	O4'-C4'-C5'-SD
4	A	801	SAH	C3'-C4'-C5'-SD
4	A	801	SAH	OXT-C-CA-N
4	A	801	SAH	CA-CB-CG-SD
4	A	801	SAH	CB-CG-SD-C5'
4	A	801	SAH	C-CA-CB-CG

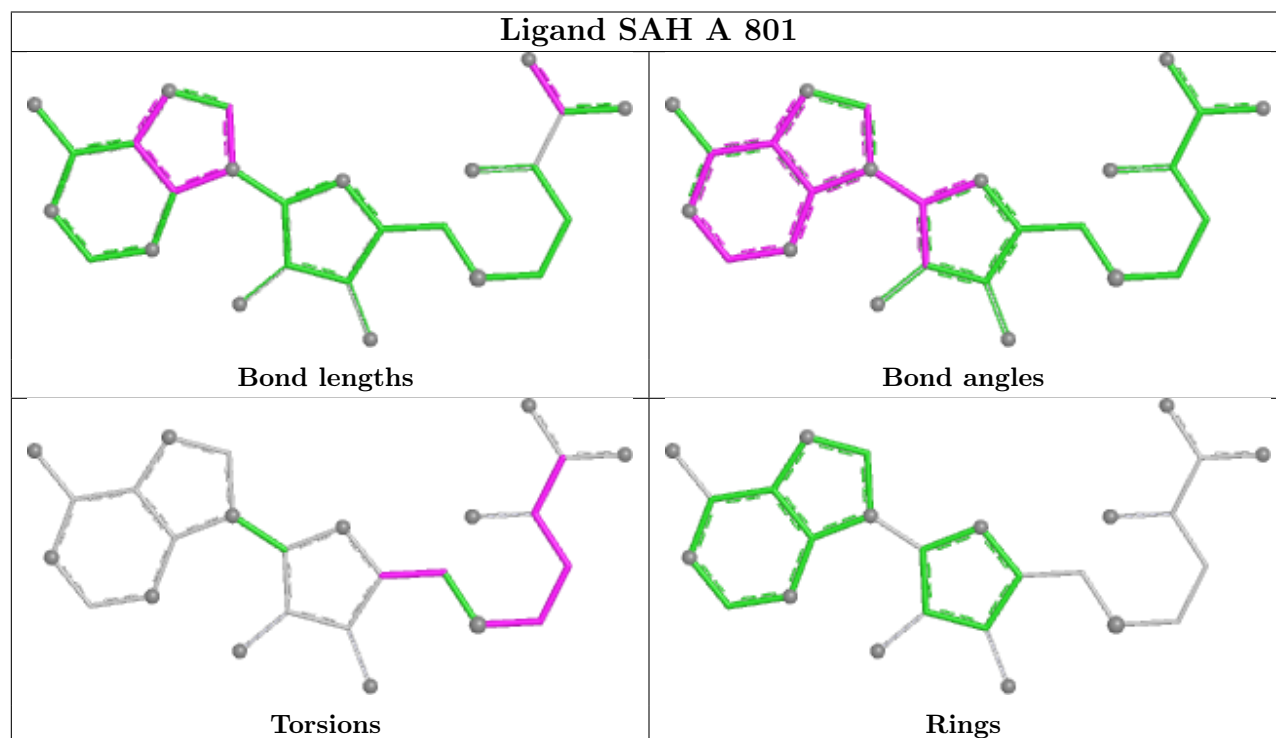
There are no ring outliers.

1 monomer is involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	801	SAH	9	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	467/533 (87%)	0.11	12 (2%) 57 36	16, 36, 76, 105	0
2	C	10/15 (66%)	0.58	1 (10%) 12 6	31, 49, 93, 147	0
3	D	13/15 (86%)	0.61	1 (7%) 19 10	38, 46, 108, 111	0
All	All	490/563 (87%)	0.14	14 (2%) 53 32	16, 36, 78, 147	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	356	VAL	4.1
1	A	499	CYS	3.3
1	A	568	GLN	3.2
2	C	11	DA	3.2
1	A	361	GLY	2.8
3	D	14	DA	2.7
1	A	455	LYS	2.6
1	A	612	CYS	2.6
1	A	614	CYS	2.6
1	A	203	MET	2.5
1	A	382	GLY	2.4
1	A	559	PHE	2.4
1	A	397	CYS	2.1
1	A	613	TYR	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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
2	5CM	C	7	20/21	0.95	0.09	25,27,32,32	0

6.3 Carbohydrates [i](#)

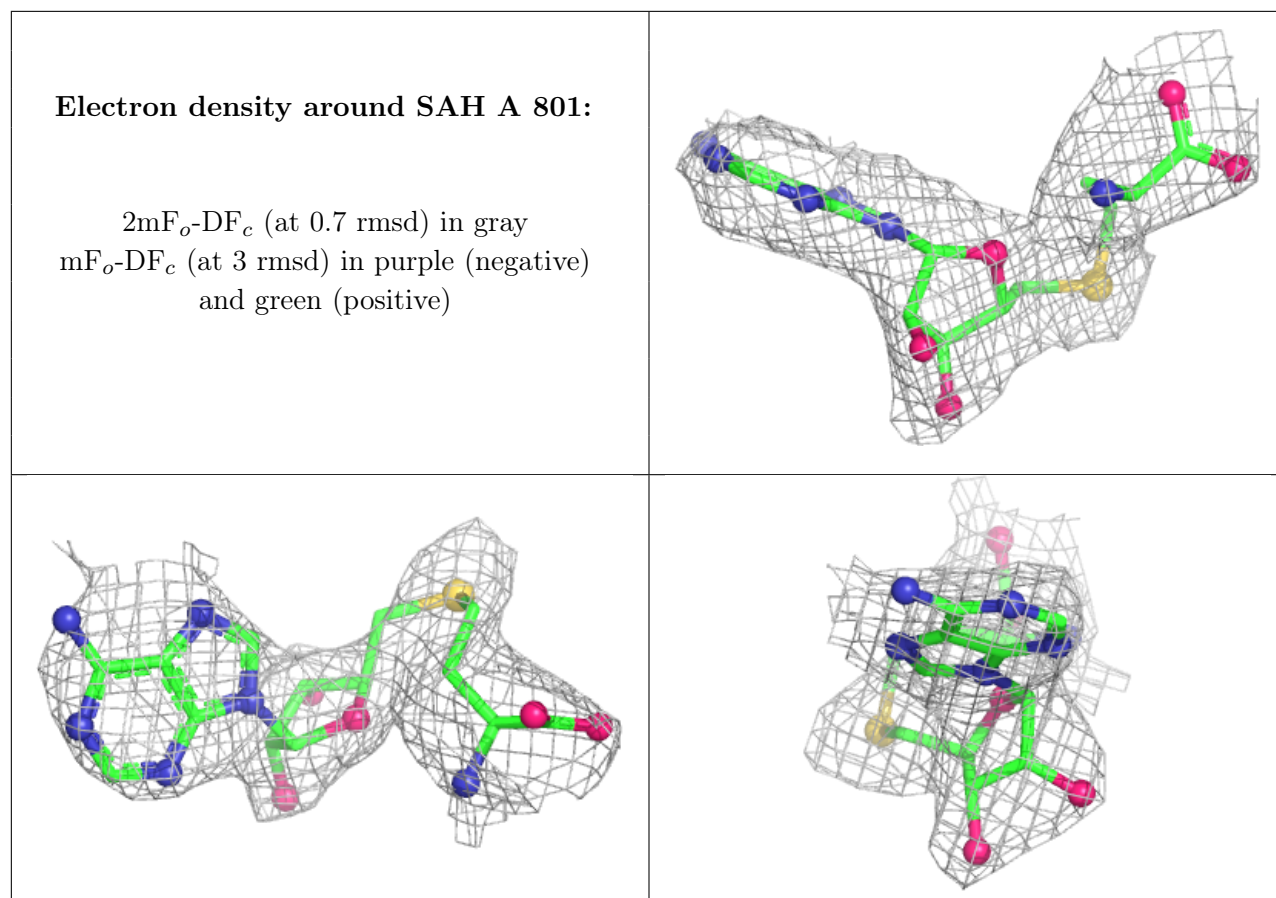
There are no oligosaccharides in this entry.

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
5	ZN	A	805	1/1	0.82	0.14	71,71,71,71	0
4	SAH	A	801	26/26	0.87	0.12	54,85,88,89	0
5	ZN	A	802	1/1	0.94	0.10	28,28,28,28	0
5	ZN	A	804	1/1	0.98	0.03	35,35,35,35	0
5	ZN	A	803	1/1	0.98	0.05	44,44,44,44	0

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



6.5 Other polymers ⓘ

There are no such residues in this entry.