



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 01:03 PM UTC

PDB ID : 4D7T / pdb_00004d7t
Title : Structure of the SthK Carboxy-Terminal Region in complex with cAMP
Authors : Kesters, D.; Brams, M.; Nys, M.; Wijckmans, E.; Spurny, R.; Voets, T.;
Tytgat, J.; Ulens, C.
Deposited on : 2014-11-27
Resolution : 2.58 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

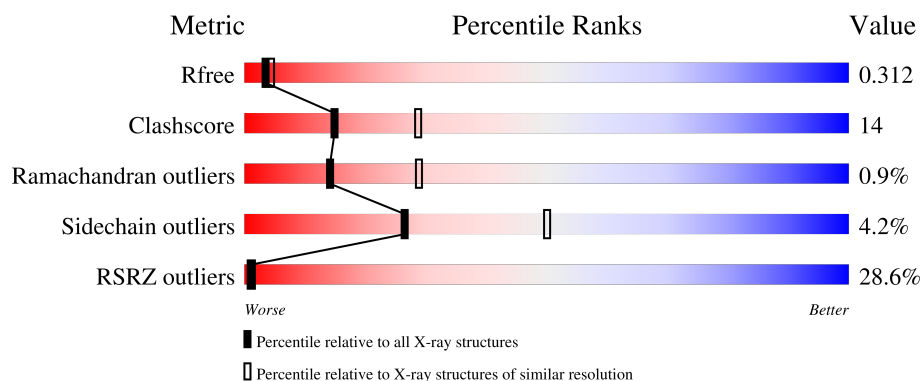
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.58 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	4770 (2.60-2.56)
Clashscore	190562	5124 (2.60-2.56)
Ramachandran outliers	187476	5046 (2.60-2.56)
Sidechain outliers	187428	5046 (2.60-2.56)
RSRZ outliers	180081	4770 (2.60-2.56)

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	198	15% 78% 22%
1	B	198	38% 57% 25% 5% 13%

2 Entry composition [i](#)

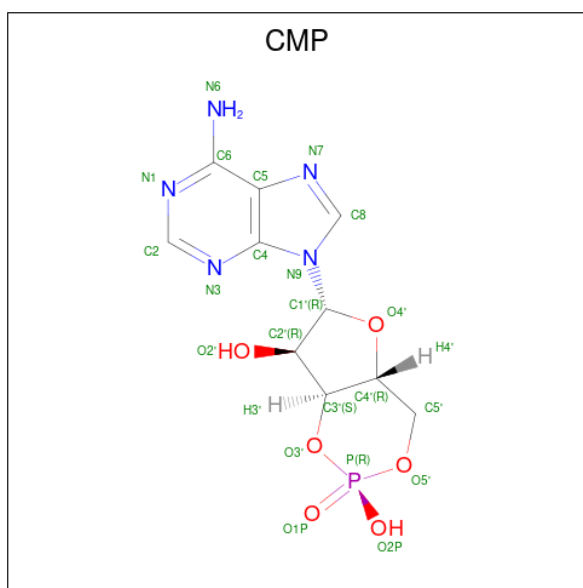
There are 3 unique types of molecules in this entry. The entry contains 2975 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 STHK_CNBD_CAMP.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	197	Total	C	N	O	S	0	0	0
			1590	1020	279	287	4			
1	B	173	Total	C	N	O	S	0	0	0
			1340	870	223	244	3			

- Molecule 2 is ADENOSINE-3',5'-CYCLIC-MONOPHOSPHATE (CCD ID: CMP) (formula: $C_{10}H_{12}N_5O_6P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			22	10	5	6	1		
2	B	1	Total	C	N	O	P	0	0
			22	10	5	6	1		

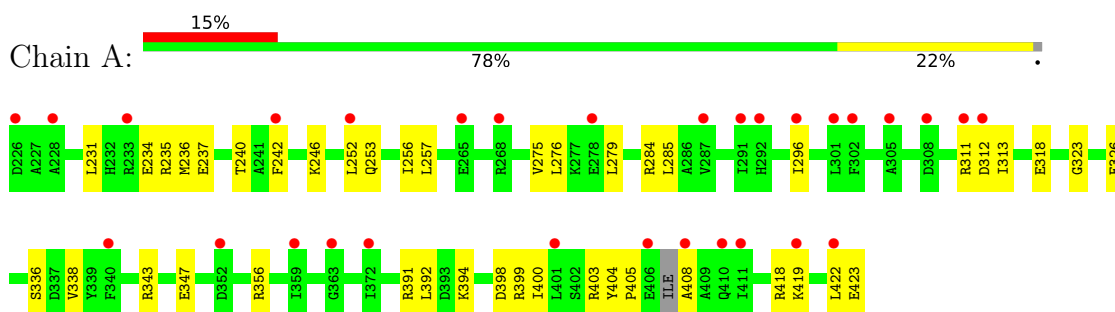
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	O	0	0
			1	1		

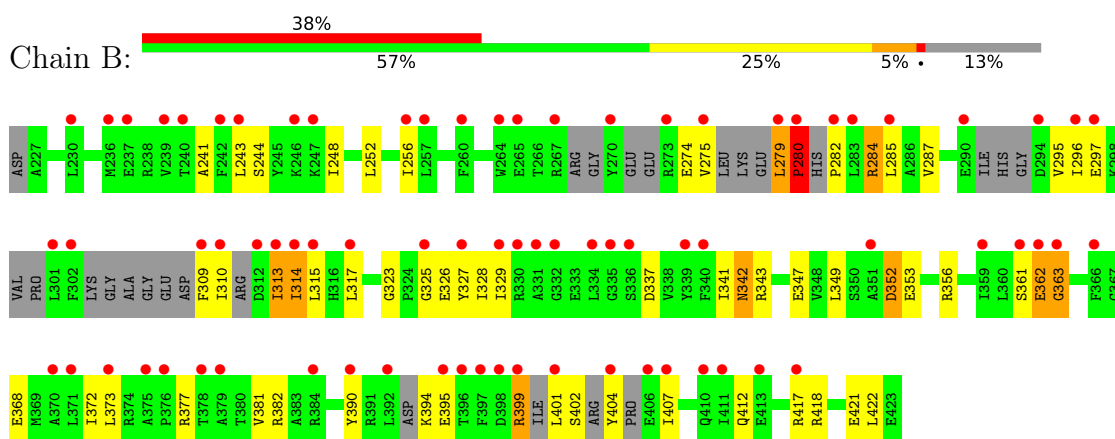
3 Residue-property plots [i](#)

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: STHK_CNBD_CAMP



• Molecule 1: STHK_CNBD_CAMP



4 Data and refinement statistics

Property	Value	Source
Space group	P 4	Depositor
Cell constants a, b, c, α , β , γ	87.44Å 87.44Å 79.47Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.80 – 2.58 48.80 – 2.58	Depositor EDS
% Data completeness (in resolution range)	98.9 (48.80-2.58) 91.8 (48.80-2.58)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.01 (at 2.58Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.265 , 0.310 0.273 , 0.312	Depositor DCC
R_{free} test set	971 reflections (5.18%)	wwPDB-VP
Wilson B-factor (Å ²)	64.3	Xtriage
Anisotropy	0.168	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 54.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.063 for h,-k,-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	2975	wwPDB-VP
Average B, all atoms (Å ²)	99.0	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 29.02 % 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 1.6649e-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: CMP

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	0.34	0/1620	0.78	0/2185
1	B	0.38	0/1356	0.96	9/1826 (0.5%)
All	All	0.36	0/2976	0.87	9/4011 (0.2%)

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	363	GLY	N-CA-C	-8.99	103.27	113.79
1	B	314	ILE	N-CA-C	7.63	118.43	110.72
1	B	395	GLU	N-CA-C	-7.28	103.84	114.39
1	B	342	ASN	N-CA-C	-6.18	103.67	112.13
1	B	315	LEU	N-CA-C	-5.70	106.09	113.16
1	B	279	LEU	CA-C-N	5.29	125.83	120.38
1	B	279	LEU	C-N-CA	5.29	125.83	120.38
1	B	280	PRO	CA-C-N	5.03	126.13	119.84
1	B	280	PRO	C-N-CA	5.03	126.13	119.84

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	1590	0	1584	23	0
1	B	1340	0	1271	56	0
2	A	22	0	11	1	0
2	B	22	0	11	3	0
3	B	1	0	0	0	0
All	All	2975	0	2877	81	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 14.

All (81) 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:A:1424:CMP:H2	2:A:1424:CMP:C2	0.97	1.49
2:B:1424:CMP:H2	2:B:1424:CMP:C2	0.97	1.48
1:B:399:ARG:HE	1:B:402:SER:CA	1.51	1.23
1:B:399:ARG:NE	1:B:402:SER:HB2	1.55	1.19
1:B:399:ARG:C	1:B:401:LEU:N	2.03	1.17
1:B:399:ARG:CD	1:B:402:SER:HB2	1.76	1.14
1:B:402:SER:HG	1:B:404:TYR:N	1.47	1.10
1:B:399:ARG:HE	1:B:402:SER:CB	1.69	1.06
1:B:399:ARG:NE	1:B:402:SER:CB	2.24	0.97
1:B:399:ARG:HD3	1:B:402:SER:HB2	1.50	0.93
1:B:399:ARG:HG3	1:B:402:SER:H	1.38	0.88
1:B:399:ARG:NE	1:B:402:SER:CA	2.36	0.87
1:B:325:GLY:O	1:B:382:ARG:NH1	2.07	0.87
1:B:399:ARG:HE	1:B:402:SER:HA	1.41	0.83
1:B:399:ARG:CG	1:B:402:SER:H	1.93	0.82
1:B:353:GLU:OE1	1:B:382:ARG:NH2	2.16	0.77
1:B:399:ARG:HG3	1:B:401:LEU:N	2.01	0.76
1:B:402:SER:OG	1:B:404:TYR:N	2.18	0.74
1:B:337:ASP:OD1	1:B:394:LYS:N	2.21	0.73
1:B:399:ARG:CD	1:B:402:SER:CB	2.60	0.73
1:B:399:ARG:HG3	1:B:402:SER:N	2.02	0.73
1:B:399:ARG:HE	1:B:402:SER:N	1.86	0.72
1:B:342:ASN:HD22	1:B:390:TYR:HE1	1.37	0.72
1:B:377:ARG:NH1	2:B:1424:CMP:O2P	2.24	0.69
1:B:285:LEU:HD12	1:B:314:ILE:HG12	1.77	0.66
1:B:362:GLU:CD	1:B:363:GLY:H	2.05	0.64
1:A:336:SER:HB2	1:A:394:LYS:HB3	1.80	0.63
1:A:313:ILE:HD13	1:A:400:ILE:HG13	1.80	0.63
1:B:342:ASN:O	1:B:362:GLU:HG3	1.99	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:399:ARG:NE	1:B:402:SER:N	2.45	0.63
1:A:275:VAL:HG13	1:A:276:LEU:HG	1.81	0.62
1:B:243:LEU:HB3	1:B:248:ILE:HG13	1.83	0.60
1:B:309:PHE:HB3	1:B:404:TYR:CZ	2.38	0.59
1:B:399:ARG:CA	1:B:401:LEU:N	2.65	0.59
1:B:323:GLY:N	1:B:326:GLU:OE2	2.35	0.59
1:A:231:LEU:O	1:A:235:ARG:HG2	2.03	0.59
1:A:398:ASP:OD1	1:A:399:ARG:N	2.35	0.58
1:A:318:GLU:OE1	1:A:391:ARG:NH1	2.37	0.58
1:B:368:GLU:OE2	1:B:418:ARG:NH1	2.37	0.57
1:A:312:ASP:OD2	1:A:404:TYR:OH	2.23	0.56
1:B:349:LEU:HB3	1:B:353:GLU:HA	1.91	0.53
1:B:252:LEU:O	1:B:256:ILE:HG12	2.10	0.52
1:B:317:LEU:HD22	1:B:390:TYR:HB3	1.92	0.52
1:B:399:ARG:N	1:B:401:LEU:N	2.57	0.52
1:B:399:ARG:CG	1:B:402:SER:N	2.65	0.51
1:B:372:ILE:HG23	1:B:412:GLN:HG2	1.93	0.51
1:B:352:ASP:OD1	1:B:352:ASP:N	2.31	0.50
1:A:235:ARG:HH11	1:A:235:ARG:HG3	1.77	0.49
1:B:241:ALA:O	1:B:244:SER:OG	2.30	0.49
1:B:313:ILE:HG23	1:B:317:LEU:HG	1.95	0.49
1:A:418:ARG:O	1:A:422:LEU:HB2	2.13	0.48
1:A:276:LEU:HA	1:A:279:LEU:HD13	1.96	0.47
1:B:399:ARG:HD3	1:B:402:SER:CB	2.35	0.47
1:B:284:ARG:H	1:B:284:ARG:HG2	1.49	0.47
1:B:279:LEU:HA	1:B:280:PRO:HD3	1.56	0.47
1:B:280:PRO:O	1:B:284:ARG:HG2	2.15	0.47
1:A:312:ASP:CG	1:A:403:ARG:HH22	2.23	0.46
1:B:295:VAL:HG21	1:B:342:ASN:HA	1.98	0.46
1:A:276:LEU:O	1:A:284:ARG:HD3	2.15	0.46
1:A:285:LEU:HD21	1:A:311:ARG:HG3	1.98	0.46
1:B:296:ILE:HD13	1:B:310:ILE:HG13	1.97	0.46
1:A:419:LYS:HE3	1:A:423:GLU:OE2	2.17	0.45
1:A:338:VAL:HG13	1:A:392:LEU:HB3	1.98	0.45
1:B:328:ILE:O	1:B:329:ILE:HD13	2.17	0.44
1:B:284:ARG:HA	1:B:287:VAL:HG22	1.99	0.44
1:B:368:GLU:OE1	2:B:1424:CMP:O2'	2.34	0.44
1:A:253:GLN:O	1:A:257:LEU:HG	2.18	0.44
1:A:347:GLU:OE2	1:A:356:ARG:NH2	2.49	0.44
1:B:399:ARG:NE	1:B:402:SER:HA	2.20	0.43
1:B:327:TYR:OH	1:B:382:ARG:NH2	2.51	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:418:ARG:O	1:B:422:LEU:HB2	2.20	0.42
1:A:405:PRO:HA	1:A:408:ALA:HB3	2.02	0.41
1:B:417:ARG:O	1:B:421:GLU:HG2	2.20	0.41
1:A:234:GLU:C	1:A:236:MET:H	2.28	0.41
1:A:252:LEU:O	1:A:256:ILE:HG12	2.21	0.41
1:B:347:GLU:O	1:B:381:VAL:HA	2.21	0.41
1:A:237:GLU:HA	1:A:240:THR:HB	2.03	0.41
1:B:342:ASN:HB3	1:B:343:ARG:HG2	2.03	0.41
1:A:242:PHE:O	1:A:246:LYS:HG2	2.21	0.40
1:A:323:GLY:H	1:A:326:GLU:CD	2.25	0.40
1:B:341:ILE:HD12	1:B:361:SER:O	2.21	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	193/198 (98%)	184 (95%)	9 (5%)	0	100	100
1	B	151/198 (76%)	135 (89%)	13 (9%)	3 (2%)	6	11
All	All	344/396 (87%)	319 (93%)	22 (6%)	3 (1%)	14	29

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	280	PRO
1	B	282	PRO
1	B	297	GLU

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	161/171 (94%)	159 (99%)	2 (1%)	63	82
1	B	128/171 (75%)	118 (92%)	10 (8%)	11	25
All	All	289/342 (84%)	277 (96%)	12 (4%)	26	50

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	296	ILE
1	A	343	ARG
1	B	274	GLU
1	B	275	VAL
1	B	284	ARG
1	B	313	ILE
1	B	352	ASP
1	B	356	ARG
1	B	362	GLU
1	B	373	LEU
1	B	399	ARG
1	B	407	ILE

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

Mol	Chain	Res	Type
1	B	342	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

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	CMP	A	1424	-	25,25,25	2.16	7 (28%)	37,39,39	1.78	11 (29%)
2	CMP	B	1424	-	25,25,25	2.10	6 (24%)	37,39,39	1.84	11 (29%)

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	CMP	A	1424	-	-	0/4/31/31	0/4/4/4
2	CMP	B	1424	-	-	0/4/31/31	0/4/4/4

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1424	CMP	P-O5'	6.69	1.65	1.57
2	B	1424	CMP	P-O5'	6.34	1.64	1.57
2	A	1424	CMP	P-O3'	3.83	1.64	1.57
2	B	1424	CMP	P-O3'	3.65	1.63	1.57
2	A	1424	CMP	O5'-C5'	-3.45	1.41	1.46
2	B	1424	CMP	O5'-C5'	-3.12	1.41	1.46
2	B	1424	CMP	C5-C4	3.12	1.44	1.39
2	A	1424	CMP	C5-C4	3.03	1.44	1.39
2	B	1424	CMP	O3'-C3'	-3.01	1.40	1.44
2	A	1424	CMP	O3'-C3'	-2.67	1.40	1.44
2	A	1424	CMP	O4'-C4'	-2.65	1.39	1.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1424	CMP	O4'-C4'	-2.19	1.40	1.45
2	A	1424	CMP	C2-N1	2.02	1.37	1.33

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1424	CMP	C4-N9-C8	5.15	111.15	105.74
2	A	1424	CMP	C4-N9-C8	5.10	111.10	105.74
2	B	1424	CMP	O5'-P-O3'	-4.31	99.93	105.70
2	A	1424	CMP	O5'-P-O3'	-3.92	100.46	105.70
2	A	1424	CMP	C5-C4-N9	-2.97	102.57	105.81
2	B	1424	CMP	C5-C4-N9	-2.93	102.62	105.81
2	B	1424	CMP	O2P-P-O5'	2.92	114.27	107.16
2	A	1424	CMP	N3-C4-N9	2.74	131.82	127.17
2	B	1424	CMP	C4'-O4'-C1'	-2.67	103.58	109.47
2	B	1424	CMP	N3-C4-N9	2.64	131.66	127.17
2	A	1424	CMP	O2P-P-O5'	2.64	113.61	107.16
2	A	1424	CMP	C2'-C1'-N9	-2.62	106.79	113.30
2	B	1424	CMP	C5'-C4'-C3'	2.60	117.45	112.56
2	A	1424	CMP	O4'-C1'-N9	-2.58	103.14	108.09
2	A	1424	CMP	C4'-O4'-C1'	-2.45	104.06	109.47
2	B	1424	CMP	O4'-C1'-N9	-2.44	103.41	108.09
2	A	1424	CMP	N9-C8-N7	-2.32	110.64	113.94
2	B	1424	CMP	N9-C8-N7	-2.31	110.66	113.94
2	A	1424	CMP	O3'-C3'-C2'	2.31	117.87	115.61
2	B	1424	CMP	C2'-C3'-C4'	-2.29	99.23	103.24
2	A	1424	CMP	C5'-C4'-C3'	2.13	116.57	112.56
2	B	1424	CMP	C2'-C1'-N9	-2.12	108.03	113.30

There are no chirality outliers.

There are no torsion outliers.

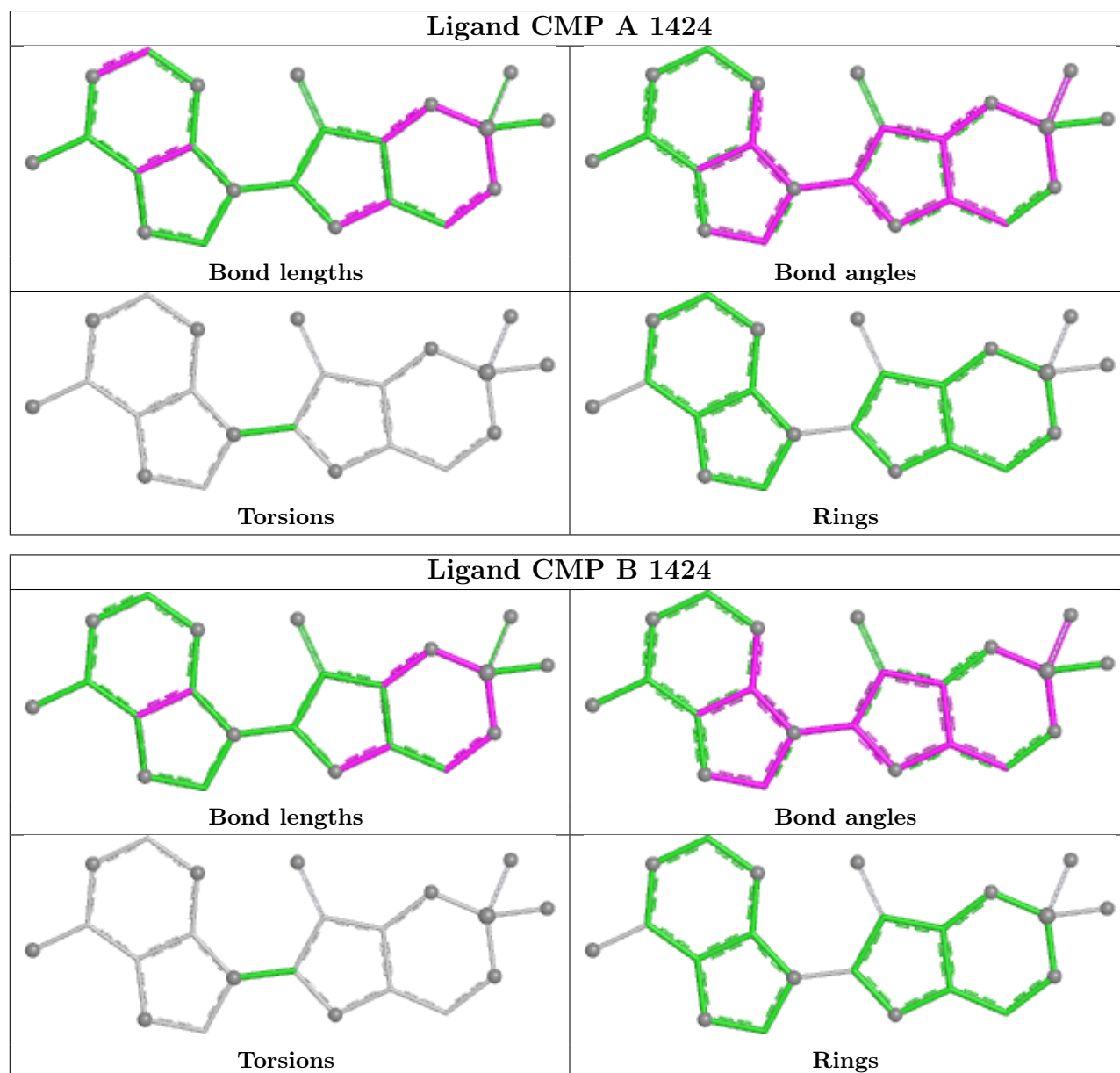
There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1424	CMP	1	0
2	B	1424	CMP	3	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 ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	197/198 (99%)	1.08	30 (15%) 5 4	51, 79, 136, 192	0
1	B	173/198 (87%)	2.10	76 (43%) 0 0	79, 112, 166, 228	0
All	All	370/396 (93%)	1.56	106 (28%) 1 1	51, 98, 151, 228	0

All (106) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	397	PHE	12.3
1	B	313	ILE	9.6
1	B	314	ILE	8.7
1	B	256	ILE	8.6
1	B	398	ASP	6.4
1	B	378	THR	5.3
1	B	242	PHE	5.3
1	B	373	LEU	4.9
1	B	310	ILE	4.8
1	B	334	LEU	4.8
1	A	302	PHE	4.7
1	B	392	LEU	4.6
1	A	406	GLU	4.4
1	B	396	THR	4.4
1	B	406	GLU	4.4
1	B	309	PHE	4.3
1	B	407	ILE	4.2
1	B	312	ASP	4.1
1	B	302	PHE	4.1
1	B	379	ALA	4.0
1	B	275	VAL	3.8
1	B	279	LEU	3.8
1	B	294	ASP	3.8
1	B	282	PRO	3.8

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Mol	Chain	Res	Type	RSRZ
1	A	340	PHE	3.7
1	B	362	GLU	3.7
1	B	325	GLY	3.6
1	B	257	LEU	3.6
1	B	317	LEU	3.6
1	B	363	GLY	3.6
1	A	226	ASP	3.6
1	B	361	SER	3.5
1	B	283	LEU	3.5
1	A	311	ARG	3.4
1	B	301	LEU	3.4
1	B	243	LEU	3.3
1	B	315	LEU	3.2
1	B	366	PHE	3.2
1	A	308	ASP	3.2
1	B	395	GLU	3.2
1	B	411	ILE	3.2
1	A	228	ALA	3.1
1	B	237	GLU	3.1
1	B	401	LEU	3.1
1	B	351	ALA	3.0
1	B	376	PRO	3.0
1	B	260	PHE	3.0
1	A	265	GLU	2.9
1	B	330	ARG	2.9
1	B	270	TYR	2.9
1	A	411	ILE	2.9
1	B	265	GLU	2.8
1	A	296	ILE	2.8
1	B	329	ILE	2.8
1	B	404	TYR	2.8
1	B	246	LYS	2.7
1	A	291	ILE	2.7
1	A	301	LEU	2.7
1	B	399	ARG	2.7
1	B	240	THR	2.7
1	A	352	ASP	2.6
1	B	296	ILE	2.6
1	B	359	ILE	2.6
1	A	410	GLN	2.6
1	B	280	PRO	2.6
1	A	408	ALA	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	327	TYR	2.6
1	B	410	GLN	2.5
1	B	264	TRP	2.5
1	B	267	ARG	2.5
1	B	417	ARG	2.5
1	A	305	ALA	2.5
1	B	384	ARG	2.4
1	A	278	GLU	2.4
1	B	273	ARG	2.3
1	A	252	LEU	2.3
1	A	419	LYS	2.3
1	B	297	GLU	2.3
1	B	375	ALA	2.3
1	B	390	TYR	2.3
1	B	340	PHE	2.3
1	A	287	VAL	2.3
1	B	339	TYR	2.3
1	A	292	HIS	2.3
1	B	371	LEU	2.3
1	B	239	VAL	2.3
1	B	332	GLY	2.2
1	B	290	GLU	2.2
1	A	372	ILE	2.2
1	B	336	SER	2.2
1	A	233	ARG	2.2
1	A	363	GLY	2.1
1	A	401	LEU	2.1
1	B	335	GLY	2.1
1	B	331	ALA	2.1
1	B	230	LEU	2.1
1	B	370	ALA	2.1
1	B	285	LEU	2.1
1	B	247	LYS	2.1
1	A	422	LEU	2.1
1	A	359	ILE	2.0
1	B	236	MET	2.0
1	A	242	PHE	2.0
1	A	268	ARG	2.0
1	B	413	GLU	2.0
1	A	312	ASP	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](#)

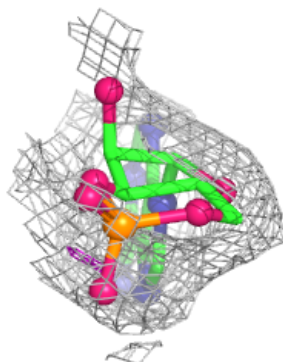
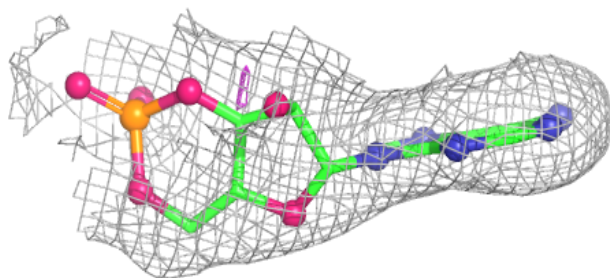
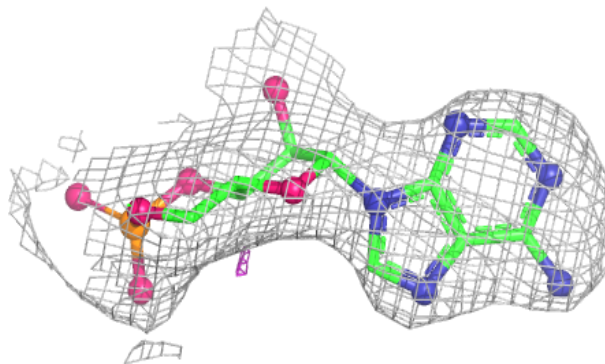
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
2	CMP	B	1424	22/22	0.91	0.11	69,78,106,135	0
2	CMP	A	1424	22/22	0.95	0.09	55,68,80,113	0

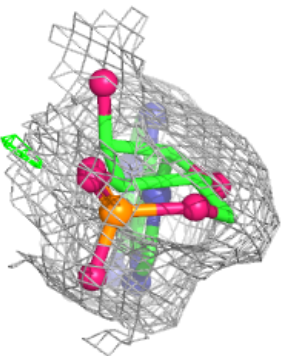
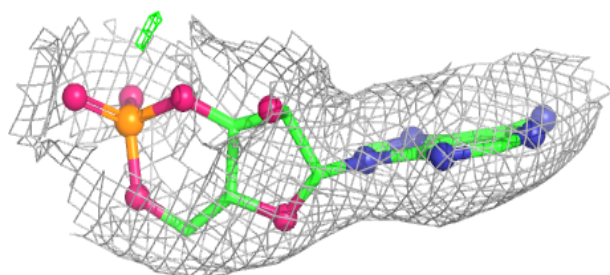
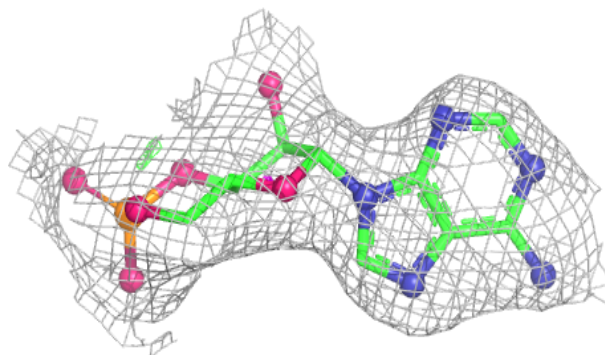
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around CMP B 1424:

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

**Electron density around CMP A 1424:**

$2mF_o-DF_c$ (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.