



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 29, 2024 – 06:16 pm GMT

PDB ID : 8CNB  
Title : Crystal structure of CREBBP-Y1503C histone acetyltransferase domain in complex with Coenzyme A  
Authors : Mechaly, A.E.; Zhang, W.; Haouz, A.; Green, M.; Rodrigues-Lima, F.  
Deposited on : 2023-02-22  
Resolution : 1.99 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](mailto: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 : 4.02b-467  
Mogul : 1.8.4, CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.36  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.36

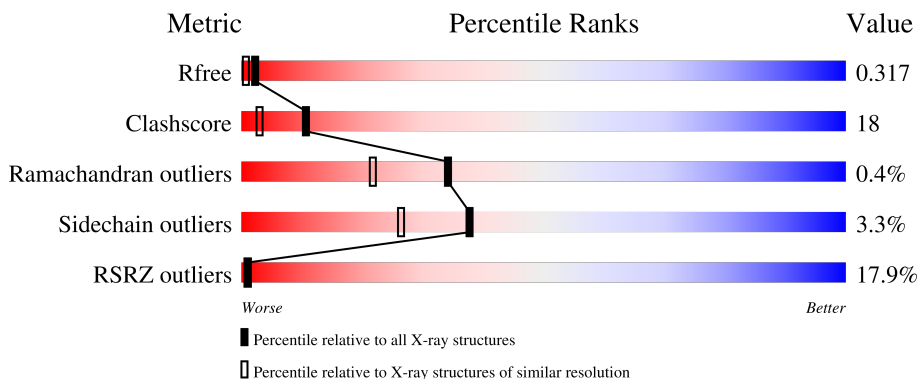
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.99 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	11647 (2.00-1.96)
Clashscore	141614	1014 (1.98-1.98)
Ramachandran outliers	138981	1006 (1.98-1.98)
Sidechain outliers	138945	1006 (1.98-1.98)
RSRZ outliers	127900	11410 (2.00-1.96)

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  $\geq 3$ , 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	619	

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 4514 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 acetyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	531	4409	2821	759	794	35	0	1	0

There are 11 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1075	MET	-	initiating methionine	UNP F8VPR5
A	1076	GLY	-	expression tag	UNP F8VPR5
A	1077	SER	-	expression tag	UNP F8VPR5
A	1503	CYS	TYR	engineered mutation	UNP F8VPR5
A	1555	GLU	-	linker	UNP F8VPR5
A	1613	SER	-	linker	UNP F8VPR5
A	1614	GLY	-	linker	UNP F8VPR5
A	1615	GLY	-	linker	UNP F8VPR5
A	1616	SER	-	linker	UNP F8VPR5
A	1617	GLY	-	linker	UNP F8VPR5
A	1618	SER	-	linker	UNP F8VPR5

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

- Molecule 3 is COENZYME A (three-letter code: COA) (formula: C<sub>21</sub>H<sub>36</sub>N<sub>7</sub>O<sub>16</sub>P<sub>3</sub>S) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
			Total	C	N	O	P			S
3	A	1	48	21	7	16	3	1	0	0

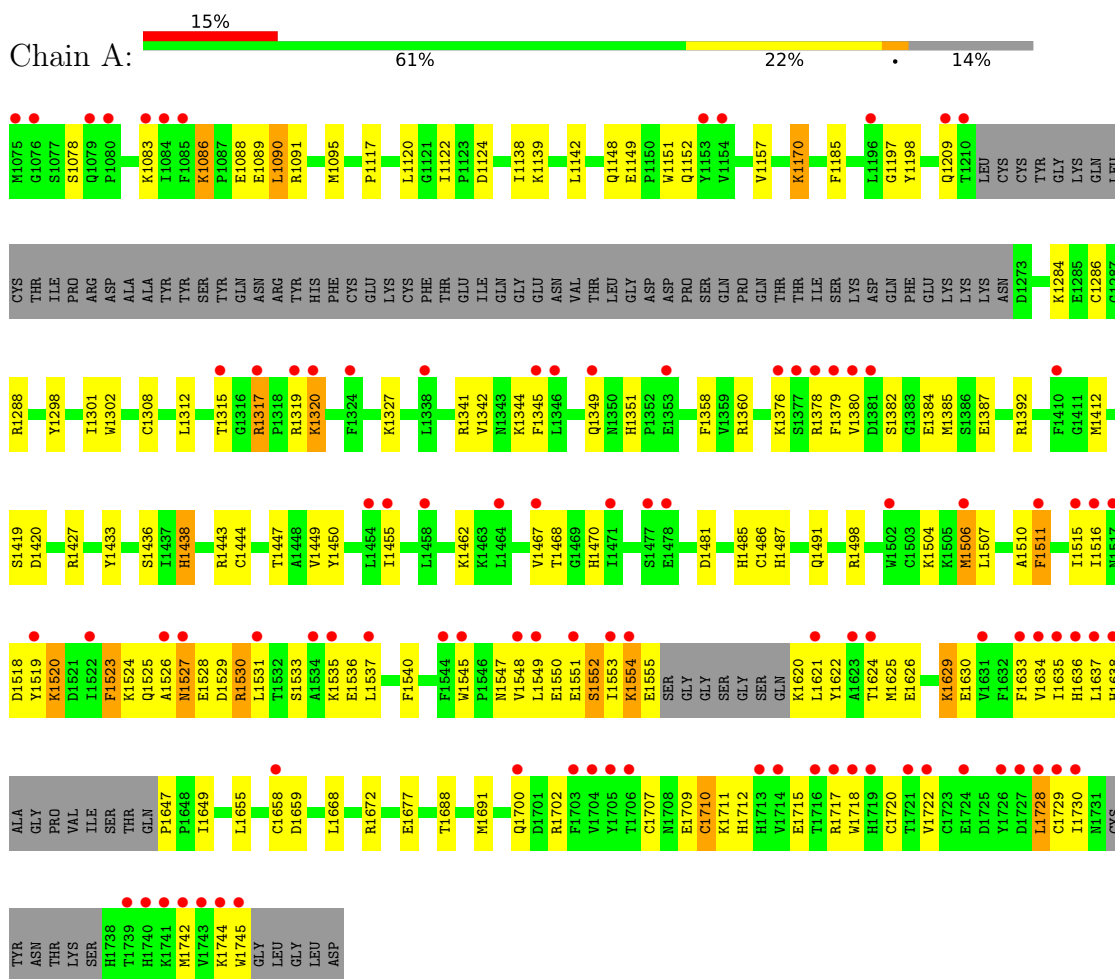
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	53	Total	O	0	0
			53	53		

### 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: histone acetyltransferase



## 4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	86.67Å 144.82Å 155.44Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.10 – 1.99 105.96 – 1.98	Depositor EDS
% Data completeness (in resolution range)	49.9 (46.10-1.99) 49.9 (105.96-1.98)	Depositor EDS
$R_{merge}$	0.11	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.17 (at 1.98Å)	Xtrriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, $R_{free}$	0.262 , 0.316 0.262 , 0.317	Depositor DCC
$R_{free}$ test set	1705 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	45.0	Xtrriage
Anisotropy	0.094	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 44.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	4514	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	67.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.91% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: ZN, COA

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	0.64	3/4534 (0.1%)	0.98	17/6127 (0.3%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1530	ARG	CG-CD	-6.09	1.36	1.51
1	A	1523	PHE	CB-CG	-5.62	1.41	1.51
1	A	1523	PHE	CD1-CE1	-5.49	1.28	1.39

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1530	ARG	CG-CD-NE	-9.41	92.03	111.80
1	A	1378	ARG	NE-CZ-NH1	9.23	124.92	120.30
1	A	1728	LEU	CA-CB-CG	8.33	134.46	115.30
1	A	1710	CYS	CA-CB-SG	7.61	127.69	114.00
1	A	1523	PHE	CB-CG-CD1	-7.39	115.63	120.80
1	A	1378	ARG	NE-CZ-NH2	-7.29	116.66	120.30
1	A	1715	GLU	CB-CA-C	6.68	123.76	110.40
1	A	1530	ARG	CA-CB-CG	-6.25	99.65	113.40
1	A	1537	LEU	CA-CB-CG	-6.21	101.01	115.30
1	A	1715	GLU	CA-CB-CG	-6.10	99.99	113.40
1	A	1317	ARG	CA-CB-CG	5.93	126.44	113.40
1	A	1090	LEU	CB-CG-CD2	-5.61	101.47	111.00
1	A	1120	LEU	CA-CB-CG	5.60	128.18	115.30
1	A	1527	ASN	CB-CA-C	5.12	120.64	110.40
1	A	1520	LYS	CA-CB-CG	5.10	124.63	113.40
1	A	1481	ASP	CB-CG-OD2	-5.04	113.76	118.30
1	A	1344	LYS	CB-CG-CD	-5.02	98.56	111.60

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	4409	0	4277	159	0
2	A	4	0	0	0	0
3	A	48	0	32	4	0
4	A	53	0	0	18	0
All	All	4514	0	4309	161	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 18.

All (161) 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:1302:TRP:HD1	1:A:1320:LYS:HE2	1.15	1.03
1:A:1302:TRP:CD1	1:A:1320:LYS:HE2	1.97	0.98
1:A:1385:MET:HE3	1:A:1535:LYS:HB2	1.49	0.94
1:A:1462:LYS:HB2	1:A:1637:LEU:HD23	1.51	0.92
1:A:1707:CYS:SG	4:A:1937:HOH:O	2.28	0.90
1:A:1298:TYR:HB3	1:A:1301:ILE:HD12	1.56	0.87
1:A:1729:CYS:SG	4:A:1937:HOH:O	2.34	0.84
1:A:1710:CYS:SG	4:A:1937:HOH:O	2.35	0.83
1:A:1717:ARG:HA	1:A:1730:ILE:HD11	1.60	0.83
1:A:1095:MET:HE2	1:A:1139:LYS:HG3	1.62	0.81
1:A:1711:LYS:O	1:A:1711:LYS:HD2	1.80	0.80
1:A:1523:PHE:CE1	1:A:1625:MET:HG2	2.16	0.80
1:A:1467:VAL:HG13	1:A:1468:THR:HG23	1.64	0.79
1:A:1717:ARG:O	4:A:1901:HOH:O	2.00	0.78
1:A:1523:PHE:HE1	1:A:1625:MET:HG2	1.47	0.78
1:A:1342:VAL:HG21	1:A:1449:VAL:HG13	1.66	0.78
1:A:1718:TRP:CD1	1:A:1730:ILE:HG13	2.20	0.77
1:A:1485:HIS:O	1:A:1485:HIS:ND1	2.19	0.75
1:A:1312:LEU:HA	1:A:1315:THR:HB	1.68	0.74

*Continued on next page...*



*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1523:PHE:CE2	1:A:1626:GLU:HA	2.22	0.74
1:A:1707:CYS:SG	1:A:1729:CYS:N	2.61	0.72
1:A:1138:ILE:HD13	1:A:1157:VAL:HG12	1.75	0.69
1:A:1518:ASP:OD1	1:A:1519:TYR:N	2.24	0.69
1:A:1718:TRP:HD1	1:A:1730:ILE:HG13	1.59	0.68
1:A:1436:SER:HB3	3:A:1805:COA:O9P	1.94	0.68
1:A:1504:LYS:NZ	4:A:1904:HOH:O	2.27	0.68
1:A:1523:PHE:CE2	1:A:1629:LYS:HB3	2.28	0.68
1:A:1444[A]:CYS:SG	4:A:1942:HOH:O	2.52	0.67
1:A:1444[B]:CYS:SG	4:A:1942:HOH:O	2.52	0.67
1:A:1485:HIS:O	1:A:1485:HIS:CG	2.48	0.67
1:A:1447:THR:OG1	3:A:1805:COA:O4A	2.12	0.66
1:A:1523:PHE:CD2	1:A:1629:LYS:HB3	2.30	0.66
1:A:1545:TRP:HA	1:A:1548:VAL:HG12	1.78	0.66
1:A:1545:TRP:HH2	1:A:1625:MET:HA	1.61	0.66
3:A:1805:COA:O7A	4:A:1902:HOH:O	2.14	0.66
1:A:1419:SER:OG	1:A:1649:ILE:HG13	1.97	0.64
1:A:1717:ARG:HA	1:A:1730:ILE:CD1	2.26	0.64
1:A:1455:ILE:HD13	1:A:1510:ALA:HB2	1.79	0.64
1:A:1525:GLN:NE2	1:A:1529:ASP:OD1	2.24	0.64
1:A:1379:PHE:CD2	1:A:1535:LYS:HD3	2.33	0.63
1:A:1745:TRP:N	4:A:1901:HOH:O	2.32	0.62
1:A:1385:MET:CE	1:A:1535:LYS:HB2	2.28	0.62
1:A:1320:LYS:HD3	1:A:1320:LYS:H	1.64	0.62
1:A:1349:GLN:N	1:A:1349:GLN:OE1	2.33	0.62
1:A:1149:GLU:HG2	1:A:1151:TRP:H	1.65	0.61
1:A:1649:ILE:HD12	1:A:1649:ILE:O	2.01	0.61
1:A:1148:GLN:N	1:A:1152:GLN:OE1	2.34	0.61
1:A:1327:LYS:O	1:A:1360:ARG:NH2	2.30	0.60
1:A:1486:CYS:HB3	1:A:1677:GLU:HB3	1.82	0.60
1:A:1622:TYR:O	1:A:1626:GLU:HG2	2.01	0.60
1:A:1327:LYS:HB2	1:A:1358:PHE:CE1	2.36	0.60
1:A:1688:THR:HA	1:A:1691:MET:HE3	1.84	0.59
1:A:1551:GLU:C	1:A:1553:ILE:H	2.05	0.59
1:A:1315:THR:HG22	1:A:1317:ARG:HB3	1.85	0.59
1:A:1547:ASN:HB2	4:A:1924:HOH:O	2.04	0.58
1:A:1709:GLU:HB3	4:A:1937:HOH:O	2.02	0.58
1:A:1717:ARG:HH21	1:A:1745:TRP:HE3	1.49	0.58
1:A:1702:ARG:HA	1:A:1702:ARG:HE	1.68	0.58
1:A:1462:LYS:HB2	1:A:1637:LEU:CD2	2.29	0.58
1:A:1524:LYS:O	1:A:1528:GLU:HB2	2.03	0.57

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1438:HIS:CE1	1:A:1443:ARG:HH22	2.22	0.57
1:A:1170:LYS:H	1:A:1170:LYS:HD2	1.70	0.57
1:A:1209:GLN:OE1	4:A:1903:HOH:O	2.17	0.57
1:A:1436:SER:HG	1:A:1450:TYR:HH	1.46	0.57
1:A:1551:GLU:O	1:A:1553:ILE:N	2.38	0.56
1:A:1376:LYS:O	1:A:1380:VAL:HG12	2.06	0.56
1:A:1462:LYS:NZ	4:A:1911:HOH:O	2.38	0.56
1:A:1545:TRP:CH2	1:A:1625:MET:HA	2.40	0.56
1:A:1376:LYS:HG2	1:A:1380:VAL:CG1	2.37	0.55
1:A:1302:TRP:HD1	1:A:1320:LYS:CE	2.04	0.54
1:A:1523:PHE:CZ	1:A:1625:MET:O	2.60	0.54
1:A:1455:ILE:CD1	1:A:1510:ALA:HB2	2.38	0.54
1:A:1718:TRP:CZ2	1:A:1744:LYS:HD2	2.43	0.53
1:A:1519:TYR:OH	4:A:1904:HOH:O	2.19	0.53
1:A:1170:LYS:H	1:A:1170:LYS:CD	2.21	0.53
1:A:1384:GLU:HB3	1:A:1535:LYS:HE2	1.89	0.53
1:A:1379:PHE:HD2	1:A:1535:LYS:HD3	1.74	0.53
1:A:1487:HIS:HD2	1:A:1491:GLN:NE2	2.07	0.53
1:A:1341:ARG:HH12	1:A:1455:ILE:HB	1.73	0.53
1:A:1523:PHE:HZ	1:A:1625:MET:O	1.90	0.53
1:A:1545:TRP:CE3	1:A:1548:VAL:HG11	2.44	0.53
1:A:1629:LYS:HG2	1:A:1630:GLU:N	2.24	0.52
1:A:1088:GLU:OE2	4:A:1905:HOH:O	2.19	0.52
1:A:1462:LYS:HG3	1:A:1638:HIS:NE2	2.24	0.52
1:A:1523:PHE:CZ	1:A:1625:MET:C	2.83	0.52
1:A:1507:LEU:HD11	1:A:1633:PHE:HD2	1.74	0.52
1:A:1527:ASN:O	1:A:1530:ARG:HG3	2.09	0.52
1:A:1630:GLU:O	1:A:1633:PHE:HE1	1.92	0.52
1:A:1658:CYS:HA	1:A:1700:GLN:OE1	2.09	0.52
1:A:1095:MET:HE3	1:A:1142:LEU:HD23	1.91	0.52
1:A:1717:ARG:CA	1:A:1730:ILE:HD11	2.37	0.52
1:A:1455:ILE:HG12	1:A:1506:MET:HE2	1.91	0.51
1:A:1526:ALA:O	1:A:1531:LEU:HD22	2.09	0.51
1:A:1286:CYS:CB	1:A:1288:ARG:HG2	2.41	0.51
1:A:1551:GLU:O	1:A:1553:ILE:HG13	2.11	0.51
1:A:1157:VAL:HG21	1:A:1185:PHE:CZ	2.47	0.50
1:A:1341:ARG:HH12	1:A:1455:ILE:CB	2.23	0.50
1:A:1520:LYS:HG3	1:A:1524:LYS:HD2	1.93	0.50
1:A:1198:TYR:CE1	1:A:1288:ARG:HB2	2.47	0.50
1:A:1548:VAL:HG21	1:A:1624:THR:HG21	1.94	0.49
1:A:1533:SER:OG	1:A:1536:GLU:HG3	2.12	0.49

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1376:LYS:HG2	1:A:1380:VAL:HG11	1.95	0.49
1:A:1427:ARG:NH2	1:A:1647:PRO:O	2.45	0.49
1:A:1659:ASP:N	1:A:1700:GLN:OE1	2.42	0.49
1:A:1720:CYS:SG	1:A:1722:VAL:HG12	2.54	0.48
1:A:1302:TRP:CD1	1:A:1320:LYS:HG2	2.49	0.48
1:A:1380:VAL:HG21	1:A:1387:GLU:HG3	1.95	0.48
1:A:1284:LYS:NZ	4:A:1918:HOH:O	2.47	0.47
1:A:1634:VAL:C	1:A:1635:ILE:HD13	2.35	0.47
1:A:1520:LYS:CG	1:A:1524:LYS:HD2	2.44	0.47
1:A:1341:ARG:HH22	1:A:1455:ILE:HG21	1.78	0.47
1:A:1523:PHE:CD1	1:A:1523:PHE:N	2.75	0.47
1:A:1462:LYS:HD3	4:A:1911:HOH:O	2.15	0.47
1:A:1548:VAL:CG2	1:A:1624:THR:HG21	2.45	0.47
1:A:1485:HIS:CE1	1:A:1672:ARG:HG2	2.49	0.47
1:A:1717:ARG:NH2	1:A:1745:TRP:HE3	2.12	0.47
1:A:1730:ILE:N	1:A:1730:ILE:HD12	2.30	0.47
1:A:1550:GLU:O	1:A:1553:ILE:HG23	2.14	0.46
1:A:1086:LYS:N	1:A:1089:GLU:OE1	2.43	0.46
1:A:1511:PHE:HA	1:A:1516:ILE:O	2.16	0.45
1:A:1555:GLU:CD	1:A:1620:LYS:HE3	2.37	0.45
1:A:1197:GLY:O	1:A:1288:ARG:HD3	2.16	0.45
1:A:1668:LEU:O	1:A:1672:ARG:HG3	2.17	0.45
1:A:1553:ILE:HD12	1:A:1554:LYS:N	2.32	0.45
1:A:1702:ARG:HA	1:A:1702:ARG:NE	2.32	0.45
1:A:1345:PHE:CZ	1:A:1349:GLN:NE2	2.85	0.45
1:A:1470:HIS:HD2	1:A:1540:PHE:CZ	2.35	0.45
1:A:1552:SER:HB2	1:A:1621:LEU:HD21	1.99	0.44
1:A:1485:HIS:ND1	1:A:1672:ARG:HG2	2.33	0.44
1:A:1392:ARG:CZ	1:A:1655:LEU:HD21	2.47	0.44
1:A:1086:LYS:HG3	1:A:1089:GLU:HG3	1.99	0.44
1:A:1419:SER:H	1:A:1649:ILE:CG1	2.30	0.44
1:A:1717:ARG:NH1	1:A:1745:TRP:O	2.51	0.44
1:A:1516:ILE:HD13	1:A:1516:ILE:HA	1.93	0.44
1:A:1549:LEU:HD22	1:A:1621:LEU:HD13	2.00	0.43
1:A:1707:CYS:SG	1:A:1729:CYS:HB3	2.58	0.43
1:A:1083:LYS:HG2	1:A:1288:ARG:HH21	1.84	0.43
1:A:1462:LYS:HD2	1:A:1637:LEU:HB3	2.00	0.43
1:A:1515:ILE:HD13	1:A:1515:ILE:HA	1.83	0.43
1:A:1531:LEU:HD12	1:A:1536:GLU:CD	2.39	0.43
1:A:1157:VAL:HG21	1:A:1185:PHE:CE2	2.53	0.43
1:A:1148:GLN:HG2	1:A:1152:GLN:HE22	1.83	0.42

*Continued on next page...*

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1342:VAL:CG2	1:A:1449:VAL:HG13	2.45	0.42
1:A:1341:ARG:HH22	1:A:1455:ILE:CG2	2.33	0.42
1:A:1525:GLN:OE1	1:A:1636:HIS:NE2	2.30	0.42
1:A:1659:ASP:H	1:A:1700:GLN:CD	2.21	0.42
1:A:1385:MET:HE3	1:A:1385:MET:HB3	1.80	0.42
1:A:1090:LEU:HA	1:A:1090:LEU:HD23	1.75	0.42
1:A:1710:CYS:C	1:A:1712:HIS:H	2.22	0.42
1:A:1527:ASN:C	1:A:1530:ARG:H	2.23	0.42
1:A:1551:GLU:C	1:A:1553:ILE:N	2.71	0.42
1:A:1086:LYS:HG2	1:A:1089:GLU:CD	2.40	0.41
1:A:1728:LEU:HD12	1:A:1742:MET:CE	2.50	0.41
1:A:1117:PRO:HB3	1:A:1122:ILE:O	2.21	0.41
1:A:1455:ILE:HD11	1:A:1506:MET:O	2.21	0.40
1:A:1531:LEU:HD12	1:A:1536:GLU:HB3	2.03	0.40
3:A:1805:COA:P3B	4:A:1902:HOH:O	2.79	0.40
1:A:1341:ARG:HH12	1:A:1455:ILE:CG2	2.35	0.40
1:A:1412:MET:HA	1:A:1433:TYR:O	2.21	0.40
1:A:1523:PHE:CZ	1:A:1625:MET:HG2	2.56	0.40
1:A:1545:TRP:HA	1:A:1548:VAL:CG1	2.46	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	522/619 (84%)	505 (97%)	15 (3%)	2 (0%)	34 22

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1552	SER
1	A	1078	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	491/567 (87%)	475 (97%)	16 (3%)	38 26

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1086	LYS
1	A	1091	ARG
1	A	1124	ASP
1	A	1170	LYS
1	A	1308	CYS
1	A	1319	ARG
1	A	1320	LYS
1	A	1351	HIS
1	A	1382	SER
1	A	1420	ASP
1	A	1438	HIS
1	A	1498	ARG
1	A	1506	MET
1	A	1511	PHE
1	A	1554	LYS
1	A	1629	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1470	HIS
1	A	1487	HIS
1	A	1527	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 monosaccharides 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
3	COA	A	1805	1	41,50,50	0.57	0	52,75,75	0.87	3 (5%)

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	COA	A	1805	1	-	11/44/64/64	0/3/3/3

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1805	COA	C5A-C6A-N6A	2.42	124.03	120.35
3	A	1805	COA	O5B-P1A-O1A	2.19	117.63	109.07
3	A	1805	COA	O5A-P2A-O4A	2.06	122.42	112.24

There are no chirality outliers.

All (11) torsion outliers are listed below:

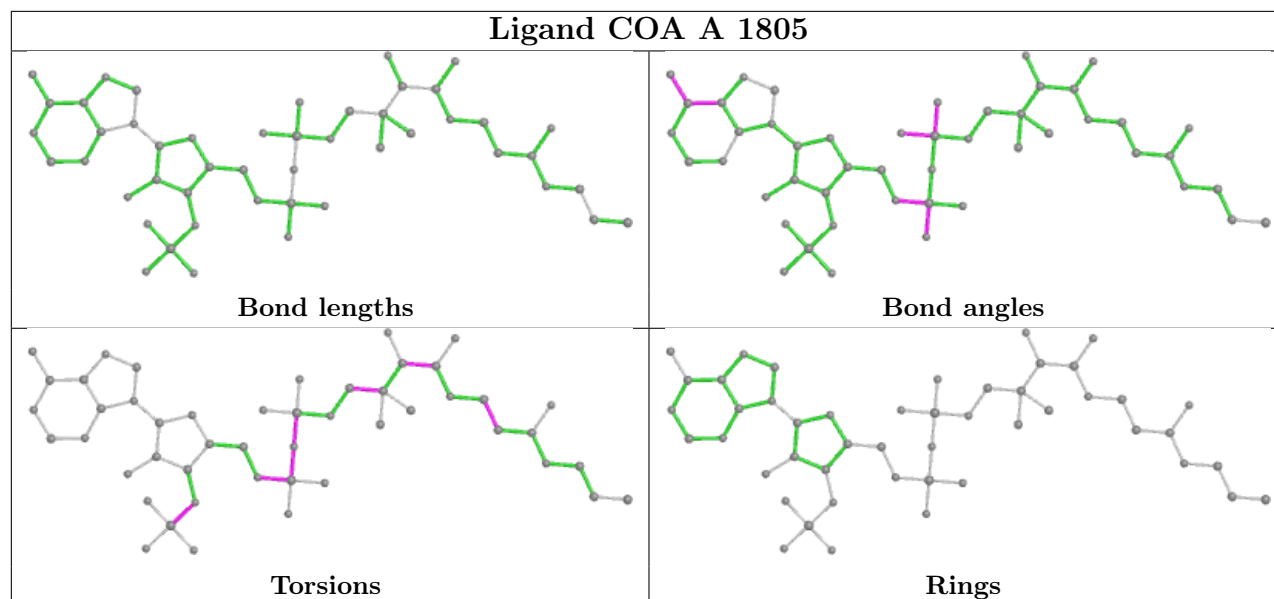
Mol	Chain	Res	Type	Atoms
3	A	1805	COA	C3B-O3B-P3B-O8A
3	A	1805	COA	C5B-O5B-P1A-O1A
3	A	1805	COA	C5B-O5B-P1A-O3A
3	A	1805	COA	O9P-C9P-CAP-OAP
3	A	1805	COA	N8P-C9P-CAP-OAP
3	A	1805	COA	C5P-C6P-C7P-N8P
3	A	1805	COA	P1A-O3A-P2A-O4A
3	A	1805	COA	P1A-O3A-P2A-O6A
3	A	1805	COA	C5B-O5B-P1A-O2A
3	A	1805	COA	P2A-O3A-P1A-O2A
3	A	1805	COA	CEP-CBP-CCP-O6A

There are no ring outliers.

1 monomer is involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1805	COA	4	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	531/619 (85%)	1.33	95 (17%) <b>1</b> <b>1</b>	27, 60, 121, 200	0

All (95) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1714	VAL	9.5
1	A	1705	TYR	8.0
1	A	1346	LEU	7.4
1	A	1380	VAL	7.1
1	A	1713	HIS	6.9
1	A	1743	VAL	6.9
1	A	1718	TRP	6.8
1	A	1516	ILE	6.6
1	A	1531	LEU	6.4
1	A	1319	ARG	6.2
1	A	1317	ARG	6.1
1	A	1379	PHE	6.1
1	A	1728	LEU	5.8
1	A	1079	GLN	5.5
1	A	1717	ARG	5.5
1	A	1553	ILE	5.4
1	A	1381	ASP	5.4
1	A	1703	PHE	5.2
1	A	1534	ALA	4.9
1	A	1624	THR	4.7
1	A	1621	LEU	4.7
1	A	1740	HIS	4.6
1	A	1537	LEU	4.6
1	A	1745	TRP	4.5
1	A	1455	ILE	4.5
1	A	1722	VAL	4.3
1	A	1515	ILE	4.2

*Continued on next page...*

*Continued from previous page...*

<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
1	A	1635	ILE	4.2
1	A	1730	ILE	4.2
1	A	1721	THR	4.2
1	A	1554	LYS	4.2
1	A	1741	LYS	4.1
1	A	1076	GLY	4.1
1	A	1526	ALA	4.1
1	A	1502	TRP	4.0
1	A	1377	SER	4.0
1	A	1719	HIS	3.8
1	A	1535	LYS	3.8
1	A	1716	THR	3.7
1	A	1631	VAL	3.7
1	A	1700	GLN	3.5
1	A	1742	MET	3.5
1	A	1637	LEU	3.4
1	A	1196	LEU	3.4
1	A	1548	VAL	3.4
1	A	1634	VAL	3.4
1	A	1545	TRP	3.4
1	A	1636	HIS	3.2
1	A	1464	LEU	3.1
1	A	1633	PHE	3.1
1	A	1345	PHE	3.0
1	A	1519	TYR	3.0
1	A	1511	PHE	2.9
1	A	1744	LYS	2.9
1	A	1704	VAL	2.8
1	A	1376	LYS	2.8
1	A	1084	ILE	2.8
1	A	1477	SER	2.8
1	A	1410	PHE	2.8
1	A	1378	ARG	2.8
1	A	1210	THR	2.8
1	A	1551	GLU	2.8
1	A	1315	THR	2.7
1	A	1727	ASP	2.7
1	A	1739	THR	2.7
1	A	1471	ILE	2.6
1	A	1458	LEU	2.6
1	A	1522	ILE	2.6
1	A	1544	PHE	2.6

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
1	A	1209	GLN	2.6
1	A	1517	ASN	2.5
1	A	1658	CYS	2.5
1	A	1353	GLU	2.5
1	A	1467	VAL	2.4
1	A	1549	LEU	2.4
1	A	1506	MET	2.4
1	A	1706	THR	2.4
1	A	1320	LYS	2.4
1	A	1724	GLU	2.4
1	A	1638	HIS	2.3
1	A	1154	VAL	2.3
1	A	1454	LEU	2.3
1	A	1729	CYS	2.3
1	A	1324	PHE	2.2
1	A	1623	ALA	2.2
1	A	1527	ASN	2.2
1	A	1726	TYR	2.2
1	A	1478	GLU	2.2
1	A	1349	GLN	2.1
1	A	1338	LEU	2.1
1	A	1080	PRO	2.1
1	A	1085	PHE	2.1
1	A	1075	MET	2.0
1	A	1083	LYS	2.0
1	A	1153	TYR	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 [\(i\)](#)

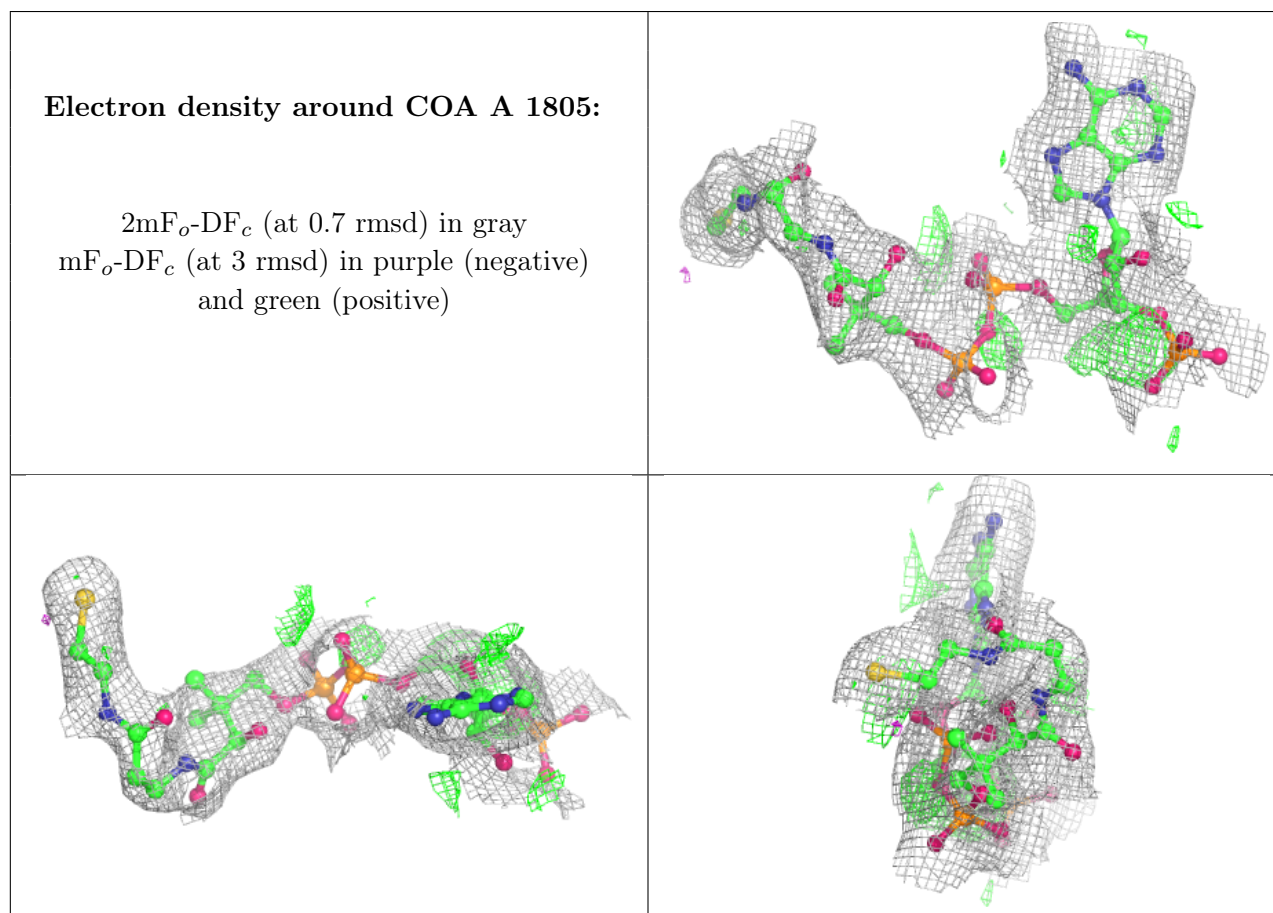
There are no monosaccharides in this entry.

## 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, 95<sup>th</sup> 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	ZN	A	1804	1/1	0.66	0.10	202,202,202,202	0
3	COA	A	1805	48/48	0.93	0.19	73,80,98,103	0
2	ZN	A	1803	1/1	0.94	0.36	280,280,280,280	0
2	ZN	A	1802	1/1	0.99	0.16	75,75,75,75	0
2	ZN	A	1801	1/1	1.00	0.23	53,53,53,53	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 [i](#)

There are no such residues in this entry.