



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 10:07 AM GMT

PDB ID : 1D6H
Title : CHALONE SYNTHASE (N336A MUTANT COMPLEXED WITH COA)
Authors : Jez, J.M.; Ferrer, J.L.; Bowman, M.E.; Dixon, R.A.; Noel, J.P.
Deposited on : 1999-10-13
Resolution : 2.15 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

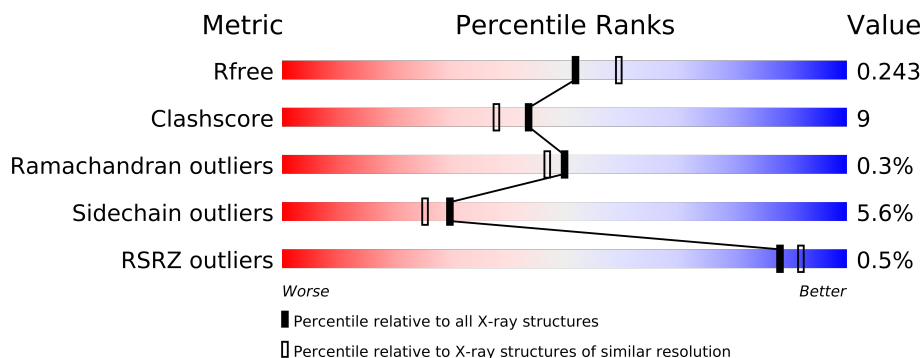
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance


The reported resolution of this entry is 2.15 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}	66092	1094 (2.18-2.14)
Clashscore	79885	1299 (2.18-2.14)
Ramachandran outliers	78287	1272 (2.18-2.14)
Sidechain outliers	78261	1272 (2.18-2.14)
RSRZ outliers	66119	1094 (2.18-2.14)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	387	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
3	COA	A	390	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3224 atoms, of which 0 are hydrogen and 0 are deuterium.

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 CHALCONE SYNTHASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	387	2987	1899	503	565	20	0	1	0

There is a discrepancy between the modelled and reference sequences:

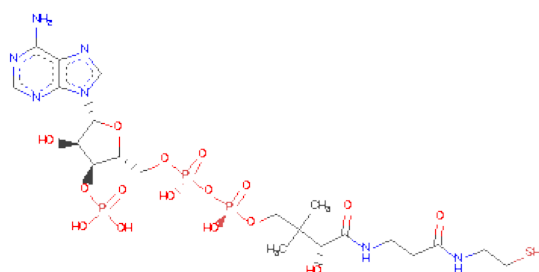
Chain	Residue	Modelled	Actual	Comment	Reference
A	336	ALA	ASN	ENGINEERED	UNP P30074

- Molecule 2 is SULFATE ION (three-letter code: SO₄) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	S		
2	A	1	5	4	1	0	0

- Molecule 3 is COENZYME A (three-letter code: COA) (formula: C₂₁H₃₆N₇O₁₆P₃S).



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

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	184	Total	O		
			184	184	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	97.36Å 97.36Å 65.21Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	85.00 – 2.15 20.06 – 2.15	Depositor EDS
% Data completeness (in resolution range)	(Not available) (85.00-2.15) 92.0 (20.06-2.15)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.16 (at 2.15Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.184 , 0.257 0.174 , 0.243	Depositor DCC
R_{free} test set	932 reflections (5.42%)	DCC
Wilson B-factor (Å ²)	29.1	Xtriage
Anisotropy	0.091	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 56.0	EDS
Estimated twinning fraction	0.043 for -h,-k,l	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 18142 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3224	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.34% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: COA, CSD, SO4

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	0.83	0/3036	1.71	51/4108 (1.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	11

There are no bond length outliers.

All (51) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	151	ARG	NE-CZ-NH2	-12.06	114.27	120.30
1	A	311	ASP	CB-CG-OD2	-11.84	107.64	118.30
1	A	185	ARG	NE-CZ-NH2	-10.09	115.25	120.30
1	A	68	ARG	NE-CZ-NH1	-9.98	115.31	120.30
1	A	67	ARG	NE-CZ-NH1	9.97	125.28	120.30
1	A	94	ARG	CD-NE-CZ	9.47	136.86	123.60
1	A	376	GLY	O-C-N	-9.24	107.92	122.70
1	A	67	ARG	NE-CZ-NH2	-8.66	115.97	120.30
1	A	376	GLY	CA-C-N	8.54	135.99	117.20
1	A	207	ASP	CB-CG-OD2	-8.30	110.83	118.30
1	A	61	ASP	CB-CG-OD2	-7.73	111.34	118.30
1	A	389	ILE	CA-CB-CG1	-7.61	96.54	111.00
1	A	342	VAL	N-CA-CB	-7.47	95.06	111.50
1	A	38	ASP	CB-CG-OD1	7.45	125.00	118.30
1	A	348	GLU	OE1-CD-OE2	-7.25	114.59	123.30
1	A	138	PRO	N-CA-CB	6.96	111.65	103.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	255	ASP	CB-CG-OD2	6.94	124.55	118.30
1	A	350	ARG	NE-CZ-NH1	6.92	123.76	120.30
1	A	172	ARG	NE-CZ-NH2	-6.82	116.89	120.30
1	A	203	ASP	CB-CG-OD2	6.54	124.18	118.30
1	A	165	PHE	O-C-N	6.38	132.91	122.70
1	A	58	ARG	NE-CZ-NH2	6.25	123.43	120.30
1	A	270	ASP	CB-CG-OD1	6.06	123.75	118.30
1	A	54	GLU	CA-CB-CG	6.03	126.66	113.40
1	A	165	PHE	C-N-CA	-5.98	106.76	121.70
1	A	137	MET	CA-C-O	-5.94	107.62	120.10
1	A	199	ARG	NE-CZ-NH1	-5.92	117.34	120.30
1	A	259	ARG	NE-CZ-NH1	5.83	123.21	120.30
1	A	8	ARG	NE-CZ-NH2	5.79	123.20	120.30
1	A	96	ASP	CB-CG-OD1	5.72	123.45	118.30
1	A	376	GLY	N-CA-C	-5.64	99.00	113.10
1	A	138	PRO	CA-N-CD	-5.58	103.69	111.50
1	A	358	LEU	O-C-N	-5.52	113.88	122.70
1	A	199	ARG	NE-CZ-NH2	5.49	123.05	120.30
1	A	185	ARG	NE-CZ-NH1	5.42	123.01	120.30
1	A	47	GLU	OE1-CD-OE2	5.41	129.80	123.30
1	A	385	ARG	NE-CZ-NH2	5.41	123.01	120.30
1	A	294	ASP	CB-CG-OD1	5.41	123.17	118.30
1	A	58	ARG	NE-CZ-NH1	-5.38	117.61	120.30
1	A	388	ALA	N-CA-CB	5.36	117.60	110.10
1	A	156	ARG	NE-CZ-NH2	5.35	122.97	120.30
1	A	165	PHE	CB-CG-CD2	5.35	124.54	120.80
1	A	141	ASP	CB-CG-OD2	-5.35	113.49	118.30
1	A	315	GLN	CA-CB-CG	5.33	125.13	113.40
1	A	179	GLU	OE1-CD-OE2	5.26	129.62	123.30
1	A	12	ARG	CD-NE-CZ	5.21	130.90	123.60
1	A	199	ARG	CD-NE-CZ	5.12	130.77	123.60
1	A	237	PHE	CB-CG-CD2	-5.06	117.25	120.80
1	A	170	VAL	CA-CB-CG2	5.06	118.49	110.90
1	A	227	ASP	CB-CG-OD1	5.06	122.85	118.30
1	A	10	ALA	N-CA-CB	-5.05	103.03	110.10

There are no chirality outliers.

All (11) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	188	VAL	Mainchain
1	A	218	GLY	Mainchain

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Mol	Chain	Res	Type	Group
1	A	229	VAL	Mainchain
1	A	289	PRO	Mainchain
1	A	305	GLY	Mainchain,Peptide
1	A	345	ILE	Mainchain
1	A	348	GLU	Mainchain
1	A	360	THR	Mainchain
1	A	378	THR	Mainchain
1	A	64[B]	MET	Mainchain

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2987	0	3029	52	0
2	A	5	0	0	0	0
3	A	48	0	31	5	0
4	A	184	0	0	6	2
All	All	3224	0	3060	53	2

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 9.

All (53) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:342:VAL:HG13	1:A:370:LEU:HD11	1.55	0.89
1:A:51:GLU:HG2	4:A:530:HOH:O	1.78	0.83
1:A:104:ARG:NH2	1:A:108:GLU:OE1	2.16	0.74
1:A:58:ARG:HB3	1:A:62:LYS:HZ1	1.53	0.74
1:A:58:ARG:HB3	1:A:62:LYS:NZ	2.02	0.73
1:A:38:ASP:O	1:A:42:LYS:HG3	1.88	0.72
1:A:42:LYS:HG2	1:A:47:GLU:OE1	1.91	0.70
1:A:62:LYS:HE3	3:A:390:COA:P1A	2.38	0.64
1:A:267:LEU:HD21	3:A:390:COA:S1P	2.40	0.62
1:A:271:VAL:HB	1:A:272:PRO:HD3	1.82	0.60
1:A:42:LYS:HG2	1:A:47:GLU:CD	2.23	0.59

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:50:THR:O	1:A:54:GLU:HG2	2.03	0.58
1:A:82:ASN:ND2	4:A:534:HOH:O	2.38	0.56
1:A:102:VAL:HB	1:A:103:PRO:HD3	1.87	0.55
1:A:352:LYS:HD3	4:A:504:HOH:O	2.05	0.55
1:A:230:PRO:O	1:A:231:GLU:HB2	2.07	0.54
1:A:292:ILE:HD12	1:A:293:SER:N	2.23	0.54
1:A:126:HIS:HE1	4:A:446:HOH:O	1.90	0.54
1:A:348:GLU:OE2	1:A:352:LYS:HD2	2.10	0.52
1:A:38:ASP:OD2	1:A:53:LYS:NZ	2.34	0.51
1:A:55:LYS:NZ	3:A:390:COA:O7A	2.38	0.51
1:A:14:GLU:HG3	1:A:227:ASP:OD2	2.11	0.50
3:A:390:COA:H52A	3:A:390:COA:H122	1.93	0.50
1:A:301:ILE:HG22	1:A:342:VAL:HG22	1.97	0.47
1:A:256:GLY:HA2	1:A:264:THR:O	2.15	0.47
1:A:197:THR:CG2	1:A:263:LEU:HD23	2.44	0.47
1:A:292:ILE:HD12	1:A:292:ILE:C	2.34	0.47
1:A:20:LEU:HD23	1:A:235:PRO:HA	1.97	0.47
1:A:241:TRP:HH2	1:A:285:GLU:HG3	1.80	0.46
1:A:24:THR:HB	1:A:344:PHE:CZ	2.50	0.45
1:A:5:SER:O	1:A:9:LYS:HG3	2.17	0.45
1:A:62:LYS:HE3	3:A:390:COA:O1A	2.17	0.45
1:A:100:VAL:O	1:A:103:PRO:HD2	2.17	0.45
1:A:129:VAL:HG21	1:A:141:ASP:HA	1.99	0.44
1:A:271:VAL:CB	1:A:272:PRO:HD3	2.48	0.44
1:A:197:THR:HG22	1:A:263:LEU:HD23	1.99	0.44
1:A:389:ILE:HA	1:A:389:ILE:HD12	1.54	0.44
1:A:194:THR:HG23	1:A:217:ASP:OD1	2.18	0.44
1:A:29:ASN:HB3	1:A:70:MET:O	2.17	0.43
1:A:299:PHE:CZ	1:A:368:GLY:HA3	2.53	0.43
1:A:188:VAL:O	1:A:221:ALA:HA	2.18	0.43
1:A:316:LYS:HE2	4:A:505:HOH:O	2.18	0.43
1:A:36:TYR:N	1:A:37:PRO:CD	2.81	0.43
1:A:240:VAL:HG21	1:A:367:TRP:HZ3	1.84	0.43
1:A:58:ARG:O	1:A:62:LYS:HD3	2.19	0.42
1:A:241:TRP:CH2	1:A:285:GLU:HG3	2.54	0.42
1:A:151:ARG:HD2	1:A:151:ARG:HH11	1.62	0.42
1:A:125:THR:OG1	1:A:126:HIS:HD2	2.03	0.41
1:A:172:ARG:HA	1:A:242:THR:HG21	2.02	0.41
1:A:342:VAL:HB	4:A:410:HOH:O	2.19	0.41
1:A:286:ALA:HB1	1:A:383:VAL:CG2	2.51	0.41
1:A:95:GLN:OE1	1:A:134:GLY:HA2	2.20	0.41
1:A:137:MET:HA	1:A:138:PRO:C	2.42	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
4:A:400:HOH:O	4:A:506:HOH:O[4_557]	2.10	0.10
4:A:470:HOH:O	4:A:543:HOH:O[4_557]	2.17	0.03

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

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	385/387 (100%)	374 (97%)	10 (3%)	1 (0%)	50	47

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	90	SER

5.3.2 Protein sidechains ⓘ

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	322/321 (100%)	304 (94%)	18 (6%)	30	25

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	SER
1	A	14	GLU
1	A	54	GLU
1	A	67	ARG
1	A	75	GLU

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Mol	Chain	Res	Type
1	A	133	SER
1	A	231	GLU
1	A	284	VAL
1	A	285	GLU
1	A	292	ILE
1	A	314	GLU
1	A	320	LYS
1	A	342	VAL
1	A	350	ARG
1	A	352	LYS
1	A	355	GLN
1	A	359	LYS
1	A	389	ILE

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

Mol	Chain	Res	Type
1	A	126	HIS
1	A	212	GLN
1	A	355	GLN

5.3.3 RNA ⓘ

There are no RNA chains 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	$\# Z > 2$
1	CSD	A	164	1	7,7,8	7.11	4 (57%)	6,8,10	7.96	5 (83%)

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
1	CSD	A	164	1	-	0/3/6/8	0/0/0/0

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	164	CSD	OD1-SG	-14.22	1.28	1.47
1	A	164	CSD	O-C	10.70	1.18	1.11
1	A	164	CSD	CB-SG	-5.03	1.62	1.80
1	A	164	CSD	CA-C	2.98	1.53	1.48

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	164	CSD	OD2-SG-CB	17.10	166.25	97.42
1	A	164	CSD	OD1-SG-CB	-5.97	89.42	105.25
1	A	164	CSD	OD2-SG-OD1	-4.98	101.12	109.39
1	A	164	CSD	C-CA-N	3.60	117.42	113.83
1	A	164	CSD	CB-CA-N	-3.49	104.56	110.16

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.5 Carbohydrates ⓘ

There are no carbohydrates 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	$\# Z > 2$
3	COA	A	390	-	50,50,50	1.33	10 (20%)	75,75,75	2.35	22 (29%)
2	SO4	A	391	-	4,4,4	1.15	0	6,6,6	0.44	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	COA	A	390	-	-	0/48/64/64	0/1/3/3
2	SO4	A	391	-	-	0/0/0/0	0/0/0/0

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	390	COA	OAP-CAP	-3.86	1.35	1.42
3	A	390	COA	C3P-N4P	3.02	1.53	1.46
3	A	390	COA	C7P-N8P	2.62	1.52	1.46
3	A	390	COA	P3B-O8A	-2.59	1.45	1.54
3	A	390	COA	P3B-O9A	-2.48	1.45	1.54
3	A	390	COA	CCP-CBP	2.26	1.56	1.52
3	A	390	COA	P3B-O7A	2.21	1.58	1.51
3	A	390	COA	C2A-N1A	2.10	1.38	1.33
3	A	390	COA	P2A-O6A	-2.06	1.49	1.59
3	A	390	COA	O4B-C1B	-2.01	1.38	1.41

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	390	COA	CDP-CBP-CAP	8.39	123.36	108.82
3	A	390	COA	C7P-N8P-C9P	-6.01	110.24	122.57
3	A	390	COA	O9A-P3B-O7A	-5.65	91.96	110.44
3	A	390	COA	C3P-N4P-C5P	-5.36	111.92	122.84
3	A	390	COA	C6P-C5P-N4P	4.54	124.76	116.50
3	A	390	COA	O5P-C5P-N4P	-4.07	114.87	122.94
3	A	390	COA	CBP-CAP-C9P	-3.99	108.86	112.73
3	A	390	COA	O9P-C9P-N8P	3.94	131.05	123.05
3	A	390	COA	C4B-O4B-C1B	3.74	113.81	109.75
3	A	390	COA	CCP-CBP-CAP	-3.70	103.28	108.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	390	COA	O8A-P3B-O3B	3.69	117.71	107.09
3	A	390	COA	O4B-C4B-C3B	-3.43	97.08	104.88
3	A	390	COA	OAP-CAP-CBP	-3.11	103.47	110.71
3	A	390	COA	CEP-CBP-CDP	-3.04	104.75	109.08
3	A	390	COA	O9A-P3B-O3B	2.92	115.50	107.09
3	A	390	COA	O4B-C1B-N9A	2.82	111.06	108.44
3	A	390	COA	O4B-C1B-C2B	-2.70	102.63	106.77
3	A	390	COA	C2B-C3B-C4B	2.57	108.16	103.16
3	A	390	COA	O9P-C9P-CAP	-2.49	115.95	120.48
3	A	390	COA	O6A-CCP-CBP	-2.07	107.11	110.57
3	A	390	COA	C5A-C6A-N6A	2.04	125.33	120.72
3	A	390	COA	C8A-N9A-C4A	-2.02	105.35	106.90

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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	387/387 (100%)	-0.34	1 (0%) 91 95	24, 32, 45, 59	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	231	GLU	3.0

6.2 Non-standard residues in protein, DNA, RNA chains

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	RSR	LLDF	B-factors(Å ²)	Q<0.9
1	CSD	A	164	8/9	0.10	0.17	30,43,55,56	0

6.3 Carbohydrates

There are no carbohydrates in this entry.

6.4 Ligands

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	COA	A	390	48/48	0.38	6.91	24,40,64,65	48
2	SO4	A	391	5/5	0.13	-0.72	52,53,53,55	5

6.5 Other polymers ⓘ

There are no such residues in this entry.