



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 02:35 AM UTC

PDB ID : 2CTS / pdb_00002cts
Title : CRYSTALLOGRAPHIC REFINEMENT AND ATOMIC MODELS OF
TWO DIFFERENT FORMS OF CITRATE SYNTHASE AT 2.7 AND 1.7
ANGSTROMS RESOLUTION
Authors : Remington, S.; Wiegand, G.; Huber, R.
Deposited on : 1984-01-27
Resolution : 2.00 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) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

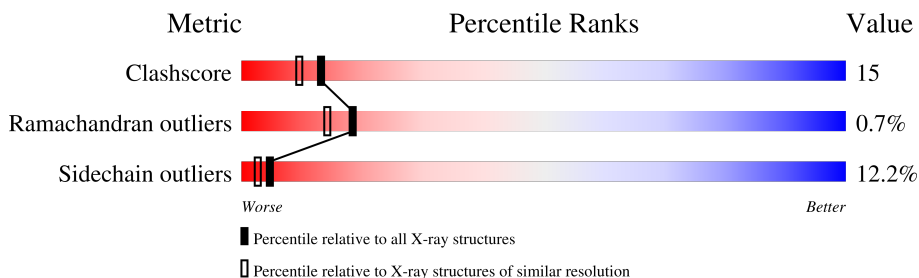
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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))
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	437	

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
2	COA	A	438	X	-	-	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 3598 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.

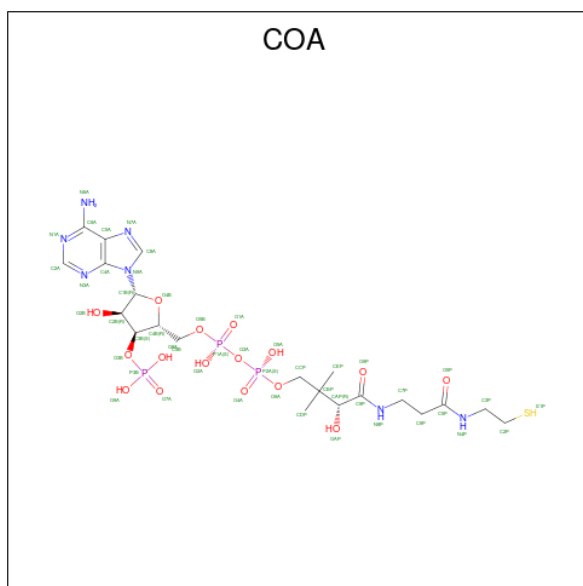
- Molecule 1 is a protein called CITRATE SYNTHASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	437	3442	2198	591	634	19	132	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	32	ALA	VAL	conflict	UNP P00889

- Molecule 2 is COENZYME A (CCD ID: COA) (formula: $C_{21}H_{36}N_7O_{16}P_3S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
2	A	1	48	21	7	16	3	6	0

- Molecule 3 is CITRIC ACID (CCD ID: CIT) (formula: $C_6H_8O_7$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			13	6	7		

- Molecule 4 is water.

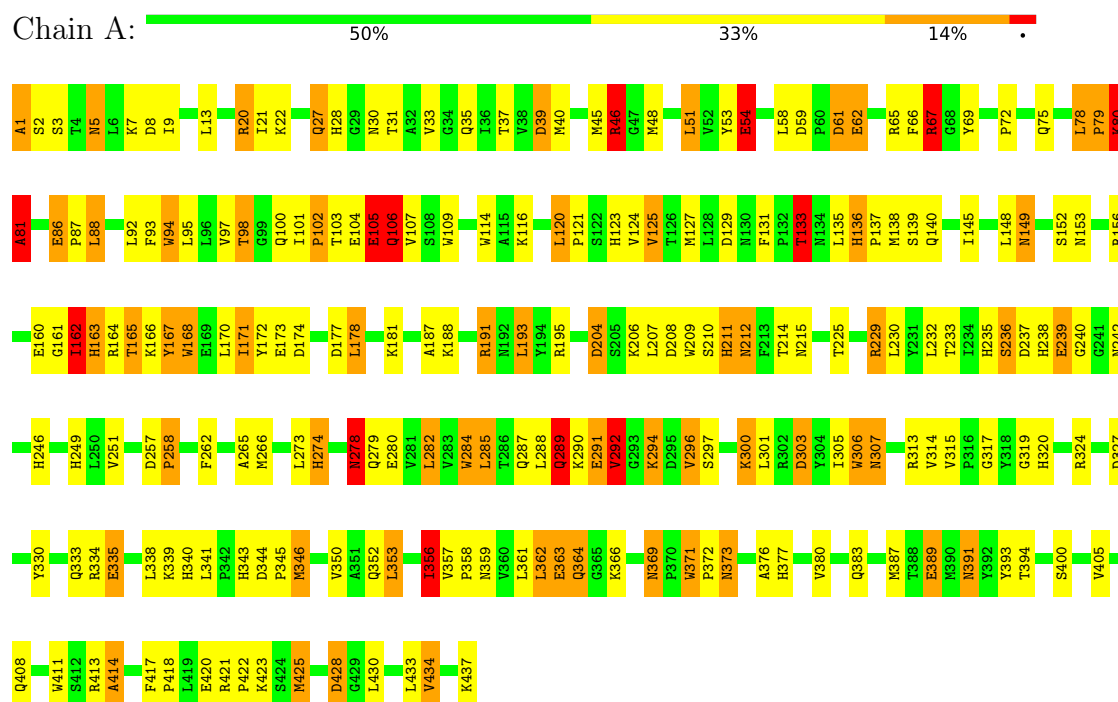
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	95	Total	O	0	0
			95	95		

3 Residue-property plots

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. 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. 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.

Note EDS was not executed.

• Molecule 1: CITRATE SYNTHASE



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	104.14Å 78.25Å 58.40Å 90.00° 78.50° 90.00°	Depositor
Resolution (Å)	5.00 – 2.00	Depositor
% Data completeness (in resolution range)	(Not available) (5.00-2.00)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	EREF	Depositor
R, R_{free}	0.161 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3598	wwPDB-VP
Average B, all atoms (Å ²)	14.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

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

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	1.47	40/3526 (1.1%)	1.91	77/4783 (1.6%)

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	1	48

All (40) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	114	TRP	NE1-CE2	-10.38	1.26	1.37
1	A	306	TRP	NE1-CE2	-9.85	1.26	1.37
1	A	94	TRP	NE1-CE2	-9.72	1.26	1.37
1	A	168	TRP	NE1-CE2	-9.66	1.26	1.37
1	A	209	TRP	NE1-CE2	-9.44	1.27	1.37
1	A	284	TRP	NE1-CE2	-9.26	1.27	1.37
1	A	274	HIS	ND1-CE1	9.21	1.41	1.32
1	A	238	HIS	CE1-NE2	9.05	1.41	1.32
1	A	211	HIS	ND1-CE1	9.04	1.41	1.32
1	A	28	HIS	ND1-CE1	9.01	1.41	1.32
1	A	320	HIS	CE1-NE2	8.97	1.41	1.32
1	A	411	TRP	NE1-CE2	-8.95	1.27	1.37
1	A	28	HIS	CE1-NE2	8.91	1.41	1.32
1	A	109	TRP	NE1-CE2	-8.87	1.27	1.37
1	A	163	HIS	ND1-CE1	8.78	1.41	1.32
1	A	371	TRP	NE1-CE2	-8.66	1.27	1.37
1	A	211	HIS	CE1-NE2	8.64	1.41	1.32
1	A	274	HIS	CE1-NE2	8.61	1.41	1.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	249	HIS	CE1-NE2	8.58	1.41	1.32
1	A	249	HIS	ND1-CE1	8.57	1.41	1.32
1	A	123	HIS	CE1-NE2	8.56	1.41	1.32
1	A	235	HIS	ND1-CE1	8.55	1.41	1.32
1	A	340	HIS	ND1-CE1	8.54	1.41	1.32
1	A	377	HIS	ND1-CE1	8.41	1.41	1.32
1	A	163	HIS	CE1-NE2	8.41	1.41	1.32
1	A	246	HIS	CE1-NE2	8.19	1.40	1.32
1	A	340	HIS	CE1-NE2	8.10	1.40	1.32
1	A	235	HIS	CE1-NE2	8.09	1.40	1.32
1	A	136	HIS	CE1-NE2	8.09	1.40	1.32
1	A	377	HIS	CE1-NE2	8.08	1.40	1.32
1	A	238	HIS	ND1-CE1	8.04	1.40	1.32
1	A	246	HIS	ND1-CE1	8.00	1.40	1.32
1	A	320	HIS	ND1-CE1	7.85	1.40	1.32
1	A	123	HIS	ND1-CE1	7.56	1.40	1.32
1	A	81	ALA	N-CA	-7.33	1.36	1.46
1	A	136	HIS	ND1-CE1	7.29	1.39	1.32
1	A	291	GLU	C-N	-6.78	1.25	1.34
1	A	80	LYS	C-N	-6.76	1.24	1.33
1	A	343	HIS	CE1-NE2	5.66	1.38	1.32
1	A	289	GLN	C-O	-5.64	1.17	1.24

All (77) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	80	LYS	CA-C-N	-15.29	100.80	122.06
1	A	80	LYS	C-N-CA	-15.29	100.80	122.06
1	A	300	LYS	CA-C-N	13.03	138.23	120.38
1	A	300	LYS	C-N-CA	13.03	138.23	120.38
1	A	46	ARG	NE-CZ-NH2	-9.90	110.29	119.20
1	A	39	ASP	CA-CB-CG	-9.56	103.04	112.60
1	A	30	ASN	OD1-CG-ND2	-8.42	114.18	122.60
1	A	162	ILE	CB-CA-C	-8.20	102.26	111.45
1	A	278	ASN	CB-CA-C	-7.98	98.24	110.92
1	A	131	PHE	O-C-N	7.23	127.73	121.30
1	A	81	ALA	CA-C-O	-7.21	112.86	122.51
1	A	173	GLU	CA-C-N	7.13	130.17	120.54
1	A	173	GLU	C-N-CA	7.13	130.17	120.54
1	A	20	ARG	CD-NE-CZ	7.07	134.29	124.40
1	A	131	PHE	CA-C-N	7.01	127.04	119.89
1	A	131	PHE	C-N-CA	7.01	127.04	119.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	279	GLN	O-C-N	6.96	130.08	122.15
1	A	46	ARG	CD-NE-CZ	6.83	133.95	124.40
1	A	251	VAL	CA-C-N	6.76	127.63	119.99
1	A	251	VAL	C-N-CA	6.76	127.63	119.99
1	A	343	HIS	ND1-CE1-NE2	-6.69	101.71	108.40
1	A	79	PRO	N-CA-CB	6.63	109.16	103.19
1	A	129	ASP	CA-CB-CG	-6.60	106.00	112.60
1	A	191	ARG	CD-NE-CZ	6.59	133.62	124.40
1	A	356	ILE	N-CA-C	6.55	117.20	111.56
1	A	417	PHE	CA-C-N	6.42	126.37	120.21
1	A	417	PHE	C-N-CA	6.42	126.37	120.21
1	A	78	LEU	CA-C-N	6.36	126.73	119.93
1	A	78	LEU	C-N-CA	6.36	126.73	119.93
1	A	278	ASN	O-C-N	6.21	128.73	122.08
1	A	356	ILE	N-CA-CB	-6.16	104.64	111.41
1	A	420	GLU	CA-C-N	-6.12	116.26	122.74
1	A	420	GLU	C-N-CA	-6.12	116.26	122.74
1	A	102	PRO	CA-C-N	6.10	131.73	122.65
1	A	102	PRO	C-N-CA	6.10	131.73	122.65
1	A	152	SER	N-CA-C	6.09	119.09	109.96
1	A	344	ASP	CA-CB-CG	-6.06	106.54	112.60
1	A	67	ARG	CD-NE-CZ	6.00	132.80	124.40
1	A	30	ASN	CB-CG-ND2	5.99	125.39	116.40
1	A	121	PRO	N-CA-CB	5.97	108.62	103.31
1	A	149	ASN	CA-C-N	5.88	128.44	120.38
1	A	149	ASN	C-N-CA	5.88	128.44	120.38
1	A	257	ASP	CA-C-N	5.82	125.50	119.56
1	A	257	ASP	C-N-CA	5.82	125.50	119.56
1	A	22	LYS	N-CA-CB	5.82	118.77	110.16
1	A	2	SER	CA-C-N	5.78	132.58	121.54
1	A	2	SER	C-N-CA	5.78	132.58	121.54
1	A	173	GLU	N-CA-C	5.77	117.25	111.07
1	A	61	ASP	CA-CB-CG	-5.72	106.88	112.60
1	A	27	GLN	N-CA-CB	-5.71	100.50	110.44
1	A	120	LEU	CA-C-N	5.61	125.54	119.76
1	A	120	LEU	C-N-CA	5.61	125.54	119.76
1	A	153	ASN	CA-CB-CG	-5.58	107.02	112.60
1	A	289	GLN	CG-CD-NE2	-5.48	108.18	116.40
1	A	54	GLU	CB-CG-CD	-5.45	103.33	112.60
1	A	177	ASP	CB-CA-C	-5.45	100.73	110.70
1	A	210	SER	O-C-N	5.42	128.33	122.15
1	A	434	VAL	N-CA-C	5.35	115.97	110.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	8	ASP	CA-CB-CG	-5.33	107.27	112.60
1	A	294	LYS	N-CA-C	5.30	122.09	110.80
1	A	87	PRO	N-CA-CB	5.26	107.88	103.25
1	A	102	PRO	N-CA-CB	5.24	108.31	103.39
1	A	389	GLU	CA-C-N	5.21	128.94	120.60
1	A	389	GLU	C-N-CA	5.21	128.94	120.60
1	A	59	ASP	CA-CB-CG	-5.21	107.39	112.60
1	A	320	HIS	CA-CB-CG	-5.19	108.61	113.80
1	A	104	GLU	CA-C-N	5.17	127.21	120.28
1	A	104	GLU	C-N-CA	5.17	127.21	120.28
1	A	72	PRO	N-CA-CB	5.16	109.14	103.26
1	A	422	PRO	N-CA-CB	5.14	108.65	103.25
1	A	51	LEU	CB-CA-C	-5.10	99.34	109.79
1	A	335	GLU	O-C-N	5.08	127.58	122.09
1	A	28	HIS	CA-CB-CG	-5.07	108.73	113.80
1	A	258	PRO	N-CA-CB	5.04	108.52	103.48
1	A	80	LYS	O-C-N	5.03	128.58	122.14
1	A	54	GLU	N-CA-C	5.02	118.55	112.23
1	A	173	GLU	O-C-N	5.01	127.23	122.07

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	214	THR	CB

All (48) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1	ALA	Mainchain
1	A	105	GLU	Sidechain
1	A	106	GLN	Sidechain
1	A	116	LYS	Mainchain
1	A	125	VAL	Mainchain
1	A	133	THR	Mainchain
1	A	145	ILE	Mainchain
1	A	162	ILE	Mainchain
1	A	167	TYR	Sidechain
1	A	181	LYS	Mainchain
1	A	204	ASP	Sidechain
1	A	208	ASP	Sidechain
1	A	237	ASP	Sidechain
1	A	239	GLU	Sidechain

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Mol	Chain	Res	Type	Group
1	A	242	ASN	Sidechain
1	A	278	ASN	Sidechain,Mainchain
1	A	280	GLU	Mainchain
1	A	287	GLN	Sidechain
1	A	289	GLN	Sidechain
1	A	290	LYS	Mainchain
1	A	292	VAL	Mainchain
1	A	303	ASP	Sidechain
1	A	307	ASN	Mainchain
1	A	315	VAL	Mainchain
1	A	333	GLN	Mainchain
1	A	35	GLN	Sidechain
1	A	359	ASN	Sidechain
1	A	363	GLU	Sidechain
1	A	364	GLN	Sidechain
1	A	369	ASN	Sidechain
1	A	373	ASN	Sidechain
1	A	39	ASP	Sidechain
1	A	391	ASN	Sidechain
1	A	414	ALA	Mainchain
1	A	418	PRO	Mainchain
1	A	421	ARG	Sidechain
1	A	428	ASP	Sidechain
1	A	46	ARG	Sidechain
1	A	51	LEU	Mainchain
1	A	54	GLU	Mainchain
1	A	61	ASP	Sidechain
1	A	66	PHE	Mainchain
1	A	67	ARG	Mainchain,Peptide
1	A	81	ALA	Mainchain
1	A	86	GLU	Sidechain
1	A	9	ILE	Mainchain

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	3442	0	3432	99	12

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	48	0	32	1	0
3	A	13	0	5	1	0
4	A	95	0	0	3	4
All	All	3598	0	3469	99	12

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 15.

All (99) 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:86:GLU:HG2	1:A:230:LEU:HB2	1.28	1.07
1:A:350:VAL:HG21	1:A:380:VAL:HG21	1.39	1.04
1:A:67:ARG:HB3	1:A:69:TYR:CD1	2.19	0.78
1:A:81:ALA:HB2	1:A:88:LEU:CD1	2.15	0.76
1:A:81:ALA:HB2	1:A:88:LEU:HD13	1.69	0.74
1:A:366:LYS:HE2	2:A:438:COA:O3B	1.89	0.73
1:A:334:ARG:O	1:A:338:LEU:HG	1.90	0.71
1:A:103:THR:HB	1:A:106:GLN:HG2	1.71	0.70
1:A:67:ARG:HB3	1:A:69:TYR:HD1	1.57	0.68
1:A:79:PRO:HG2	1:A:107:VAL:HG21	1.76	0.68
1:A:137:PRO:HG2	1:A:391:ASN:O	1.93	0.67
1:A:300:LYS:O	1:A:303:ASP:HB2	1.95	0.66
1:A:350:VAL:CG2	1:A:380:VAL:HG21	2.23	0.65
1:A:425:MET:HE1	1:A:433:LEU:HD22	1.78	0.64
1:A:285:LEU:HB3	1:A:346:MET:CE	2.28	0.63
1:A:187:ALA:O	1:A:191:ARG:HB2	2.00	0.62
1:A:62:GLU:HG2	1:A:65:ARG:HH22	1.65	0.60
1:A:352:GLN:O	1:A:356:ILE:HD12	2.01	0.60
1:A:67:ARG:H	1:A:69:TYR:H	1.49	0.59
1:A:163:HIS:ND1	1:A:165:THR:HB	2.17	0.58
1:A:233:THR:O	1:A:236:SER:HB2	2.04	0.57
1:A:380:VAL:HG23	4:A:484:HOH:O	2.03	0.57
1:A:288:LEU:HD12	1:A:292:VAL:HB	1.87	0.57
1:A:86:GLU:HG2	1:A:230:LEU:CB	2.19	0.56
1:A:162:ILE:HD13	1:A:170:LEU:HD13	1.86	0.56
1:A:92:LEU:HD23	1:A:236:SER:OG	2.06	0.55
1:A:94:TRP:HB3	1:A:102:PRO:HG3	1.89	0.55
1:A:274:HIS:CE1	3:A:439:CIT:H22	2.41	0.55
1:A:78:LEU:HB3	4:A:453:HOH:O	2.06	0.55
1:A:301:LEU:HD23	1:A:356:ILE:HD13	1.88	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:296:VAL:HG12	1:A:300:LYS:HB3	1.89	0.54
1:A:285:LEU:HB3	1:A:346:MET:HE2	1.89	0.54
1:A:5:ASN:C	1:A:5:ASN:HD22	2.14	0.54
1:A:306:TRP:HZ2	1:A:363:GLU:HG2	1.73	0.53
1:A:211:HIS:HA	1:A:214:THR:HG22	1.91	0.53
1:A:317:GLY:O	1:A:376:ALA:HB2	2.09	0.53
1:A:69:TYR:CE2	1:A:101:ILE:HD11	2.44	0.53
1:A:357:VAL:HB	1:A:358:PRO:HD3	1.90	0.52
1:A:204:ASP:H	1:A:212:ASN:HD21	1.57	0.52
1:A:81:ALA:HB2	1:A:88:LEU:HD11	1.90	0.51
1:A:124:VAL:HG21	1:A:148:LEU:HD23	1.92	0.51
1:A:301:LEU:CD2	1:A:356:ILE:HD13	2.40	0.51
1:A:297:SER:O	1:A:301:LEU:HB2	2.10	0.51
1:A:282:LEU:HD22	1:A:393:TYR:HE2	1.76	0.51
1:A:327:ASP:HB2	1:A:373:ASN:HA	1.91	0.50
1:A:371:TRP:HB3	1:A:372:PRO:HD2	1.93	0.50
1:A:358:PRO:O	1:A:362:LEU:HB2	2.11	0.50
1:A:164:ARG:HA	1:A:167:TYR:CE1	2.47	0.50
1:A:335:GLU:O	1:A:339:LYS:HE2	2.11	0.50
1:A:288:LEU:CD1	1:A:292:VAL:HG11	2.42	0.49
1:A:357:VAL:HG12	1:A:361:LEU:HD12	1.94	0.49
1:A:282:LEU:HD22	1:A:393:TYR:CE2	2.48	0.49
1:A:266:MET:HA	1:A:266:MET:HE2	1.95	0.49
1:A:171:ILE:HG13	1:A:258:PRO:HG3	1.95	0.48
1:A:53:TYR:CE2	1:A:408:GLN:HG2	2.49	0.48
1:A:168:TRP:CB	1:A:414:ALA:HB2	2.44	0.47
1:A:301:LEU:O	1:A:305:ILE:HG13	2.14	0.47
1:A:301:LEU:HD21	1:A:353:LEU:HD12	1.95	0.47
1:A:92:LEU:HD22	1:A:233:THR:HG23	1.97	0.46
1:A:136:HIS:HD2	1:A:138:MET:H	1.63	0.46
1:A:225:THR:O	1:A:229:ARG:HG2	2.16	0.46
1:A:330:TYR:CD2	1:A:372:PRO:HB2	2.50	0.46
1:A:53:TYR:CD2	1:A:240:GLY:HA3	2.51	0.46
1:A:37:THR:O	1:A:40:MET:HB2	2.17	0.45
1:A:262:PHE:O	1:A:265:ALA:HB3	2.17	0.45
1:A:125:VAL:HG13	1:A:188:LYS:HE2	1.99	0.45
1:A:161:GLY:HA2	4:A:513:HOH:O	2.17	0.45
1:A:288:LEU:HD11	1:A:292:VAL:HG11	1.98	0.44
1:A:324:ARG:HH11	1:A:369:ASN:HB2	1.83	0.44
1:A:5:ASN:C	1:A:5:ASN:ND2	2.76	0.44
1:A:135:LEU:HD23	1:A:140:GLN:HG2	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:166:LYS:O	1:A:170:LEU:HD12	2.19	0.43
1:A:215:ASN:HD22	1:A:215:ASN:HA	1.58	0.43
1:A:124:VAL:HG21	1:A:148:LEU:CD2	2.49	0.43
1:A:278:ASN:HD22	1:A:394:THR:HA	1.83	0.42
1:A:207:LEU:HD22	1:A:211:HIS:ND1	2.33	0.42
1:A:288:LEU:HD12	1:A:292:VAL:CB	2.50	0.42
1:A:324:ARG:HD3	1:A:324:ARG:HA	1.87	0.41
1:A:327:ASP:O	1:A:330:TYR:HB3	2.20	0.41
1:A:232:LEU:HA	1:A:400:SER:HB2	2.02	0.41
1:A:178:LEU:HD12	1:A:178:LEU:O	2.20	0.41
1:A:193:LEU:HD12	1:A:193:LEU:HA	1.86	0.41
1:A:284:TRP:HZ3	1:A:353:LEU:HD21	1.86	0.41
1:A:430:LEU:HD23	1:A:430:LEU:HA	1.87	0.41
1:A:53:TYR:CD2	1:A:408:GLN:HG2	2.55	0.41
1:A:94:TRP:O	1:A:98:THR:HB	2.21	0.41
1:A:133:THR:HA	1:A:193:LEU:HD11	2.02	0.41
1:A:364:GLN:HE21	1:A:364:GLN:HB3	1.44	0.41
1:A:7:LYS:HB2	1:A:172:TYR:CE2	2.56	0.41
1:A:45:MET:HB3	1:A:48:MET:HB2	2.03	0.41
1:A:80:LYS:HB3	1:A:81:ALA:H	1.55	0.41
1:A:156:ARG:O	1:A:160:GLU:HG3	2.21	0.41
1:A:346:MET:O	1:A:350:VAL:HG23	2.22	0.40
1:A:93:PHE:O	1:A:97:VAL:HG23	2.22	0.40
1:A:135:LEU:HD11	1:A:139:SER:HB3	2.02	0.40
1:A:174:ASP:HB2	1:A:258:PRO:HG2	2.03	0.40
1:A:21:ILE:HD13	1:A:21:ILE:HG21	1.86	0.40
1:A:95:LEU:HA	1:A:100:GLN:O	2.21	0.40
1:A:319:GLY:HA2	1:A:369:ASN:O	2.21	0.40

All (12) 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	Interatomic distance (Å)	Clash overlap (Å)
1:A:1:ALA:N	4:A:528:HOH:O[1_554]	0.85	1.35
1:A:3:SER:OG	1:A:291:GLU:OE2[1_554]	1.38	0.82
1:A:292:VAL:CG2	4:A:490:HOH:O[1_556]	1.40	0.80
1:A:1:ALA:CA	4:A:528:HOH:O[1_554]	1.54	0.66
1:A:127:MET:CE	1:A:127:MET:CE[2_555]	1.83	0.37
1:A:3:SER:O	1:A:313:ARG:NH2[1_554]	1.94	0.26
1:A:1:ALA:CB	4:A:528:HOH:O[1_554]	1.98	0.22

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:31:THR:CG2	1:A:434:VAL:CG2[2_555]	2.09	0.11
1:A:105:GLU:OE1	1:A:291:GLU:CG[1_554]	2.11	0.09
1:A:58:LEU:O	1:A:423:LYS:NZ[2_555]	2.12	0.08
1:A:45:MET:CE	1:A:45:MET:CE[2_555]	2.14	0.06
1:A:345:PRO:CG	1:A:437:LYS:NZ[3_555]	2.15	0.05

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	435/437 (100%)	408 (94%)	24 (6%)	3 (1%)	18 14

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	294	LYS
1	A	239	GLU
1	A	5	ASN

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	370/370 (100%)	325 (88%)	45 (12%)	5 3

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	LEU
1	A	20	ARG
1	A	27	GLN
1	A	33	VAL
1	A	46	ARG
1	A	54	GLU
1	A	62	GLU
1	A	75	GLN
1	A	80	LYS
1	A	88	LEU
1	A	98	THR
1	A	105	GLU
1	A	106	GLN
1	A	120	LEU
1	A	133	THR
1	A	149	ASN
1	A	165	THR
1	A	171	ILE
1	A	178	LEU
1	A	193	LEU
1	A	195	ARG
1	A	206	LYS
1	A	212	ASN
1	A	229	ARG
1	A	236	SER
1	A	273	LEU
1	A	282	LEU
1	A	285	LEU
1	A	289	GLN
1	A	292	VAL
1	A	296	VAL
1	A	307	ASN
1	A	314	VAL
1	A	341	LEU
1	A	346	MET
1	A	353	LEU
1	A	356	ILE
1	A	362	LEU
1	A	383	GLN
1	A	387	MET
1	A	389	GLU
1	A	405	VAL
1	A	413	ARG

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Mol	Chain	Res	Type
1	A	425	MET
1	A	428	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	5	ASN
1	A	27	GLN
1	A	28	HIS
1	A	130	ASN
1	A	136	HIS
1	A	140	GLN
1	A	192	ASN
1	A	212	ASN
1	A	215	ASN
1	A	223	GLN
1	A	267	ASN
1	A	289	GLN
1	A	310	ASN
1	A	352	GLN
1	A	359	ASN
1	A	364	GLN
1	A	383	GLN
1	A	391	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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	COA	A	438	-	47,50,50	2.10	10 (21%)	69,75,75	2.03	25 (36%)
3	CIT	A	439	-	12,12,12	1.29	1 (8%)	17,17,17	1.75	3 (17%)

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	COA	A	438	-	1/1/11/13	22/48/64/64	0/3/3/3
3	CIT	A	439	-	-	6/16/16/16	-

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	438	COA	P2A-O3A	6.42	1.66	1.59
2	A	438	COA	C5A-N7A	-6.41	1.27	1.39
2	A	438	COA	C4A-N9A	-5.62	1.26	1.37
2	A	438	COA	C8A-N7A	3.87	1.39	1.31
2	A	438	COA	C4A-N3A	-3.42	1.28	1.34
2	A	438	COA	C6A-N1A	-3.25	1.26	1.35
2	A	438	COA	P3B-O3B	2.68	1.64	1.59
2	A	438	COA	O9P-C9P	2.47	1.28	1.23
3	A	439	CIT	C2-C3	2.39	1.56	1.54
2	A	438	COA	P3B-O7A	2.15	1.57	1.50
2	A	438	COA	P1A-O3A	2.01	1.61	1.59

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	438	COA	N9A-C8A-N7A	-6.35	104.93	113.94
3	A	439	CIT	O6-C6-C3	5.31	123.33	113.14
2	A	438	COA	C3B-C2B-C1B	4.48	109.74	99.89
2	A	438	COA	C5A-C4A-N3A	-4.31	120.78	126.72
2	A	438	COA	O9P-C9P-N8P	-3.67	115.23	122.98
2	A	438	COA	C6A-C5A-C4A	3.65	122.16	117.18
2	A	438	COA	C4A-N9A-C1B	-3.54	118.35	126.63
2	A	438	COA	C5A-N7A-C8A	3.53	109.00	103.45
2	A	438	COA	C4A-N9A-C8A	3.26	109.16	105.74
2	A	438	COA	C4B-O4B-C1B	2.90	115.88	109.47
2	A	438	COA	C5B-C4B-C3B	-2.80	105.05	114.38
2	A	438	COA	C5A-C4A-N9A	2.74	108.80	105.81
2	A	438	COA	N6A-C6A-N1A	2.53	124.00	118.38
2	A	438	COA	CEP-CBP-CAP	2.51	113.05	108.77
2	A	438	COA	C1B-N9A-C8A	2.42	132.47	127.09
2	A	438	COA	CDP-CBP-CAP	-2.39	104.69	108.77
2	A	438	COA	O9P-C9P-CAP	2.36	127.47	120.89
2	A	438	COA	C2A-N1A-C6A	2.30	122.50	118.73
2	A	438	COA	OAP-CAP-C9P	2.26	118.83	109.38
2	A	438	COA	O4B-C1B-N9A	2.23	112.37	108.09
2	A	438	COA	O4B-C1B-C2B	-2.21	101.88	106.62
3	A	439	CIT	O6-C6-O5	-2.21	116.80	123.86
2	A	438	COA	C4A-C5A-N7A	-2.17	108.10	110.58
2	A	438	COA	C6P-C7P-N8P	-2.12	107.50	112.00
2	A	438	COA	C7P-N8P-C9P	2.09	126.30	122.55
2	A	438	COA	O2B-C2B-C3B	-2.07	105.38	111.19
2	A	438	COA	O5P-C5P-N4P	2.04	127.03	123.03
3	A	439	CIT	O4-C5-C4	2.02	120.74	114.35

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	438	COA	CAP

All (28) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	438	COA	CDP-CBP-CCP-O6A
2	A	438	COA	CEP-CBP-CCP-O6A
2	A	438	COA	CAP-CBP-CCP-O6A
2	A	438	COA	OAP-CAP-CBP-CCP
2	A	438	COA	C9P-CAP-CBP-CCP
2	A	438	COA	OAP-CAP-CBP-CDP

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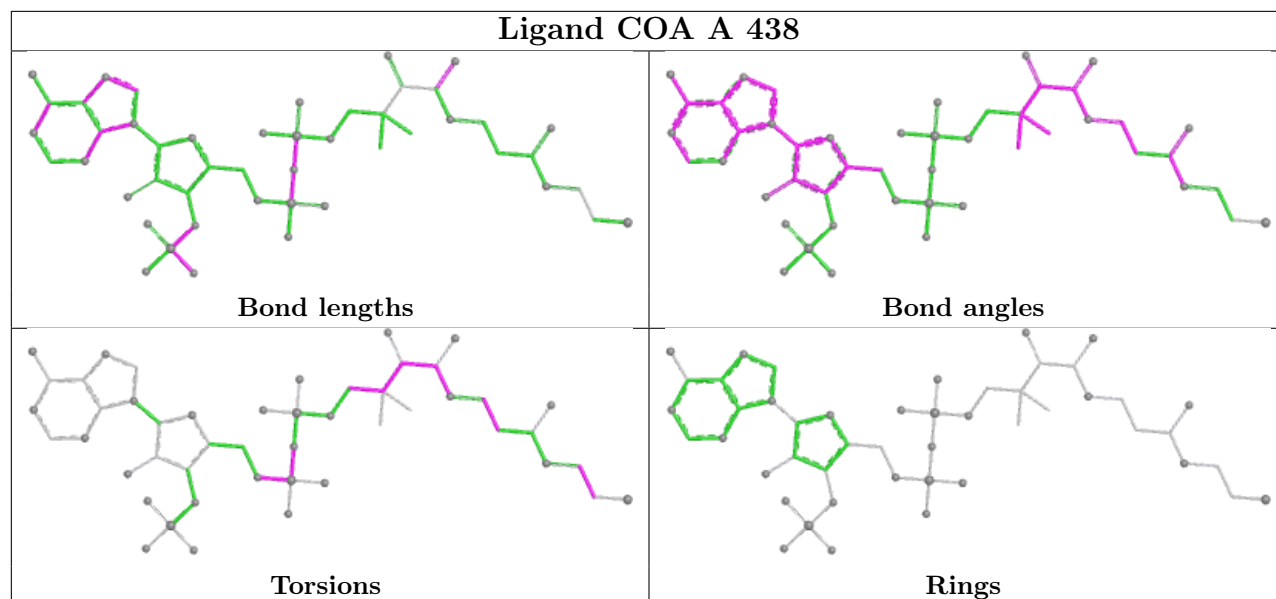
Mol	Chain	Res	Type	Atoms
2	A	438	COA	C9P-CAP-CBP-CDP
2	A	438	COA	OAP-CAP-CBP-CEP
2	A	438	COA	C9P-CAP-CBP-CEP
2	A	438	COA	CAP-C9P-N8P-C7P
2	A	438	COA	C5P-C6P-C7P-N8P
3	A	439	CIT	C2-C3-C6-O5
3	A	439	CIT	C2-C3-C6-O6
3	A	439	CIT	O7-C3-C6-O5
3	A	439	CIT	O7-C3-C6-O6
2	A	438	COA	O9P-C9P-N8P-C7P
2	A	438	COA	O9P-C9P-CAP-OAP
2	A	438	COA	C5B-O5B-P1A-O1A
2	A	438	COA	C5B-O5B-P1A-O2A
2	A	438	COA	C5B-O5B-P1A-O3A
2	A	438	COA	P2A-O3A-P1A-O2A
2	A	438	COA	O9P-C9P-CAP-CBP
2	A	438	COA	N8P-C9P-CAP-CBP
3	A	439	CIT	C2-C3-C4-C5
2	A	438	COA	P2A-O3A-P1A-O1A
3	A	439	CIT	C4-C3-C6-O6
2	A	438	COA	N8P-C9P-CAP-OAP
2	A	438	COA	S1P-C2P-C3P-N4P

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	438	COA	1	0
3	A	439	CIT	1	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

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.