



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 12:56 PM UTC

PDB ID : 6CTS / pdb_00006cts
Title : PROPOSED MECHANISM FOR THE CONDENSATION REACTION OF CITRATE SYNTHASE. 1.9-ANGSTROMS STRUCTURE OF THE TERNARY COMPLEX WITH OXALOACETATE AND CARBOXYMETHYL COENZYME A
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Deposited on : 1989-11-16
Resolution : 2.50 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)
Xtrriage (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

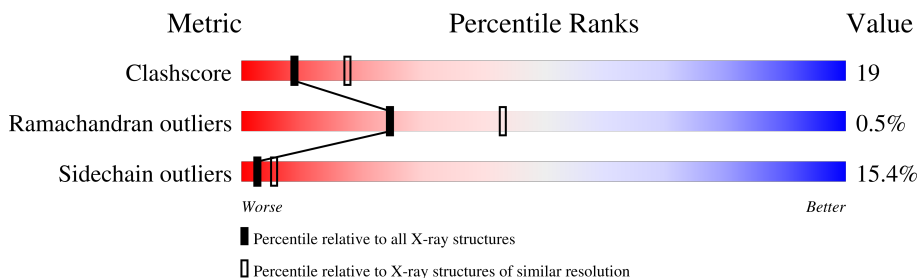
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)

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	433	 54% 33% 9% . .

2 Entry composition [i](#)

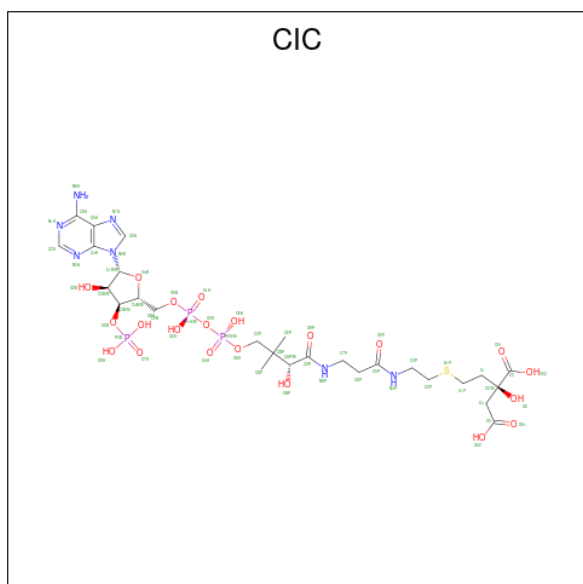
There are 3 unique types of molecules in this entry. The entry contains 3466 atoms, of which 0 are hydrogens and 0 are deuteriums.

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	429	3306	2115	571	603	17	0	0	0

- Molecule 2 is CITRYL-THIOETHER-COENZYME *A (CCD ID: CIC) (formula: $C_{27}H_{44}N_7O_{21}P_3S$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
			Total	C	N	O	P	S		
2	A	1	59	27	7	21	3	1	0	0

- Molecule 3 is water.

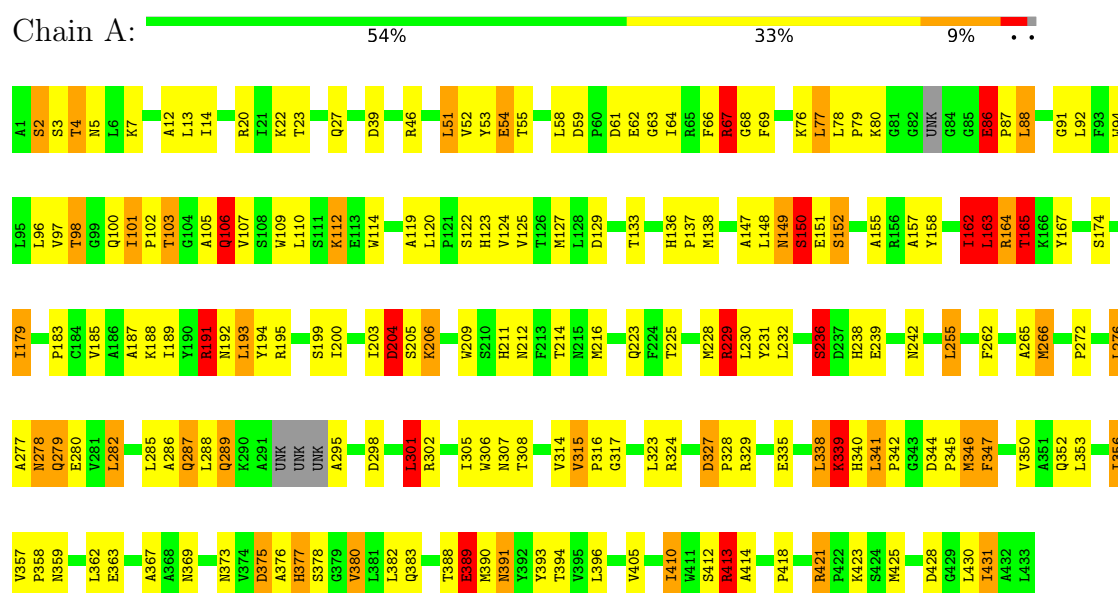
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	101	Total	O	0	0
			101	101		

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	104.00Å 78.10Å 58.30Å 90.00° 78.90° 90.00°	Depositor
Resolution (Å)	6.00 – 2.50	Depositor
% Data completeness (in resolution range)	(Not available) (6.00-2.50)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	TNT	Depositor
R, R_{free}	0.110 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3466	wwPDB-VP
Average B, all atoms (Å ²)	18.0	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: CIC

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.95	5/3386 (0.1%)	1.99	117/4598 (2.5%)

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	1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	214	THR	CB-OG1	6.01	1.53	1.43
1	A	103	THR	CB-OG1	5.91	1.53	1.43
1	A	315	VAL	N-CA	-5.89	1.42	1.46
1	A	133	THR	CB-OG1	5.56	1.52	1.43
1	A	98	THR	CB-OG1	5.46	1.52	1.43

All (117) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	229	ARG	CD-NE-CZ	12.36	141.70	124.40
1	A	431	ILE	N-CA-CB	11.76	123.57	110.51
1	A	191	ARG	CD-NE-CZ	10.06	138.48	124.40
1	A	302	ARG	CD-NE-CZ	-9.90	110.53	124.40
1	A	229	ARG	NE-CZ-NH2	-9.66	110.50	119.20
1	A	39	ASP	CA-CB-CG	-9.54	103.06	112.60
1	A	12	ALA	CA-C-N	9.25	132.87	120.65
1	A	12	ALA	C-N-CA	9.25	132.87	120.65

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	3	SER	N-CA-C	8.66	121.50	109.29
1	A	22	LYS	N-CA-CB	8.48	122.30	109.91
1	A	204	ASP	CA-CB-CG	8.47	121.07	112.60
1	A	359	ASN	CA-CB-CG	8.42	121.02	112.60
1	A	315	VAL	CA-C-O	8.40	123.82	119.12
1	A	66	PHE	CB-CA-C	8.12	124.29	111.39
1	A	289	GLN	CB-CG-CD	-8.05	98.91	112.60
1	A	431	ILE	CB-CA-C	7.71	121.74	111.88
1	A	22	LYS	N-CA-C	7.66	119.32	110.97
1	A	229	ARG	NE-CZ-NH1	7.63	129.13	121.50
1	A	51	LEU	CB-CA-C	-7.63	94.16	109.79
1	A	391	ASN	CA-CB-CG	7.61	120.20	112.60
1	A	340	HIS	N-CA-C	7.51	122.57	113.41
1	A	106	GLN	N-CA-CB	7.36	121.05	110.16
1	A	278	ASN	CA-CB-CG	-7.09	105.51	112.60
1	A	54	GLU	CB-CG-CD	-6.89	100.89	112.60
1	A	378	SER	N-CA-C	6.85	118.63	111.03
1	A	162	ILE	CB-CA-C	-6.80	101.08	111.08
1	A	308	THR	N-CA-CB	6.75	119.80	110.01
1	A	112	LYS	N-CA-CB	6.72	119.76	110.01
1	A	298	ASP	CA-CB-CG	6.67	119.27	112.60
1	A	64	ILE	N-CA-C	6.62	118.98	109.51
1	A	2	SER	N-CA-C	6.54	124.74	110.80
1	A	431	ILE	CA-CB-CG1	6.50	121.45	110.40
1	A	174	SER	CB-CA-C	6.46	123.07	110.67
1	A	191	ARG	NE-CZ-NH2	-6.38	113.46	119.20
1	A	418	PRO	CB-CA-C	-6.36	100.40	110.96
1	A	375	ASP	CA-CB-CG	6.27	118.87	112.60
1	A	388	THR	CA-CB-OG1	-6.25	100.22	109.60
1	A	4	THR	CB-CA-C	-6.23	99.68	109.84
1	A	106	GLN	CB-CA-C	-6.17	100.36	110.85
1	A	327	ASP	CA-C-N	6.14	125.82	119.56
1	A	327	ASP	C-N-CA	6.14	125.82	119.56
1	A	152	SER	N-CA-CB	6.13	119.20	109.51
1	A	286	ALA	CA-C-N	6.12	129.61	120.31
1	A	286	ALA	C-N-CA	6.12	129.61	120.31
1	A	278	ASN	CB-CA-C	-6.09	100.74	110.84
1	A	431	ILE	CA-CB-CG2	6.08	120.84	110.50
1	A	67	ARG	N-CA-C	-6.05	97.92	110.80
1	A	12	ALA	CB-CA-C	6.01	121.70	110.70
1	A	78	LEU	CA-C-N	6.00	125.96	119.78
1	A	78	LEU	C-N-CA	6.00	125.96	119.78

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	27	GLN	CB-CG-CD	-5.91	102.55	112.60
1	A	295	ALA	CA-C-N	5.91	129.06	122.15
1	A	295	ALA	C-N-CA	5.91	129.06	122.15
1	A	428	ASP	O-C-N	5.86	128.11	122.07
1	A	101	ILE	N-CA-CB	5.86	115.53	110.08
1	A	191	ARG	NE-CZ-NH1	5.82	127.32	121.50
1	A	163	LEU	N-CA-CB	-5.80	101.06	109.83
1	A	149	ASN	CA-C-N	5.78	132.59	121.54
1	A	149	ASN	C-N-CA	5.78	132.59	121.54
1	A	344	ASP	CA-C-N	5.73	125.41	119.05
1	A	344	ASP	C-N-CA	5.73	125.41	119.05
1	A	66	PHE	O-C-N	-5.72	115.36	122.11
1	A	412	SER	CA-C-N	5.71	128.40	120.29
1	A	412	SER	C-N-CA	5.71	128.40	120.29
1	A	77	LEU	N-CA-C	5.70	120.37	112.90
1	A	165	THR	CA-CB-CG2	5.70	120.18	110.50
1	A	165	THR	N-CA-CB	-5.70	101.14	110.14
1	A	191	ARG	N-CA-CB	-5.69	101.82	110.07
1	A	209	TRP	CA-C-N	5.69	128.22	120.54
1	A	209	TRP	C-N-CA	5.69	128.22	120.54
1	A	380	VAL	O-C-N	5.63	127.33	121.87
1	A	193	LEU	N-CA-CB	-5.62	101.92	110.07
1	A	119	ALA	CB-CA-C	5.61	119.33	109.80
1	A	382	LEU	CA-C-N	5.56	127.73	120.28
1	A	382	LEU	C-N-CA	5.56	127.73	120.28
1	A	157	ALA	CA-C-N	5.54	127.64	120.44
1	A	157	ALA	C-N-CA	5.54	127.64	120.44
1	A	223	GLN	N-CA-CB	5.53	118.86	110.28
1	A	53	TYR	CA-C-N	-5.53	112.35	121.92
1	A	53	TYR	C-N-CA	-5.53	112.35	121.92
1	A	54	GLU	N-CA-C	5.51	120.28	113.50
1	A	86	GLU	N-CA-C	5.50	117.90	110.31
1	A	52	VAL	N-CA-CB	5.50	117.64	111.21
1	A	255	LEU	N-CA-CB	5.46	119.71	110.49
1	A	191	ARG	CB-CG-CD	5.44	123.82	111.30
1	A	68	GLY	N-CA-C	-5.43	100.31	113.18
1	A	67	ARG	CA-C-N	5.37	131.93	121.41
1	A	67	ARG	C-N-CA	5.37	131.93	121.41
1	A	97	VAL	N-CA-C	5.37	115.57	110.53
1	A	122	SER	N-CA-C	5.36	117.81	111.33
1	A	377	HIS	N-CA-C	5.36	120.68	113.72
1	A	61	ASP	CA-CB-CG	-5.35	107.25	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	347	PHE	N-CA-CB	5.32	117.90	109.82
1	A	150	SER	CB-CA-C	-5.30	99.87	110.42
1	A	59	ASP	CA-C-N	5.29	125.60	119.47
1	A	59	ASP	C-N-CA	5.29	125.60	119.47
1	A	80	LYS	CA-C-O	-5.26	115.23	121.60
1	A	345	PRO	N-CA-CB	5.25	109.11	103.33
1	A	338	LEU	N-CA-CB	5.25	117.93	110.06
1	A	191	ARG	CG-CD-NE	-5.24	100.48	112.00
1	A	339	LYS	CB-CA-C	-5.23	101.13	110.70
1	A	389	GLU	N-CA-CB	5.22	117.88	110.46
1	A	183	PRO	N-CA-CB	5.22	108.70	103.48
1	A	4	THR	N-CA-C	5.22	118.05	109.76
1	A	67	ARG	CA-C-O	5.21	127.95	120.51
1	A	410	ILE	CA-CB-CG1	5.16	119.17	110.40
1	A	413	ARG	CB-CG-CD	5.13	123.11	111.30
1	A	367	ALA	CB-CA-C	-5.13	101.90	109.90
1	A	327	ASP	CA-CB-CG	5.11	117.71	112.60
1	A	356	ILE	CB-CA-C	5.07	116.02	111.71
1	A	287	GLN	CB-CA-C	-5.07	100.94	110.67
1	A	301	LEU	CA-C-O	5.06	125.79	119.97
1	A	63	GLY	CA-C-N	5.05	127.46	120.49
1	A	63	GLY	C-N-CA	5.05	127.46	120.49
1	A	421	ARG	O-C-N	5.05	125.06	121.12
1	A	3	SER	CA-CB-OG	-5.01	101.08	111.10
1	A	286	ALA	CB-CA-C	5.00	120.38	110.42

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	431	ILE	CB

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	229	ARG	Sidechain

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	3306	0	3297	125	4
2	A	59	0	38	4	2
3	A	101	0	0	3	0
All	All	3466	0	3335	125	4

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 19.

All (125) 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	1.39	1.03
1:A:125:VAL:HG13	1:A:188:LYS:HE2	1.43	0.97
1:A:86:GLU:HG3	1:A:87:PRO:HD2	1.48	0.93
1:A:94:TRP:HB3	1:A:102:PRO:HG3	1.60	0.81
1:A:301:LEU:HD23	1:A:356:ILE:HD13	1.63	0.80
1:A:103:THR:H	1:A:106:GLN:HG2	1.48	0.78
1:A:14:ILE:HG12	1:A:414:ALA:HB1	1.66	0.75
1:A:67:ARG:HA	3:A:548:HOH:O	1.86	0.75
1:A:77:LEU:HB3	1:A:101:ILE:HD13	1.68	0.74
1:A:346:MET:HG2	1:A:380:VAL:HG22	1.69	0.74
1:A:305:ILE:HD13	1:A:357:VAL:HG22	1.71	0.72
1:A:92:LEU:HD23	1:A:236:SER:OG	1.89	0.71
1:A:352:GLN:O	1:A:356:ILE:HD12	1.90	0.71
1:A:194:TYR:CD2	1:A:389:GLU:HG3	2.26	0.71
1:A:136:HIS:HD2	1:A:138:MET:H	1.40	0.68
1:A:106:GLN:CA	1:A:106:GLN:HE21	2.06	0.68
1:A:77:LEU:HD13	1:A:101:ILE:CD1	2.24	0.67
1:A:301:LEU:O	1:A:305:ILE:HG13	1.94	0.67
1:A:136:HIS:CD2	1:A:137:PRO:HD2	2.30	0.67
1:A:350:VAL:HG21	1:A:380:VAL:HG21	1.77	0.66
1:A:86:GLU:CG	1:A:87:PRO:HD2	2.22	0.66
1:A:129:ASP:OD2	1:A:188:LYS:NZ	2.29	0.66
1:A:163:LEU:HD12	1:A:165:THR:H	1.62	0.65
1:A:335:GLU:O	1:A:339:LYS:HE2	1.98	0.64
1:A:266:MET:HA	1:A:266:MET:HE2	1.82	0.62
1:A:86:GLU:HG3	1:A:87:PRO:CD	2.26	0.60
1:A:327:ASP:OD2	1:A:329:ARG:HB2	2.01	0.60
1:A:88:LEU:HD12	1:A:229:ARG:HD2	1.82	0.59
1:A:301:LEU:HD23	1:A:356:ILE:CD1	2.31	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:77:LEU:HB3	1:A:101:ILE:CD1	2.33	0.58
1:A:282:LEU:HD22	1:A:393:TYR:HE2	1.67	0.58
1:A:185:VAL:O	1:A:189:ILE:HG13	2.05	0.56
1:A:136:HIS:CD2	1:A:138:MET:H	2.22	0.56
1:A:79:PRO:HG2	1:A:107:VAL:HG21	1.87	0.56
1:A:285:LEU:O	1:A:288:LEU:HB3	2.06	0.55
1:A:136:HIS:CG	1:A:137:PRO:HD2	2.41	0.55
1:A:77:LEU:HD13	1:A:101:ILE:HD11	1.87	0.55
1:A:204:ASP:H	1:A:212:ASN:HD21	1.54	0.54
1:A:341:LEU:N	1:A:342:PRO:HD3	2.22	0.54
1:A:163:LEU:HD12	1:A:164:ARG:N	2.23	0.54
1:A:277:ALA:HB3	1:A:375:ASP:OD1	2.08	0.54
1:A:279:GLN:HB2	1:A:390:MET:O	2.08	0.53
1:A:357:VAL:HB	1:A:358:PRO:HD3	1.90	0.53
1:A:103:THR:N	1:A:106:GLN:HG2	2.20	0.53
1:A:338:LEU:HD23	1:A:347:PHE:CE1	2.44	0.53
1:A:323:LEU:O	1:A:324:ARG:NH1	2.42	0.53
1:A:228:MET:HE3	1:A:232:LEU:HG	1.92	0.52
1:A:317:GLY:O	1:A:376:ALA:HB2	2.09	0.52
1:A:187:ALA:O	1:A:191:ARG:HB2	2.11	0.51
1:A:106:GLN:HE21	1:A:106:GLN:HA	1.76	0.51
1:A:123:HIS:HE1	1:A:147:ALA:O	1.93	0.51
1:A:225:THR:O	1:A:229:ARG:HG2	2.11	0.51
1:A:262:PHE:O	1:A:265:ALA:HB3	2.11	0.51
1:A:413:ARG:HD2	3:A:588:HOH:O	2.11	0.50
1:A:194:TYR:O	1:A:195:ARG:HD3	2.11	0.50
1:A:67:ARG:HB2	1:A:69:PHE:CD1	2.47	0.50
1:A:162:ILE:HD12	1:A:167:TYR:CE1	2.47	0.49
1:A:94:TRP:CD1	1:A:98:THR:HG1	2.30	0.49
1:A:338:LEU:HD23	1:A:347:PHE:HE1	1.77	0.49
1:A:282:LEU:HD22	1:A:393:TYR:CE2	2.47	0.49
1:A:163:LEU:CD1	1:A:164:ARG:N	2.76	0.49
1:A:306:TRP:HZ2	1:A:363:GLU:HG2	1.79	0.48
1:A:162:ILE:HD12	1:A:167:TYR:CD1	2.48	0.48
1:A:152:SER:HB3	1:A:155:ALA:HB3	1.96	0.48
1:A:77:LEU:HD13	1:A:101:ILE:HD12	1.95	0.47
1:A:106:GLN:CA	1:A:106:GLN:NE2	2.77	0.47
1:A:206:LYS:HE3	1:A:206:LYS:HB2	1.65	0.47
1:A:314:VAL:O	1:A:316:PRO:HD3	2.15	0.47
1:A:204:ASP:OD2	1:A:206:LYS:N	2.34	0.47
1:A:305:ILE:CD1	1:A:357:VAL:HG22	2.43	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:346:MET:HB3	1:A:380:VAL:HG13	1.97	0.46
1:A:162:ILE:CD1	1:A:167:TYR:CD1	2.99	0.46
1:A:272:PRO:HA	1:A:276:LEU:HB3	1.97	0.46
1:A:410:ILE:HD13	1:A:410:ILE:HG21	1.67	0.45
1:A:204:ASP:H	1:A:212:ASN:ND2	2.14	0.45
1:A:279:GLN:NE2	1:A:280:GLU:OE2	2.50	0.45
1:A:92:LEU:HD12	1:A:92:LEU:HA	1.54	0.44
1:A:164:ARG:H	1:A:164:ARG:HG3	1.62	0.44
1:A:347:PHE:HA	1:A:380:VAL:HG11	1.99	0.44
1:A:373:ASN:ND2	2:A:700:CIC:H31	2.32	0.44
1:A:103:THR:OG1	1:A:105:ALA:HB3	2.17	0.44
1:A:242:ASN:HB3	2:A:700:CIC:C5	2.47	0.44
1:A:327:ASP:HA	1:A:328:PRO:HD2	1.91	0.44
1:A:4:THR:HG22	1:A:5:ASN:N	2.32	0.44
1:A:150:SER:HB3	1:A:151:GLU:HG2	1.99	0.44
1:A:278:ASN:HD22	1:A:394:THR:HA	1.82	0.44
1:A:350:VAL:CG2	1:A:380:VAL:HG21	2.47	0.44
1:A:163:LEU:CD1	1:A:163:LEU:C	2.90	0.44
1:A:324:ARG:HA	1:A:324:ARG:HD3	1.68	0.44
1:A:204:ASP:OD2	1:A:204:ASP:C	2.60	0.43
1:A:106:GLN:HE21	1:A:106:GLN:N	2.16	0.43
1:A:192:ASN:HD22	1:A:192:ASN:HA	1.47	0.43
1:A:194:TYR:CG	1:A:389:GLU:HG3	2.54	0.43
1:A:339:LYS:HD3	1:A:339:LYS:HA	1.55	0.43
1:A:191:ARG:HD3	1:A:200:ILE:HG22	2.01	0.43
1:A:67:ARG:N	1:A:69:PHE:H	2.17	0.43
1:A:5:ASN:C	1:A:5:ASN:HD22	2.27	0.42
1:A:152:SER:HB3	1:A:155:ALA:CB	2.49	0.42
1:A:138:MET:HB3	1:A:138:MET:HE3	1.79	0.42
1:A:228:MET:HE1	1:A:396:LEU:HD22	2.02	0.42
1:A:109:TRP:CZ3	1:A:110:LEU:HD23	2.54	0.42
1:A:5:ASN:HD22	1:A:7:LYS:H	1.68	0.42
1:A:124:VAL:HG21	1:A:148:LEU:HD23	2.00	0.42
1:A:58:LEU:HD22	1:A:238:HIS:CE1	2.54	0.42
1:A:77:LEU:CB	1:A:101:ILE:HD13	2.45	0.42
1:A:86:GLU:HG2	1:A:230:LEU:HD13	2.02	0.42
1:A:231:TYR:C	1:A:231:TYR:CD1	2.97	0.42
1:A:114:TRP:CZ2	1:A:179:ILE:HG21	2.55	0.42
1:A:124:VAL:HG21	1:A:148:LEU:CD2	2.50	0.41
1:A:194:TYR:CE2	1:A:389:GLU:HG3	2.54	0.41
1:A:163:LEU:HD12	1:A:163:LEU:C	2.44	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:191:ARG:HD3	1:A:216:MET:O	2.20	0.41
1:A:211:HIS:HD2	1:A:211:HIS:O	2.02	0.41
1:A:91:GLY:HA3	3:A:514:HOH:O	2.21	0.41
1:A:137:PRO:HG2	1:A:391:ASN:O	2.21	0.41
1:A:101:ILE:HA	1:A:102:PRO:HD3	1.86	0.41
1:A:158:TYR:CD1	1:A:158:TYR:C	2.98	0.41
1:A:324:ARG:HH11	1:A:369:ASN:HB2	1.85	0.41
1:A:306:TRP:HZ2	1:A:363:GLU:CD	2.29	0.41
1:A:315:VAL:O	2:A:700:CIC:N6A	2.51	0.41
1:A:55:THR:HG22	1:A:96:LEU:O	2.22	0.40
1:A:341:LEU:N	1:A:342:PRO:CD	2.84	0.40
1:A:110:LEU:H	1:A:110:LEU:HG	1.72	0.40
1:A:373:ASN:HD21	2:A:700:CIC:H31	1.86	0.40
1:A:377:HIS:O	1:A:377:HIS:ND1	2.53	0.40

All (4) 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:127:MET:CE	1:A:127:MET:CE[2_555]	1.97	0.23
1:A:58:LEU:O	1:A:423:LYS:NZ[2_555]	2.06	0.14
1:A:164:ARG:NH2	2:A:700:CIC:O8A[2_555]	2.10	0.10
1:A:421:ARG:NH2	2:A:700:CIC:O51[2_555]	2.19	0.01

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	423/433 (98%)	402 (95%)	19 (4%)	2 (0%)	24 43

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	239	GLU
1	A	236	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	345/345 (100%)	292 (85%)	53 (15%)	2 5

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	SER
1	A	13	LEU
1	A	20	ARG
1	A	23	THR
1	A	46	ARG
1	A	51	LEU
1	A	54	GLU
1	A	62	GLU
1	A	67	ARG
1	A	76	LYS
1	A	86	GLU
1	A	88	LEU
1	A	100	GLN
1	A	106	GLN
1	A	112	LYS
1	A	120	LEU
1	A	149	ASN
1	A	150	SER
1	A	162	ILE
1	A	163	LEU
1	A	164	ARG
1	A	165	THR
1	A	179	ILE
1	A	191	ARG
1	A	193	LEU

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Mol	Chain	Res	Type
1	A	199	SER
1	A	203	ILE
1	A	204	ASP
1	A	205	SER
1	A	206	LYS
1	A	229	ARG
1	A	236	SER
1	A	255	LEU
1	A	266	MET
1	A	276	LEU
1	A	279	GLN
1	A	282	LEU
1	A	287	GLN
1	A	289	GLN
1	A	301	LEU
1	A	307	ASN
1	A	339	LYS
1	A	341	LEU
1	A	346	MET
1	A	353	LEU
1	A	362	LEU
1	A	383	GLN
1	A	389	GLU
1	A	405	VAL
1	A	413	ARG
1	A	425	MET
1	A	430	LEU
1	A	431	ILE

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	26	GLN
1	A	27	GLN
1	A	106	GLN
1	A	123	HIS
1	A	130	ASN
1	A	136	HIS
1	A	140	GLN
1	A	149	ASN
1	A	192	ASN

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Mol	Chain	Res	Type
1	A	211	HIS
1	A	212	ASN
1	A	215	ASN
1	A	267	ASN
1	A	287	GLN
1	A	310	ASN
1	A	352	GLN
1	A	383	GLN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

1 ligand is modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	CIC	A	700	-	59,61,61	1.71	8 (13%)	82,91,91	1.56	11 (13%)

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	CIC	A	700	-	-	8/65/81/81	0/3/3/3

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	700	CIC	P3B-O3B	6.57	1.71	1.59
2	A	700	CIC	P2A-O3A	5.55	1.65	1.59
2	A	700	CIC	C4-C2	-4.06	1.48	1.54
2	A	700	CIC	O9P-C9P	3.92	1.30	1.23
2	A	700	CIC	P3B-O7A	2.96	1.59	1.50
2	A	700	CIC	C5A-C6A	2.32	1.47	1.41
2	A	700	CIC	C2A-N1A	2.08	1.37	1.33
2	A	700	CIC	O4B-C4B	2.07	1.49	1.45

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	700	CIC	O8A-P3B-O3B	-5.71	83.61	105.85
2	A	700	CIC	C2-C4-C5	5.13	127.94	113.92
2	A	700	CIC	O2-C2-C4	4.32	119.23	109.38
2	A	700	CIC	C3P-N4P-C5P	-3.33	116.62	122.82
2	A	700	CIC	O4B-C1B-C2B	-3.13	99.91	106.62
2	A	700	CIC	O5A-P2A-O3A	2.89	115.07	107.27
2	A	700	CIC	O3B-P3B-O7A	2.86	119.55	109.33
2	A	700	CIC	O3B-C3B-C4B	2.83	119.99	110.03
2	A	700	CIC	CEP-CBP-CCP	-2.10	104.76	108.22
2	A	700	CIC	C2B-C3B-C4B	-2.08	99.59	103.24
2	A	700	CIC	C2A-N1A-C6A	2.06	122.12	118.73

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	700	CIC	C2-C1-C1P-S1P
2	A	700	CIC	C4-C2-C3-O3I
2	A	700	CIC	CEP-CBP-CCP-O6A
2	A	700	CIC	C3B-O3B-P3B-O8A
2	A	700	CIC	C1-C2-C3-O3I
2	A	700	CIC	C3P-C2P-S1P-C1P
2	A	700	CIC	CDP-CBP-CCP-O6A

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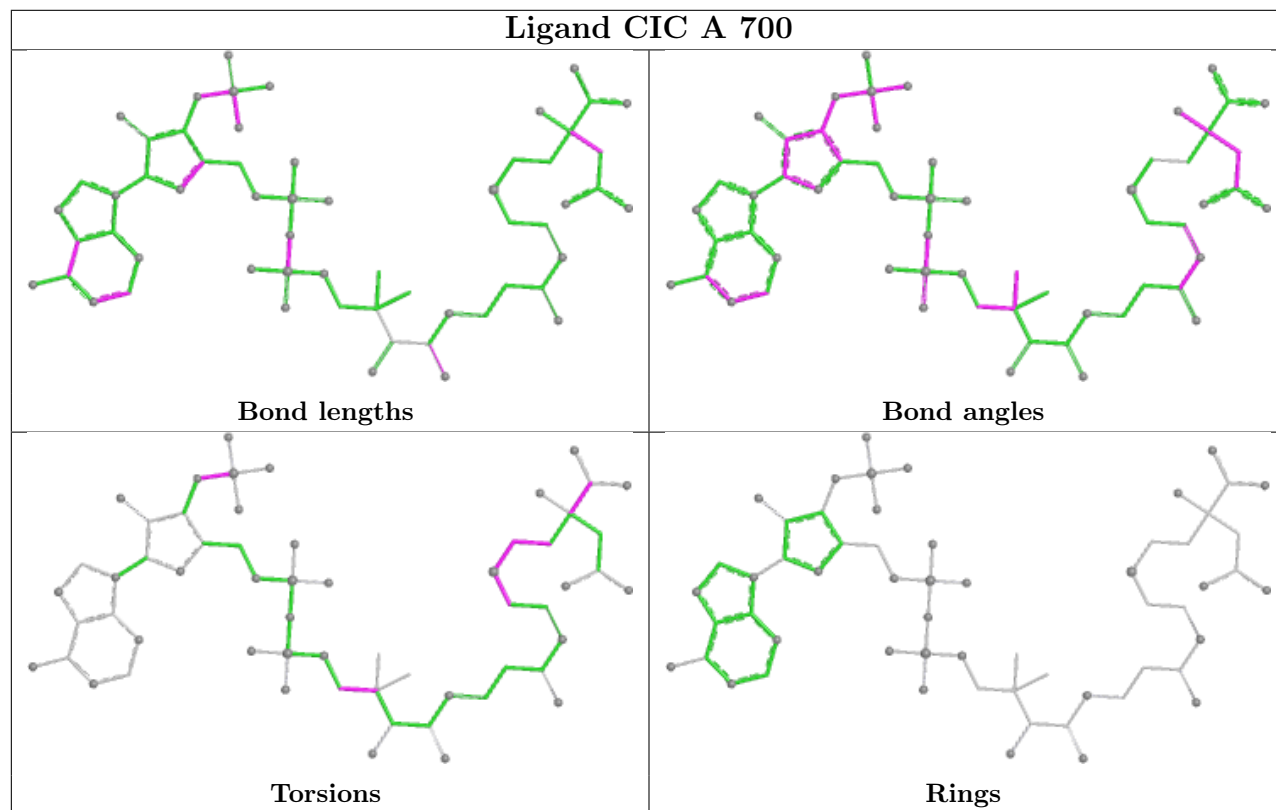
Mol	Chain	Res	Type	Atoms
2	A	700	CIC	C1-C1P-S1P-C2P

There are no ring outliers.

1 monomer is involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	700	CIC	4	2

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.