Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 01:22 AM GMT

PDB ID	:	2CTS
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
http://wwpdb.org/validation/2016/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

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

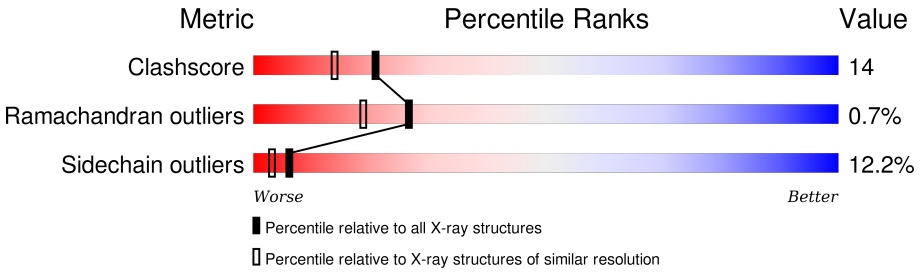
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

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	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (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 on the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria. 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 <=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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	437	57% 30% 10% •

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

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.

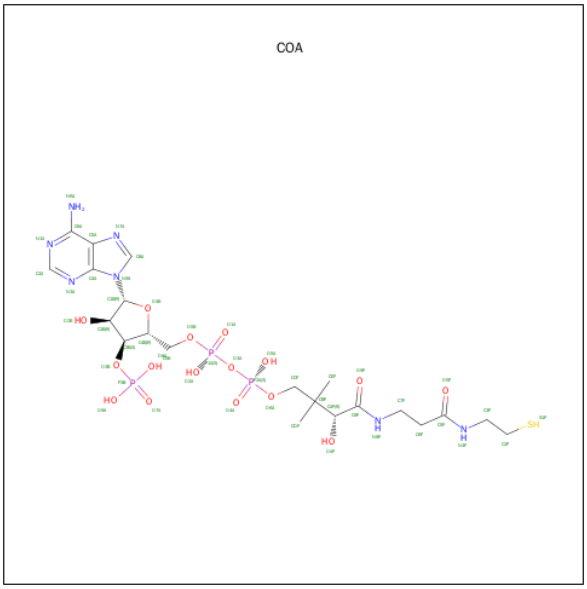
- Molecule 1 is a protein called CITRATE SYNTHASE.

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:

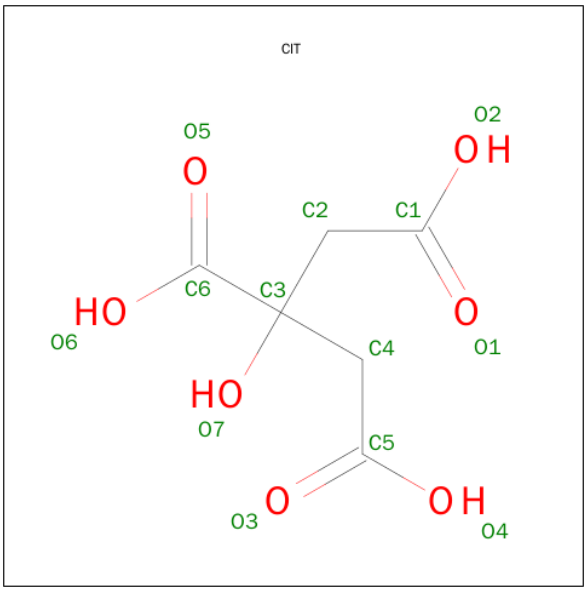
Chain	Residue	Modelled	Actual	Comment	Reference
A	32	ALA	VAL	CONFLICT	UNP P00889

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



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

- Molecule 3 is CITRIC ACID (three-letter code: CIT) (formula: $C_6H_8O_7$).



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

- Molecule 4 is water.

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

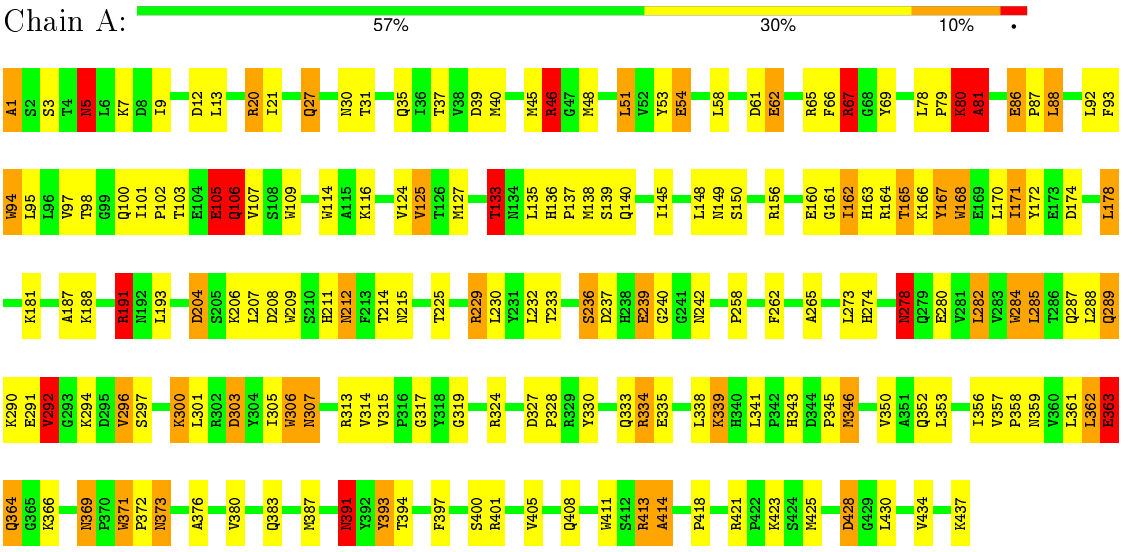
3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

Note EDS was not executed.

- Molecule 1: CITRATE SYNTHASE

Chain A:



4 Data and refinement statistics (i)

Xtriage (Phenix) and EDS were not executed - this section will therefore be 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	1.07	10/3526 (0.3%)	1.47	24/4783 (0.5%)

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	48

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	114	TRP	NE1-CE2	-8.86	1.26	1.37
1	A	306	TRP	NE1-CE2	-8.41	1.26	1.37
1	A	94	TRP	NE1-CE2	-8.30	1.26	1.37
1	A	168	TRP	NE1-CE2	-8.25	1.26	1.37
1	A	209	TRP	NE1-CE2	-8.06	1.27	1.37
1	A	284	TRP	NE1-CE2	-7.91	1.27	1.37
1	A	411	TRP	NE1-CE2	-7.65	1.27	1.37
1	A	109	TRP	NE1-CE2	-7.58	1.27	1.37
1	A	371	TRP	NE1-CE2	-7.41	1.27	1.37
1	A	81	ALA	N-CA	-5.14	1.36	1.46

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	46	ARG	NE-CZ-NH2	-20.03	110.29	120.30
1	A	20	ARG	NE-CZ-NH1	12.31	126.46	120.30
1	A	46	ARG	NE-CZ-NH1	10.36	125.48	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	20	ARG	NE-CZ-NH2	-9.35	115.62	120.30
1	A	191	ARG	NE-CZ-NH2	-9.10	115.75	120.30
1	A	67	ARG	NE-CZ-NH1	8.56	124.58	120.30
1	A	80	LYS	C-N-CA	-8.36	100.80	121.70
1	A	191	ARG	NE-CZ-NH1	7.95	124.28	120.30
1	A	20	ARG	CD-NE-CZ	7.64	134.29	123.60
1	A	12	ASP	CB-CG-OD2	7.51	125.06	118.30
1	A	46	ARG	CD-NE-CZ	7.40	133.95	123.60
1	A	191	ARG	CD-NE-CZ	7.16	133.62	123.60
1	A	300	LYS	C-N-CA	6.61	138.23	121.70
1	A	67	ARG	CD-NE-CZ	6.57	132.80	123.60
1	A	67	ARG	NE-CZ-NH2	-6.09	117.25	120.30
1	A	278	ASN	CB-CA-C	-6.08	98.24	110.40
1	A	334	ARG	NE-CZ-NH2	-6.01	117.29	120.30
1	A	51	LEU	CB-CA-C	-5.72	99.34	110.20
1	A	27	GLN	N-CA-CB	-5.61	100.50	110.60
1	A	413	ARG	NE-CZ-NH1	-5.52	117.54	120.30
1	A	393	TYR	CB-CG-CD2	-5.40	117.76	121.00
1	A	363	GLU	OE1-CD-OE2	-5.25	117.00	123.30
1	A	401	ARG	NE-CZ-NH1	5.23	122.92	120.30
1	A	150	SER	N-CA-CB	-5.13	102.81	110.50

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

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Mol	Chain	Res	Type	Group
1	A	237	ASP	Sidechain
1	A	239	GLU	Sidechain
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 (i)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3442	0	3432	96	12
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	96	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 14.

All (96) 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: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
1:A:296:VAL:HG12	1:A:300:LYS:HB3	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:285:LEU:HB3	1:A:346:MET:HE2	1.93	0.51
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:5:ASN:HD22	1:A:5:ASN:C	2.14	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:301:LEU:O	1:A:305:ILE:HG13	2.14	0.47
1:A:168:TRP:CB	1:A:414:ALA:HB2	2.44	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:135:LEU:HD23	1:A:140:GLN:HG2	2.00	0.43
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:327:ASP:O	1:A:330:TYR:HB3	2.20	0.41
1:A:324:ARG:HD3	1:A:324:ARG:HA	1.87	0.41
1:A:232:LEU:HA	1:A:400:SER:HB2	2.02	0.41
1:A:178:LEU:O	1:A:178:LEU:HD12	2.20	0.41
1:A:284:TRP:HZ3	1:A:353:LEU:HD21	1.86	0.41
1:A:193:LEU:HD12	1:A:193:LEU:HA	1.86	0.41
1:A:430:LEU:HA	1:A:430:LEU:HD23	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: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: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:346:MET:O	1:A:350:VAL:HG23	2.22	0.40
1:A:174:ASP:HB2	1:A:258:PRO:HG2	2.03	0.40
1:A:135:LEU:HD11	1:A:139:SER:HB3	2.02	0.40
1:A:93:PHE:O	1:A:97:VAL:HG23	2.22	0.40
1:A:95:LEU:HA	1:A:100:GLN:O	2.21	0.40
1:A:21:ILE:HG21	1:A:21:ILE:HD13	1.86	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
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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%)	26 19

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%)	6 3

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

Mol	Chain	Res	Type
1	A	5	ASN

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Mol	Chain	Res	Type
1	A	13	LEU
1	A	20	ARG
1	A	27	GLN
1	A	30	ASN
1	A	46	ARG
1	A	54	GLU
1	A	62	GLU
1	A	80	LYS
1	A	87	PRO
1	A	88	LEU
1	A	105	GLU
1	A	106	GLN
1	A	133	THR
1	A	149	ASN
1	A	165	THR
1	A	171	ILE
1	A	178	LEU
1	A	191	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	328	PRO
1	A	339	LYS
1	A	341	LEU
1	A	343	HIS
1	A	346	MET
1	A	362	LEU
1	A	383	GLN
1	A	387	MET
1	A	391	ASN
1	A	397	PHE
1	A	405	VAL
1	A	413	ARG

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

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	5	ASN
1	A	27	GLN
1	A	28	HIS
1	A	123	HIS
1	A	136	HIS
1	A	140	GLN
1	A	149	ASN
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	340	HIS
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 carbohydrates in this entry.

5.6 Ligand geometry (i)

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	-	40,50,50	1.50	5 (12%)	50,75,75	2.06	14 (28%)
3	CIT	A	439	-	3,12,12	1.04	0	3,17,17	1.50	1 (33%)

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	1/44/64/64	0/3/3/3
3	CIT	A	439	-	-	0/6/16/16	0/0/0/0

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	438	COA	C4A-N3A	-5.07	1.28	1.35
2	A	438	COA	C5A-N7A	-3.54	1.27	1.39
2	A	438	COA	C6A-N1A	-2.26	1.26	1.37
2	A	438	COA	C8A-N7A	2.35	1.39	1.34
2	A	438	COA	O9P-C9P	2.41	1.28	1.23

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	438	COA	C1B-N9A-C4A	-5.70	118.35	126.94
2	A	438	COA	P2A-O3A-P1A	-3.94	121.67	132.73
2	A	438	COA	O9P-C9P-N8P	-3.92	115.23	123.08
2	A	438	COA	C5B-C4B-C3B	-2.57	105.05	114.31
2	A	438	COA	CDP-CBP-CAP	-2.55	104.69	109.34
3	A	439	CIT	C3-C4-C5	-2.55	110.89	114.96

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	438	COA	CEP-CBP-CAP	2.03	113.05	109.34
2	A	438	COA	O4B-C1B-N9A	2.04	112.37	108.10
2	A	438	COA	O5P-C5P-N4P	2.06	127.03	122.94
2	A	438	COA	C2A-N1A-C6A	2.09	122.50	118.77
2	A	438	COA	N6A-C6A-N1A	2.24	124.00	119.20
2	A	438	COA	O3A-P1A-O5B	2.62	109.89	102.94
2	A	438	COA	OAP-CAP-C9P	3.68	118.83	110.38
2	A	438	COA	C3B-C2B-C1B	4.07	109.74	99.98
2	A	438	COA	C4B-O4B-C1B	5.61	115.88	109.72

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	438	COA	CAP

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	438	COA	CAP-C9P-N8P-C7P

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

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

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates

EDS was not executed - this section will therefore be empty.

6.4 Ligands

EDS was not executed - this section will therefore be empty.

6.5 Other polymers

EDS was not executed - this section will therefore be empty.