



Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 08:38 PM GMT

PDB ID : 2IFC
Title : The Structure of the Binary Complex of Oxalateacetate with Citrate Synthase
from the Thermophilic Archaeon Thermolasma acidophilum
Authors : Lehmann, C.
Deposited on : 2006-09-20
Resolution : 1.70 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>

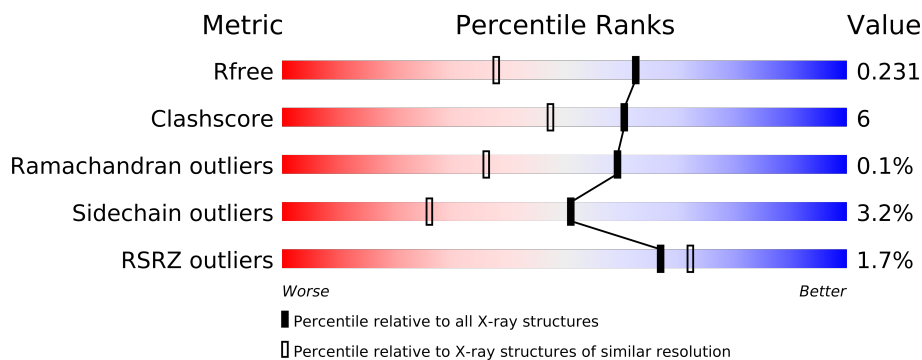
The following versions of software and data (see [references](#)) were used in the production of this report:

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 1.70 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	2456 (1.70-1.70)
Clashscore	79885	2929 (1.70-1.70)
Ramachandran outliers	78287	2878 (1.70-1.70)
Sidechain outliers	78261	2878 (1.70-1.70)
RSRZ outliers	66119	2456 (1.70-1.70)

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	385	
1	B	385	
1	C	385	
1	D	385	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 14321 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.

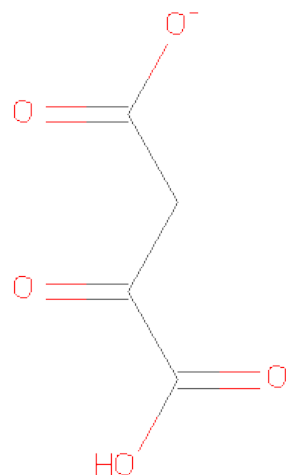
- Molecule 1 is a protein called Citrate Synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	380	Total	C	N	O	S	0	3	0
			3019	1932	511	563	13			
1	B	380	Total	C	N	O	S	0	2	0
			3012	1926	510	564	12			
1	C	382	Total	C	N	O	S	0	6	0
			3057	1950	522	573	12			
1	D	376	Total	C	N	O	S	0	1	0
			2969	1901	501	555	12			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MET	-	INITIATING METHIONINE	UNP P21553
B	0	MET	-	INITIATING METHIONINE	UNP P21553
C	0	MET	-	INITIATING METHIONINE	UNP P21553
D	0	MET	-	INITIATING METHIONINE	UNP P21553

- Molecule 2 is OXALOACETATE ION (three-letter code: OAA) (formula: C₄H₃O₅).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	C	1	Total	C	O	0	0
			9	4	5		
2	B	1	Total	C	O	0	0
			9	4	5		
2	A	1	Total	C	O	0	0
			9	4	5		
2	D	1	Total	C	O	0	0
			9	4	5		

- Molecule 3 is water.

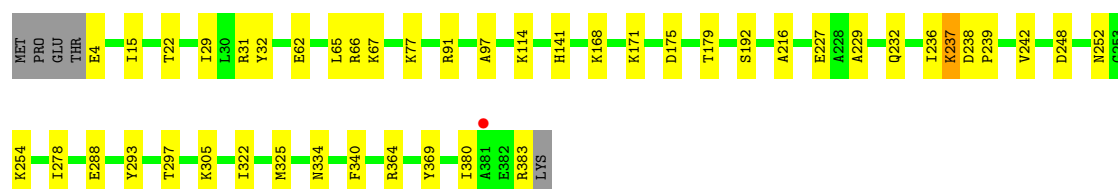
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	670	Total	O	0	0
			670	670		
3	B	472	Total	O	0	0
			472	472		
3	C	628	Total	O	0	0
			628	628		
3	D	458	Total	O	0	0
			458	458		

3 Residue-property plots

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.

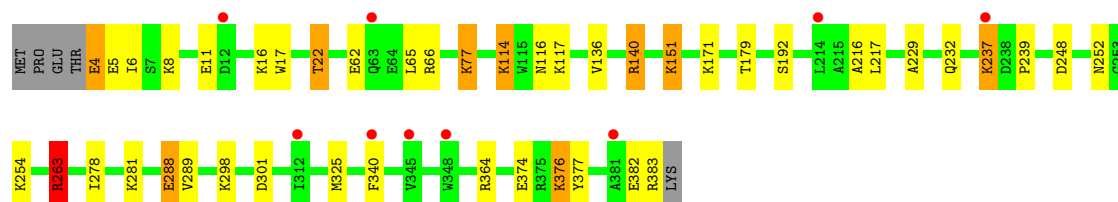
• Molecule 1: Citrate Synthase

Chain A:



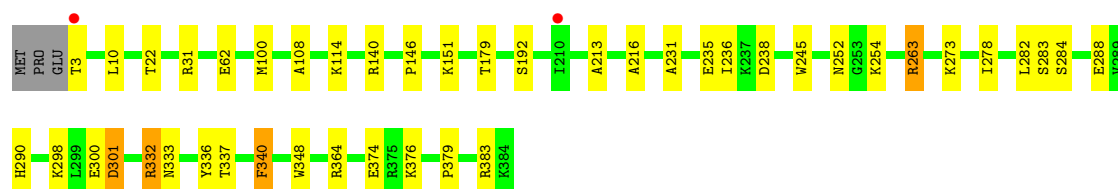
• Molecule 1: Citrate Synthase

Chain B:



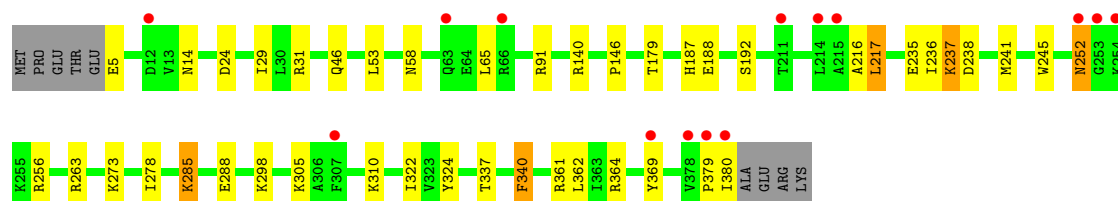
• Molecule 1: Citrate Synthase

Chain C:



• Molecule 1: Citrate Synthase

Chain D:



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	78.53Å 97.45Å 106.94Å 90.00° 93.10° 90.00°	Depositor
Resolution (Å)	41.38 – 1.70 41.38 – 1.70	Depositor EDS
% Data completeness (in resolution range)	96.8 (41.38-1.70) 96.8 (41.38-1.70)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.13 (at 1.70Å)	Xtriage
Refinement program	REFMAC 5.2	Depositor
R, R_{free}	0.181 , 0.234 0.178 , 0.231	Depositor DCC
R_{free} test set	8560 reflections (5.28%)	DCC
Wilson B-factor (Å ²)	23.4	Xtriage
Anisotropy	0.109	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 48.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.35$	Xtriage
Outliers	0 of 170708 reflections	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	14321	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.89% 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: OAA

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	1.04	1/3083 (0.0%)	0.90	3/4165 (0.1%)
1	B	0.76	0/3076	0.76	6/4156 (0.1%)
1	C	1.09	6/3133 (0.2%)	0.94	6/4232 (0.1%)
1	D	0.85	1/3033 (0.0%)	0.75	1/4100 (0.0%)
All	All	0.95	8/12325 (0.1%)	0.84	16/16653 (0.1%)

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	D	0	1

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	332[A]	ARG	C-N	-7.04	1.17	1.34
1	C	332[B]	ARG	C-N	-7.04	1.17	1.34
1	C	301[A]	ASP	C-N	-6.89	1.18	1.34
1	C	301[B]	ASP	C-N	-6.89	1.18	1.34
1	C	336	TYR	CD2-CE2	5.64	1.47	1.39
1	D	324	TYR	CD2-CE2	5.62	1.47	1.39
1	A	32	TYR	CD2-CE2	5.57	1.47	1.39
1	C	300	GLU	C-N	-5.14	1.22	1.34

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	383	ARG	NE-CZ-NH1	-10.49	115.06	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	332[A]	ARG	NE-CZ-NH1	9.23	124.91	120.30
1	C	332[B]	ARG	NE-CZ-NH1	9.23	124.91	120.30
1	A	383	ARG	NE-CZ-NH2	8.12	124.36	120.30
1	B	263	ARG	NE-CZ-NH2	-7.87	116.37	120.30
1	C	332[A]	ARG	NE-CZ-NH2	-7.84	116.38	120.30
1	C	332[B]	ARG	NE-CZ-NH2	-7.84	116.38	120.30
1	B	383	ARG	NE-CZ-NH2	-7.69	116.46	120.30
1	B	383	ARG	NE-CZ-NH1	6.77	123.69	120.30
1	B	140	ARG	NE-CZ-NH2	-6.24	117.18	120.30
1	C	336	TYR	CB-CG-CD1	-5.67	117.60	121.00
1	B	140	ARG	NE-CZ-NH1	5.51	123.06	120.30
1	C	383	ARG	NE-CZ-NH1	-5.46	117.57	120.30
1	D	217	LEU	CB-CG-CD2	5.31	120.02	111.00
1	A	248	ASP	CB-CG-OD2	5.10	122.89	118.30
1	B	263	ARG	NE-CZ-NH1	5.04	122.82	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	379	PRO	Peptide

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	3019	0	3026	37	0
1	B	3012	0	3013	41	0
1	C	3057	0	3047	41	1
1	D	2969	0	2971	37	0
2	A	9	0	2	0	0
2	B	9	0	2	0	0
2	C	9	0	2	0	0
2	D	9	0	2	0	0
3	A	670	0	0	23	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	472	0	0	13	1
3	C	628	0	0	13	0
3	D	458	0	0	13	0
All	All	14321	0	12065	142	1

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 6.

All (142) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:4:GLU:HG3	1:B:5:GLU:N	1.76	0.99
1:A:237[B]:LYS:HA	1:A:237[B]:LYS:HE3	1.45	0.94
1:A:288:GLU:HG3	3:A:822:HOH:O	1.69	0.91
1:B:252:ASN:OD1	3:B:855:HOH:O	1.90	0.89
1:B:151:LYS:HG2	3:B:821:HOH:O	1.74	0.87
1:D:46:GLN:HG2	3:D:639:HOH:O	1.74	0.87
1:A:77:LYS:HD3	3:A:774:HOH:O	1.77	0.85
1:B:4:GLU:HG3	1:B:5:GLU:H	1.36	0.84
1:C:263[B]:ARG:HH11	1:C:263[B]:ARG:HG3	1.44	0.83
1:C:332[B]:ARG:NH1	3:C:847:HOH:O	2.13	0.82
1:A:305:LYS:HD2	3:A:872:HOH:O	1.80	0.80
1:D:256:ARG:HD2	3:D:818:HOH:O	1.82	0.79
1:C:298:LYS:HG3	3:C:625:HOH:O	1.85	0.77
1:C:273:LYS:HE2	3:C:970:HOH:O	1.86	0.76
1:B:377:TYR:OH	3:B:806:HOH:O	1.98	0.75
1:C:179:THR:HG21	1:C:278:ILE:HD11	1.69	0.75
1:B:288:GLU:CD	1:B:288:GLU:H	1.93	0.72
1:B:77:LYS:HD2	3:B:623:HOH:O	1.89	0.72
1:B:22:THR:HG23	1:D:369:TYR:CG	2.25	0.72
1:A:29:ILE:HD12	3:A:930:HOH:O	1.89	0.71
1:C:140:ARG:HD3	1:C:146:PRO:O	1.91	0.71
1:C:332[B]:ARG:CZ	3:C:847:HOH:O	2.39	0.70
1:D:288:GLU:HG2	3:D:673:HOH:O	1.91	0.70
1:D:238:ASP:HB3	1:D:241:MET:HE2	1.74	0.69
1:A:66:ARG:HD3	3:A:768:HOH:O	1.93	0.69
1:B:252:ASN:HB2	3:B:815:HOH:O	1.93	0.69
1:C:179:THR:HG21	1:C:278:ILE:CD1	2.22	0.68
1:C:238:ASP:HB3	3:C:978:HOH:O	1.94	0.68
1:A:171:LYS:HD2	3:A:761:HOH:O	1.93	0.68
1:B:289:VAL:HG13	1:B:325:MET:HE3	1.76	0.67
1:C:62[B]:GLU:OE2	3:C:601:HOH:O	2.12	0.67
1:D:91:ARG:HB2	1:D:91:ARG:NH1	2.10	0.66

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:29:ILE:HG22	3:D:641:HOH:O	1.96	0.66
1:A:239:PRO:HG3	3:A:672:HOH:O	1.95	0.66
1:D:238:ASP:HB2	1:D:241:MET:HE3	1.77	0.66
1:C:288[A]:GLU:CD	1:C:288[A]:GLU:H	1.99	0.65
1:B:136:VAL:O	1:B:140:ARG:HG2	1.97	0.65
1:A:175:ASP:HB2	3:A:767:HOH:O	1.97	0.64
1:B:289:VAL:HG13	1:B:325:MET:CE	2.27	0.64
3:A:1017:HOH:O	1:C:374:GLU:HG3	1.98	0.64
1:C:301[B]:ASP:OD1	3:C:499:HOH:O	2.14	0.63
1:D:238:ASP:CB	1:D:241:MET:HE3	2.29	0.63
1:D:238:ASP:HB3	1:D:241:MET:CE	2.27	0.63
1:D:91:ARG:HB2	1:D:91:ARG:HH11	1.63	0.63
1:C:3:THR:N	3:C:972:HOH:O	2.31	0.63
1:B:22:THR:HG23	1:D:369:TYR:CD1	2.35	0.61
1:B:254:LYS:HG3	3:B:616:HOH:O	2.01	0.61
1:A:62:GLU:HG2	3:A:772:HOH:O	2.01	0.61
1:D:238:ASP:CB	1:D:241:MET:CE	2.80	0.60
1:A:22:THR:HG23	3:A:763:HOH:O	2.01	0.59
1:D:322:ILE:HG13	3:D:627:HOH:O	2.02	0.59
1:B:62:GLU:O	1:B:66:ARG:HG3	2.03	0.59
1:C:263[A]:ARG:HD2	3:C:824:HOH:O	2.02	0.58
1:D:252:ASN:O	3:D:799:HOH:O	2.17	0.58
1:A:15[B]:ILE:HG22	1:C:10:LEU:HD21	1.85	0.57
1:C:263[B]:ARG:HG3	1:C:263[B]:ARG:NH1	2.16	0.57
1:C:114:LYS:HE3	3:C:666:HOH:O	2.05	0.56
1:B:252:ASN:HB2	3:B:616:HOH:O	2.06	0.55
1:D:241:MET:HG2	3:D:726:HOH:O	2.06	0.55
1:A:369:TYR:CG	1:C:22:THR:CG2	2.90	0.55
3:A:764:HOH:O	1:C:376:LYS:HG2	2.06	0.54
1:C:263[B]:ARG:HB2	3:C:990:HOH:O	2.08	0.54
1:B:140:ARG:NH2	3:B:718:HOH:O	2.26	0.54
1:B:281:LYS:HE2	3:B:820:HOH:O	2.07	0.54
1:B:374:GLU:OE2	1:D:31:ARG:NH1	2.41	0.53
1:B:263:ARG:NH2	1:D:362:LEU:O	2.41	0.53
1:B:376:LYS:HE3	3:D:643:HOH:O	2.08	0.53
3:A:765:HOH:O	1:C:100:MET:HE2	2.08	0.53
1:B:6:ILE:HG22	1:B:8:LYS:HG3	1.90	0.52
1:B:229:ALA:O	1:B:232:GLN:HG2	2.10	0.52
1:C:231:ALA:O	1:C:235:GLU:HG3	2.09	0.52
1:B:239:PRO:HG3	3:B:617:HOH:O	2.10	0.52
1:B:171:LYS:HE3	3:B:619:HOH:O	2.09	0.51
1:A:380:ILE:HG12	1:C:62[B]:GLU:HG2	1.93	0.51

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:192:SER:HA	1:D:216:ALA:HB3	1.92	0.51
1:B:114[A]:LYS:NZ	1:B:116:ASN:HD22	2.07	0.51
1:B:239:PRO:HB3	1:B:298:LYS:HE2	1.92	0.50
1:A:114:LYS:HD2	3:A:1035:HOH:O	2.11	0.50
1:A:322:ILE:HG13	3:A:555:HOH:O	2.11	0.50
1:B:237:LYS:HB3	3:B:645:HOH:O	2.12	0.49
1:A:369:TYR:CD1	1:C:22:THR:HG23	2.48	0.49
1:D:285:LYS:HB2	3:D:572:HOH:O	2.11	0.49
1:B:192:SER:HA	1:B:216:ALA:HB3	1.95	0.49
1:C:192:SER:HA	1:C:216:ALA:HB3	1.95	0.49
1:C:31:ARG:HD2	3:C:938:HOH:O	2.12	0.48
1:C:288[B]:GLU:OE2	3:C:729:HOH:O	2.20	0.48
1:A:380:ILE:HG12	1:C:62[A]:GLU:HG3	1.96	0.48
1:D:238:ASP:HB2	1:D:241:MET:CE	2.44	0.47
1:D:263:ARG:HB2	3:D:798:HOH:O	2.14	0.47
1:A:237[B]:LYS:HE3	1:A:237[B]:LYS:CA	2.29	0.47
1:B:22:THR:CG2	1:D:369:TYR:CG	2.96	0.47
1:D:179:THR:HG21	1:D:278:ILE:CD1	2.45	0.47
1:D:236:ILE:O	1:D:237:LYS:HB2	2.15	0.47
1:A:322:ILE:HA	1:A:325[B]:MET:HE3	1.96	0.47
1:B:248:ASP:O	3:B:815:HOH:O	2.21	0.46
1:A:192:SER:HA	1:A:216:ALA:HB3	1.98	0.46
1:A:227:GLU:HG2	1:A:334:ASN:HA	1.97	0.46
1:A:252:ASN:ND2	1:A:254:LYS:HG3	2.30	0.46
3:A:764:HOH:O	1:C:376:LYS:CG	2.64	0.46
1:A:229:ALA:O	1:A:232:GLN:HG2	2.16	0.46
1:B:179:THR:HG21	1:B:278:ILE:CD1	2.46	0.46
1:A:236:ILE:HG21	1:A:242:VAL:HG22	1.98	0.46
1:B:4:GLU:CG	1:B:5:GLU:H	2.19	0.45
1:A:175:ASP:CB	3:A:767:HOH:O	2.60	0.45
3:A:762:HOH:O	1:C:100:MET:HE2	2.16	0.45
1:C:236:ILE:HG12	1:C:245:TRP:CG	2.52	0.45
1:D:235[A]:GLU:CD	1:D:245:TRP:HE1	2.20	0.45
1:B:62:GLU:CD	1:B:66:ARG:HH11	2.20	0.45
1:A:254:LYS:HB3	1:A:254:LYS:HE3	1.62	0.45
1:A:91:ARG:NH1	1:A:141:HIS:O	2.44	0.45
3:A:1004:HOH:O	1:C:379:PRO:HD3	2.17	0.45
1:B:151:LYS:HB3	1:B:151:LYS:HE2	1.68	0.44
1:C:252:ASN:OD1	1:C:254:LYS:HG2	2.17	0.44
1:D:24:ASP:HB3	1:D:29:ILE:HB	1.99	0.44
1:A:179:THR:HG21	1:A:278:ILE:CD1	2.47	0.44
1:A:114:LYS:HE3	3:A:784:HOH:O	2.17	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:293:TYR:CZ	1:A:297:THR:HG21	2.52	0.44
1:B:239:PRO:HB3	1:B:298:LYS:CE	2.48	0.44
1:D:298:LYS:HD2	3:D:740:HOH:O	2.17	0.43
1:C:332[B]:ARG:NE	1:C:333:ASN:OD1	2.51	0.43
1:B:179:THR:HG21	1:B:278:ILE:HD11	2.01	0.43
1:D:140:ARG:HD3	1:D:146:PRO:O	2.19	0.43
1:B:11:GLU:OE1	1:D:361:ARG:NH2	2.47	0.43
1:A:77:LYS:CD	3:A:774:HOH:O	2.51	0.42
1:C:337:THR:O	1:C:340:PHE:HB3	2.19	0.42
1:A:97:ALA:HB1	1:C:108:ALA:HA	2.00	0.42
1:C:263[B]:ARG:CG	1:C:263[B]:ARG:NH1	2.82	0.42
1:B:16:LYS:HE2	1:B:17:TRP:O	2.19	0.42
1:D:235[A]:GLU:HG3	3:D:784:HOH:O	2.19	0.42
1:B:4:GLU:HG2	1:D:14:ASN:OD1	2.19	0.42
1:B:4:GLU:CG	1:B:5:GLU:N	2.63	0.42
1:D:310:LYS:HB3	1:D:310:LYS:HE2	1.87	0.42
1:C:283:SER:O	1:C:290[B]:HIS:HB2	2.20	0.41
1:A:67:LYS:HD3	3:A:918:HOH:O	2.20	0.41
1:A:171:LYS:HE3	3:A:767:HOH:O	2.20	0.41
1:D:337:THR:O	1:D:340:PHE:HB3	2.21	0.41
1:C:213:ALA:HB1	1:C:348:TRP:CE2	2.56	0.40
1:D:273:LYS:HB2	1:D:273:LYS:HE2	1.83	0.40
1:A:22:THR:HG22	1:A:31:ARG:HB2	2.04	0.40
1:A:369:TYR:CG	1:C:22:THR:HG23	2.56	0.40
1:D:238:ASP:CB	1:D:241:MET:HE2	2.47	0.40
1:D:305:LYS:NZ	3:D:652:HOH:O	2.51	0.40

All (1) 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)
1:C:332[B]:ARG:NH2	3:B:706:HOH:O[2.553]	2.12	0.08

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	381/385 (99%)	368 (97%)	13 (3%)	0	100	100
1	B	380/385 (99%)	372 (98%)	8 (2%)	0	100	100
1	C	386/385 (100%)	376 (97%)	10 (3%)	0	100	100
1	D	375/385 (97%)	364 (97%)	9 (2%)	2 (0%)	38	17
All	All	1522/1540 (99%)	1480 (97%)	40 (3%)	2 (0%)	59	36

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	252	ASN
1	D	188	GLU

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	313/315 (99%)	305 (97%)	8 (3%)	59	35
1	B	312/315 (99%)	295 (95%)	17 (5%)	31	11
1	C	318/315 (101%)	311 (98%)	7 (2%)	64	43
1	D	308/315 (98%)	297 (96%)	11 (4%)	47	22
All	All	1251/1260 (99%)	1208 (97%)	43 (3%)	51	23

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	GLU
1	A	65	LEU
1	A	168	LYS
1	A	237[A]	LYS
1	A	237[B]	LYS
1	A	238	ASP
1	A	340	PHE
1	A	364	ARG
1	B	4	GLU
1	B	22	THR

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Mol	Chain	Res	Type
1	B	65	LEU
1	B	77	LYS
1	B	114[A]	LYS
1	B	114[B]	LYS
1	B	117	LYS
1	B	151	LYS
1	B	217	LEU
1	B	237	LYS
1	B	263	ARG
1	B	288	GLU
1	B	301	ASP
1	B	340	PHE
1	B	364	ARG
1	B	376	LYS
1	B	382	GLU
1	C	151	LYS
1	C	263[A]	ARG
1	C	263[B]	ARG
1	C	282	LEU
1	C	284	SER
1	C	340	PHE
1	C	364	ARG
1	D	5	GLU
1	D	53	LEU
1	D	58	ASN
1	D	65	LEU
1	D	187	HIS
1	D	217	LEU
1	D	237	LYS
1	D	285	LYS
1	D	340	PHE
1	D	364	ARG
1	D	380	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	58	ASN
1	D	247	ASN

5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

4 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	OAA	A	385	-	6,8,8	4.24	3 (50%)	7,10,10	3.48	5 (71%)
2	OAA	B	385	-	6,8,8	2.43	3 (50%)	7,10,10	1.87	1 (14%)
2	OAA	C	385	-	6,8,8	2.18	2 (33%)	7,10,10	2.01	3 (42%)
2	OAA	D	385	-	6,8,8	3.15	3 (50%)	7,10,10	1.97	3 (42%)

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	OAA	A	385	-	-	0/6/8/8	0/0/0/0
2	OAA	B	385	-	-	0/6/8/8	0/0/0/0
2	OAA	C	385	-	-	0/6/8/8	0/0/0/0
2	OAA	D	385	-	-	0/6/8/8	0/0/0/0

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	385	OAA	C2-C1	7.20	1.54	1.49
2	A	385	OAA	O3-C3	5.82	1.36	1.23
2	D	385	OAA	C2-C1	4.96	1.52	1.49
2	D	385	OAA	O3-C3	4.27	1.33	1.23
2	C	385	OAA	O3-C3	4.27	1.33	1.23
2	A	385	OAA	C2-C3	4.04	1.55	1.51
2	B	385	OAA	O3-C3	3.89	1.32	1.23
2	D	385	OAA	O4-C4	3.46	1.33	1.22
2	B	385	OAA	O4-C4	3.45	1.33	1.22
2	C	385	OAA	O4-C4	3.04	1.31	1.22
2	B	385	OAA	C2-C1	2.23	1.51	1.49

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	385	OAA	C1-C2-C3	-4.90	105.11	114.01
2	A	385	OAA	O3-C3-C2	4.74	131.36	120.84
2	B	385	OAA	C1-C2-C3	-4.06	106.63	114.01
2	A	385	OAA	C2-C3-C4	-3.81	111.71	117.75
2	D	385	OAA	O3-C3-C2	3.74	129.15	120.84
2	A	385	OAA	O4-C4-C3	-3.47	111.58	120.27
2	C	385	OAA	C1-C2-C3	-3.26	108.08	114.01
2	C	385	OAA	O3-C3-C2	2.90	127.28	120.84
2	A	385	OAA	O5-C4-C3	2.87	120.70	114.37
2	C	385	OAA	C2-C3-C4	-2.55	113.71	117.75
2	D	385	OAA	C2-C3-C4	-2.42	113.91	117.75
2	D	385	OAA	C1-C2-C3	-2.10	110.19	114.01

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	380/385 (98%)	-0.12	1 (0%) 91 95	11, 18, 31, 44	0
1	B	380/385 (98%)	0.15	9 (2%) 56 61	18, 28, 39, 47	0
1	C	382/385 (99%)	-0.14	2 (0%) 88 92	11, 18, 30, 45	0
1	D	376/385 (97%)	0.20	14 (3%) 39 44	17, 25, 40, 53	0
All	All	1518/1540 (98%)	0.02	26 (1%) 67 72	11, 23, 37, 53	0

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	253	GLY	4.4
1	C	3	THR	3.7
1	D	378	VAL	3.4
1	A	381	ALA	3.3
1	D	252	ASN	3.3
1	D	379	PRO	2.9
1	D	380	ILE	2.8
1	D	214	LEU	2.7
1	B	345	VAL	2.4
1	D	66	ARG	2.3
1	D	307	PHE	2.3
1	B	237	LYS	2.2
1	D	215	ALA	2.2
1	B	12	ASP	2.2
1	D	254	LYS	2.2
1	D	211	THR	2.1
1	B	63	GLN	2.1
1	B	214	LEU	2.1
1	D	369	TYR	2.1
1	C	210	ILE	2.1
1	D	63	GLN	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	381	ALA	2.0
1	B	340	PHE	2.0
1	B	312	ILE	2.0
1	B	348	TRP	2.0
1	D	12	ASP	2.0

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

There are no non-standard protein/DNA/RNA residues in this entry.

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(Å ²)	Q<0.9
2	OAA	D	385	9/9	0.13	0.21	20,25,28,30	0
2	OAA	A	385	9/9	0.10	-0.12	14,16,20,21	0
2	OAA	C	385	9/9	0.09	-0.51	18,19,21,23	0
2	OAA	B	385	9/9	0.07	-0.65	25,27,31,32	0

6.5 Other polymers ⓘ

There are no such residues in this entry.