



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 05:31 AM GMT

PDB ID : 2R26  
Title : The Structure of the Ternary Complex of Carboxymethyl Coenzyme A and Oxalateacetate with Citrate Synthase from the Thermophilic Archaeonthermoplasma Acidophilum  
Authors : Lehmann, C.  
Deposited on : 2007-08-24  
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](mailto:validation@mail.wwpdb.org)  
A user guide is available at  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

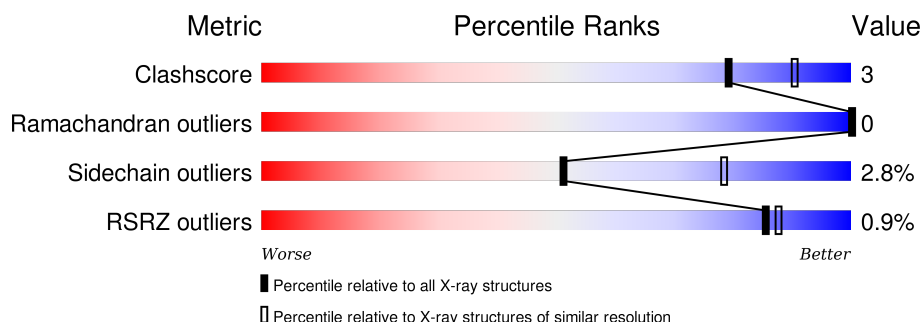
# 1 Overall quality at a glance

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	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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  $\leq 5\%$ . 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.

Mol	Chain	Length	Quality of chain
1	A	384	 89% 10% .
1	B	384	 2% 89% 9% ..
1	C	384	 1% 89% 9% ..
1	D	384	 1% 88% 11% .

## 2 Entry composition [i](#)

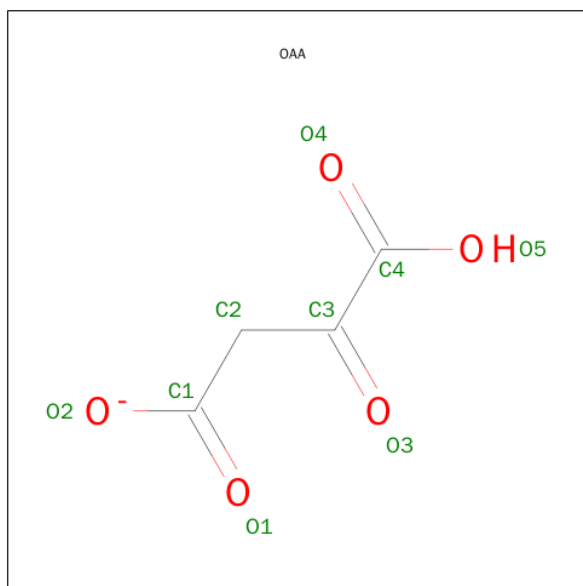
There are 4 unique types of molecules in this entry. The entry contains 13168 atoms, of which 0 are hydrogens and 0 are deuteriums.

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
1	A	381	Total	C	N	O	S	0	0	0
			3001	1919	508	562	12			
1	B	380	Total	C	N	O	S	0	0	0
			2994	1915	507	560	12			
1	C	379	Total	C	N	O	S	0	0	0
			2985	1910	506	557	12			
1	D	379	Total	C	N	O	S	0	0	0
			2985	1910	506	557	12			

- Molecule 2 is OXALOACETATE ION (three-letter code: OAA) (formula:  $C_4H_3O_5$ ).



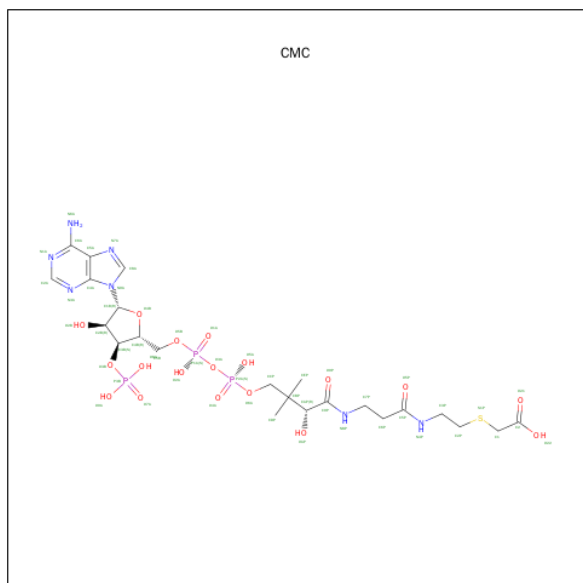
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			9	4	5		
2	B	1	Total	C	O	0	0
			9	4	5		

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

- Molecule 3 is CARBOXYMETHYL COENZYME \*A (three-letter code: CMC) (formula:  $C_{23}H_{38}N_7O_{18}P_3S$ ).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	S	0	0
			52	23	7	18	3	1		
3	B	1	Total	C	N	O	P	S	0	0
			52	23	7	18	3	1		
3	C	1	Total	C	N	O	P	S	0	0
			52	23	7	18	3	1		
3	D	1	Total	C	N	O	P	S	0	0
			52	23	7	18	3	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	255	Total	O	0	0
			255	255		
4	B	230	Total	O	0	0
			230	230		
4	C	228	Total	O	0	0
			228	228		

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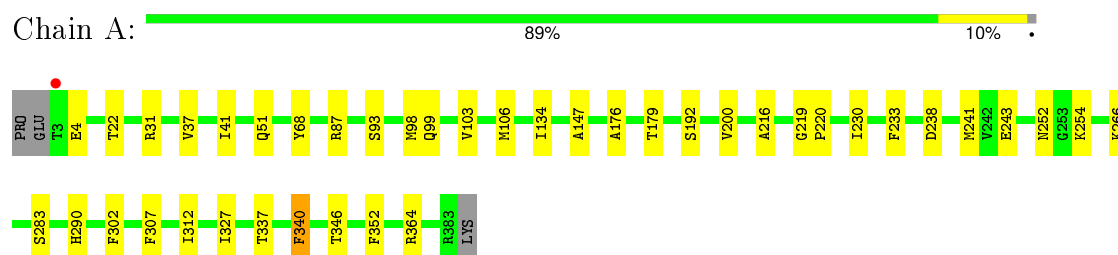
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	246	Total	O	0	0
			246	246		

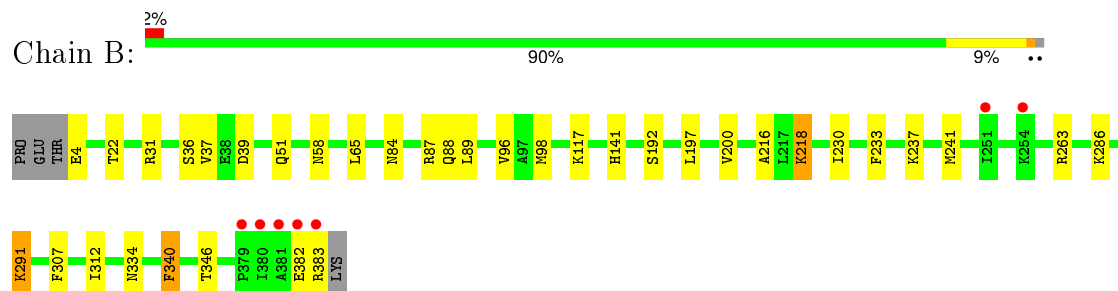
### 3 Residue-property plots

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.

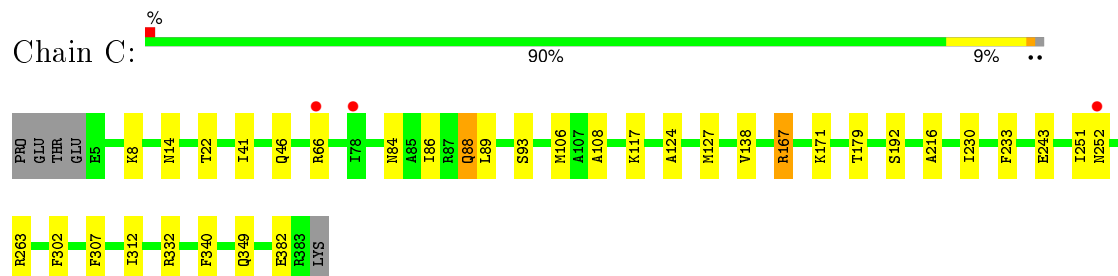
#### ● Molecule 1: Citrate Synthase



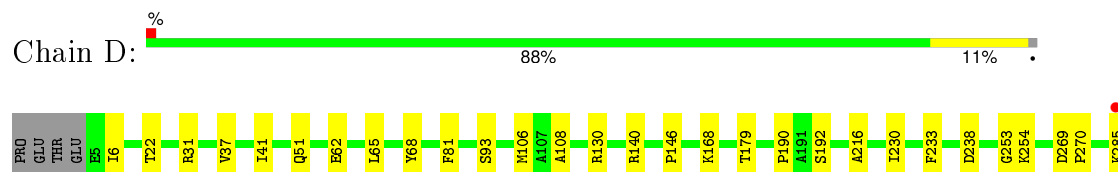
#### ● Molecule 1: Citrate Synthase

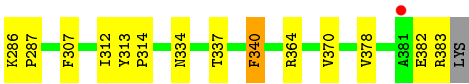


#### ● Molecule 1: Citrate Synthase



#### ● Molecule 1: Citrate Synthase





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	67.70Å 74.22Å 89.87Å 99.34° 98.34° 114.38°	Depositor
Resolution (Å)	43.09 – 2.50 43.09 – 2.40	Depositor EDS
% Data completeness (in resolution range)	97.5 (43.09-2.50) 85.5 (43.09-2.40)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.13 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.3	Depositor
R, $R_{free}$	0.200 , 0.261 0.199 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	DCC
Wilson B-factor (Å <sup>2</sup> )	30.2	Xtriage
Anisotropy	0.514	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.29 , 41.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.52$ , $\langle L^2 \rangle = 0.35$	Xtriage
Outliers	0 of 58212 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	13168	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.99% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.



## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: CMC, 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	0.34	0/3065	0.46	0/4143
1	B	0.34	0/3058	0.49	0/4133
1	C	0.34	0/3049	0.49	0/4121
1	D	0.34	0/3049	0.49	0/4121
All	All	0.34	0/12221	0.48	0/16518

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 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	3001	0	3003	25	0
1	B	2994	0	2996	19	0
1	C	2985	0	2990	20	0
1	D	2985	0	2990	26	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	1	0
3	A	52	0	31	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	52	0	31	0	0
3	C	52	0	31	0	0
3	D	52	0	31	0	0
4	A	255	0	0	0	0
4	B	230	0	0	1	0
4	C	228	0	0	2	0
4	D	246	0	0	0	0
All	All	13168	0	12111	84	0

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 3.

All (84) 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:D:382:GLU:HA	1:D:383:ARG:C	2.00	0.81
1:C:167:ARG:HG2	1:C:167:ARG:HH21	1.46	0.80
1:B:334:ASN:HB3	4:B:925:HOH:O	1.84	0.78
1:C:307:PHE:HB3	1:C:312:ILE:HB	1.71	0.70
1:C:167:ARG:HH21	1:C:167:ARG:CG	2.07	0.66
1:A:238:ASP:HB3	1:A:241:MET:HG2	1.78	0.66
1:A:176:ALA:HB3	1:A:327:ILE:HD11	1.77	0.65
1:A:307:PHE:HB3	1:A:312:ILE:HB	1.80	0.64
1:D:41:ILE:HD12	1:D:179:THR:HG23	1.80	0.63
1:C:106:MET:HE1	1:C:127:MET:SD	2.41	0.61
1:C:108:ALA:HB1	1:D:93:SER:HB3	1.84	0.59
1:D:22:THR:HG22	1:D:31:ARG:HB2	1.87	0.57
1:D:106:MET:CE	1:D:130:ARG:HD3	2.36	0.56
1:A:37:VAL:O	1:A:41:ILE:HG12	2.08	0.53
1:C:230:ILE:HA	1:C:233:PHE:CD2	2.44	0.52
1:D:37:VAL:O	1:D:41:ILE:HG12	2.10	0.52
1:B:307:PHE:HB3	1:B:312:ILE:HB	1.93	0.51
1:A:252:ASN:HD21	1:A:254:LYS:HE2	1.75	0.51
1:A:230:ILE:HA	1:A:233:PHE:CD2	2.44	0.51
1:A:243:GLU:HG2	1:A:302:PHE:CZ	2.46	0.51
1:B:382:GLU:H	1:B:383:ARG:HB2	1.76	0.50
1:D:382:GLU:CA	1:D:383:ARG:C	2.75	0.50
1:A:106:MET:CE	1:A:134:ILE:HD11	2.41	0.50
1:C:192:SER:HA	1:C:216:ALA:HB3	1.93	0.49
1:B:230:ILE:HA	1:B:233:PHE:CD2	2.47	0.49
1:C:84:ASN:O	1:C:88:GLN:HB3	2.12	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:88:GLN:HG3	1:D:81:PHE:HB2	1.94	0.48
1:A:238:ASP:HB3	1:A:241:MET:CG	2.43	0.48
1:B:237:LYS:HE2	1:B:237:LYS:HA	1.96	0.48
1:A:364:ARG:HB3	1:B:263:ARG:HG2	1.96	0.48
1:B:192:SER:HA	1:B:216:ALA:HB3	1.96	0.48
1:C:124:ALA:HA	1:C:349:GLN:OE1	2.14	0.48
1:D:190:PRO:HB3	2:D:503:OAA:C1	2.44	0.47
1:C:263:ARG:HG2	1:D:364:ARG:HB3	1.96	0.47
1:B:96:VAL:HB	1:B:218:LYS:HG3	1.98	0.46
1:D:140:ARG:HD3	1:D:146:PRO:O	2.14	0.46
1:D:192:SER:HA	1:D:216:ALA:HB3	1.96	0.46
1:D:230:ILE:HA	1:D:233:PHE:CD1	2.49	0.46
1:C:106:MET:CE	1:C:127:MET:SD	3.03	0.46
1:D:51:GLN:HB3	1:D:68:TYR:OH	2.16	0.46
1:A:93:SER:HB2	1:A:98:MET:HE3	1.97	0.45
1:C:167:ARG:NH2	1:C:167:ARG:CG	2.72	0.45
1:A:192:SER:HA	1:A:216:ALA:HB3	1.99	0.45
1:B:89:LEU:O	1:B:141:HIS:HE1	2.00	0.45
1:C:41:ILE:HD12	1:C:179:THR:HG23	1.98	0.44
1:A:200:VAL:HG11	1:A:352:PHE:HB3	2.00	0.44
1:D:106:MET:HE1	1:D:130:ARG:HH21	1.83	0.44
1:A:41:ILE:HD12	1:A:179:THR:HG23	1.98	0.44
1:C:93:SER:HB3	1:D:108:ALA:HB1	1.98	0.44
1:A:337:THR:O	1:A:340:PHE:HB3	2.17	0.44
1:A:93:SER:HB2	1:A:98:MET:CE	2.48	0.44
1:B:84:ASN:O	1:B:88:GLN:HG2	2.18	0.43
1:A:106:MET:HE3	1:A:134:ILE:HD11	1.99	0.43
1:D:307:PHE:HB3	1:D:312:ILE:HB	2.00	0.43
1:B:291:LYS:NZ	1:B:291:LYS:HB2	2.34	0.43
1:B:340:PHE:C	1:B:340:PHE:CD2	2.92	0.43
1:B:197:LEU:HA	1:B:200:VAL:HG22	2.00	0.43
1:D:253:GLY:HA3	1:D:254:LYS:HA	1.90	0.43
1:C:14:ASN:HB2	1:D:6:ILE:HA	2.01	0.42
1:D:106:MET:HE2	1:D:106:MET:HB3	1.74	0.42
1:C:46:GLN:HG3	4:C:875:HOH:O	2.18	0.42
1:B:31:ARG:HG2	1:B:36:SER:HA	2.01	0.42
1:D:337:THR:O	1:D:340:PHE:HB3	2.19	0.42
1:A:219:GLY:HA2	1:A:220:PRO:HD3	1.90	0.42
1:A:87:ARG:HD2	1:A:147:ALA:HB2	2.01	0.42
1:B:89:LEU:O	1:B:141:HIS:CE1	2.72	0.42
1:B:340:PHE:C	1:B:340:PHE:HD2	2.23	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:286:LYS:HA	1:D:287:PRO:HD2	1.93	0.41
1:D:269:ASP:HA	1:D:270:PRO:HD2	1.90	0.41
1:A:51:GLN:HA	1:A:346:THR:HG21	2.01	0.41
1:B:89:LEU:HD12	1:B:98:MET:HG2	2.03	0.41
1:C:86:ILE:HG21	1:C:138:VAL:HA	2.03	0.41
1:D:340:PHE:C	1:D:340:PHE:CD2	2.94	0.41
1:A:51:GLN:HB3	1:A:68:TYR:OH	2.20	0.41
1:A:99:GLN:O	1:A:103:VAL:HG23	2.20	0.41
1:C:243:GLU:HG2	1:C:302:PHE:CZ	2.55	0.41
1:A:22:THR:HG22	1:A:31:ARG:HB2	2.02	0.41
1:B:51:GLN:HA	1:B:346:THR:HG21	2.03	0.41
1:B:87:ARG:HA	1:B:141:HIS:CD2	2.55	0.41
1:A:340:PHE:C	1:A:340:PHE:CD2	2.94	0.41
1:D:370:VAL:HG12	1:D:370:VAL:O	2.20	0.41
1:A:283:SER:O	1:A:290:HIS:HB2	2.21	0.40
1:C:117:LYS:HE2	4:C:867:HOH:O	2.21	0.40
1:D:313:TYR:HB3	1:D:314:PRO:HD2	2.04	0.40

There are no symmetry-related clashes.

## 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	379/384 (99%)	369 (97%)	10 (3%)	0	100	100
1	B	378/384 (98%)	367 (97%)	11 (3%)	0	100	100
1	C	377/384 (98%)	364 (97%)	13 (3%)	0	100	100
1	D	377/384 (98%)	363 (96%)	14 (4%)	0	100	100
All	All	1511/1536 (98%)	1463 (97%)	48 (3%)	0	100	100

There are no Ramachandran outliers to report.

### 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	311/314 (99%)	308 (99%)	3 (1%)	82	95
1	B	310/314 (99%)	298 (96%)	12 (4%)	39	66
1	C	309/314 (98%)	297 (96%)	12 (4%)	39	66
1	D	309/314 (98%)	301 (97%)	8 (3%)	54	81
All	All	1239/1256 (99%)	1204 (97%)	35 (3%)	51	78

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	GLU
1	A	266	LYS
1	A	340	PHE
1	B	4	GLU
1	B	22	THR
1	B	37	VAL
1	B	39	ASP
1	B	58	ASN
1	B	65	LEU
1	B	117	LYS
1	B	218	LYS
1	B	241	MET
1	B	286	LYS
1	B	291	LYS
1	B	340	PHE
1	C	8	LYS
1	C	22	THR
1	C	66	ARG
1	C	88	GLN
1	C	89	LEU
1	C	167	ARG
1	C	171	LYS
1	C	251	ILE
1	C	252	ASN

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Mol	Chain	Res	Type
1	C	332	ARG
1	C	340	PHE
1	C	382	GLU
1	D	62	GLU
1	D	65	LEU
1	D	168	LYS
1	D	238	ASP
1	D	285	LYS
1	D	334	ASN
1	D	340	PHE
1	D	378	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	116	ASN
1	B	141	HIS
1	C	14	ASN
1	D	252	ASN
1	D	334	ASN

### 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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

8 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	501	-	2,8,8	3.67	1 (50%)	2,10,10	1.70	1 (50%)
3	CMC	A	701	-	41,54,54	2.85	5 (12%)	53,80,80	1.84	5 (9%)
2	OAA	B	500	-	2,8,8	3.64	1 (50%)	2,10,10	1.24	0
3	CMC	B	700	-	41,54,54	2.85	5 (12%)	53,80,80	1.90	5 (9%)
2	OAA	C	502	-	2,8,8	3.60	1 (50%)	2,10,10	1.91	1 (50%)
3	CMC	C	702	-	41,54,54	2.84	4 (9%)	53,80,80	1.92	7 (13%)
2	OAA	D	503	-	2,8,8	3.71	1 (50%)	2,10,10	1.11	0
3	CMC	D	703	-	41,54,54	2.87	5 (12%)	53,80,80	1.93	6 (11%)

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	501	-	-	0/2/8/8	0/0/0/0
3	CMC	A	701	-	-	0/46/68/68	0/3/3/3
2	OAA	B	500	-	-	0/2/8/8	0/0/0/0
3	CMC	B	700	-	-	0/46/68/68	0/3/3/3
2	OAA	C	502	-	-	0/2/8/8	0/0/0/0
3	CMC	C	702	-	-	0/46/68/68	0/3/3/3
2	OAA	D	503	-	-	0/2/8/8	0/0/0/0
3	CMC	D	703	-	-	0/46/68/68	0/3/3/3

All (23) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	703	CMC	C1-S1P	-16.18	1.55	1.82
3	B	700	CMC	C1-S1P	-16.11	1.55	1.82
3	C	702	CMC	C1-S1P	-16.06	1.56	1.82
3	A	701	CMC	C1-S1P	-16.02	1.56	1.82
3	A	701	CMC	CEP-CBP	-4.93	1.43	1.53
3	B	700	CMC	CDP-CBP	-4.92	1.43	1.53
3	C	702	CMC	CEP-CBP	-4.88	1.43	1.53
3	D	703	CMC	CEP-CBP	-4.86	1.43	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	701	CMC	CDP-CBP	-4.76	1.43	1.53
3	D	703	CMC	CDP-CBP	-4.74	1.43	1.53
3	C	702	CMC	CDP-CBP	-4.63	1.44	1.53
3	B	700	CMC	CEP-CBP	-4.62	1.44	1.53
3	B	700	CMC	O4B-C1B	2.01	1.43	1.41
3	A	701	CMC	O4B-C1B	2.10	1.43	1.41
3	D	703	CMC	O4B-C1B	2.23	1.44	1.41
3	C	702	CMC	C5A-C4A	3.12	1.47	1.40
3	A	701	CMC	C5A-C4A	3.18	1.47	1.40
3	B	700	CMC	C5A-C4A	3.19	1.47	1.40
3	D	703	CMC	C5A-C4A	3.19	1.47	1.40
2	C	502	OAA	O3-C3	5.08	1.31	1.22
2	B	500	OAA	O3-C3	5.13	1.31	1.22
2	A	501	OAA	O3-C3	5.19	1.31	1.22
2	D	503	OAA	O3-C3	5.23	1.31	1.22

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	702	CMC	N3A-C2A-N1A	-7.82	122.91	128.89
3	A	701	CMC	N3A-C2A-N1A	-7.39	123.24	128.89
3	D	703	CMC	N3A-C2A-N1A	-7.33	123.28	128.89
3	B	700	CMC	N3A-C2A-N1A	-7.28	123.32	128.89
3	B	700	CMC	P2A-O3A-P1A	-3.61	122.60	132.73
3	C	702	CMC	P2A-O3A-P1A	-3.40	123.18	132.73
3	D	703	CMC	P2A-O3A-P1A	-3.25	123.59	132.73
3	C	702	CMC	C2B-C1B-N9A	-3.11	109.53	114.29
3	D	703	CMC	C4A-C5A-N7A	-3.03	106.69	109.48
3	B	700	CMC	C4A-C5A-N7A	-2.95	106.76	109.48
3	A	701	CMC	C4A-C5A-N7A	-2.94	106.77	109.48
3	C	702	CMC	C4A-C5A-N7A	-2.92	106.79	109.48
3	D	703	CMC	C2B-C1B-N9A	-2.87	109.91	114.29
3	A	701	CMC	P2A-O3A-P1A	-2.70	125.16	132.73
2	C	502	OAA	C1-C2-C3	-2.67	110.65	115.52
2	A	501	OAA	C1-C2-C3	-2.38	111.18	115.52
3	C	702	CMC	C2A-N1A-C6A	2.01	122.36	118.77
3	C	702	CMC	C2-C1-S1P	3.64	123.07	112.89
3	A	701	CMC	C2-C1-S1P	3.83	123.62	112.89
3	D	703	CMC	C2-C1-S1P	3.91	123.83	112.89
3	B	700	CMC	C2-C1-S1P	4.25	124.78	112.89
3	A	701	CMC	C2P-S1P-C1	7.81	119.44	102.16
3	C	702	CMC	C2P-S1P-C1	7.85	119.53	102.16

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	700	CMC	C2P-S1P-C1	8.10	120.08	102.16
3	D	703	CMC	C2P-S1P-C1	8.13	120.16	102.16

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	503	OAA	1	0

## 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 ⓘ

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	381/384 (99%)	-0.34	1 (0%) 94 95	24, 36, 46, 52	0
1	B	380/384 (98%)	-0.17	7 (1%) 71 75	23, 35, 55, 69	0
1	C	379/384 (98%)	-0.25	3 (0%) 87 89	26, 37, 50, 57	0
1	D	379/384 (98%)	-0.23	2 (0%) 91 92	23, 36, 51, 66	0
All	All	1519/1536 (98%)	-0.25	13 (0%) 85 88	23, 36, 50, 69	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	382	GLU	4.6
1	B	380	ILE	3.4
1	A	3	THR	3.3
1	B	251	ILE	3.0
1	B	254	LYS	2.8
1	D	381	ALA	2.7
1	B	379	PRO	2.6
1	D	285	LYS	2.4
1	B	383	ARG	2.3
1	C	78	ILE	2.2
1	B	381	ALA	2.1
1	C	252	ASN	2.1
1	C	66	ARG	2.1

### 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 [i](#)

There are no carbohydrates in this entry.

### 6.4 Ligands [i](#)

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, 95<sup>th</sup> 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	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	OAA	B	500	9/9	0.96	0.16	1.55	31,32,34,34	0
2	OAA	A	501	9/9	0.97	0.15	0.65	32,33,33,33	0
2	OAA	C	502	9/9	0.95	0.14	0.20	29,29,29,30	0
3	CMC	D	703	52/52	0.91	0.13	-0.21	33,42,54,54	0
3	CMC	A	701	52/52	0.95	0.11	-0.58	31,33,42,44	0
3	CMC	C	702	52/52	0.94	0.11	-0.78	32,36,43,43	0
3	CMC	B	700	52/52	0.94	0.11	-0.89	33,37,44,45	0
2	OAA	D	503	9/9	0.96	0.10	-1.37	26,26,27,28	0

### 6.5 Other polymers [i](#)

There are no such residues in this entry.