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wwPDB X-ray Structure Validation Summary Report (i)

Sep 19, 2023 – 04:41 AM EDT

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PDB ID	:	5CHU
Title	:	CRYSTAL STRUCTURE OF Fox-4 cephamycinase complexed with sulfate
Authors	:	Malashkevich, V.N.; Toro, R.; Lefurgy, S.; Almo, S.C.
Deposited on	:	2015-07-10
Resolution	:	1.10 Å(reported)

1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:
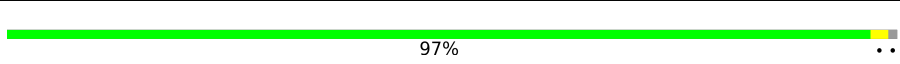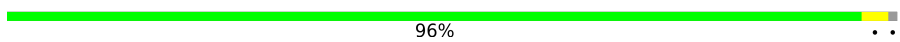*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.10 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.
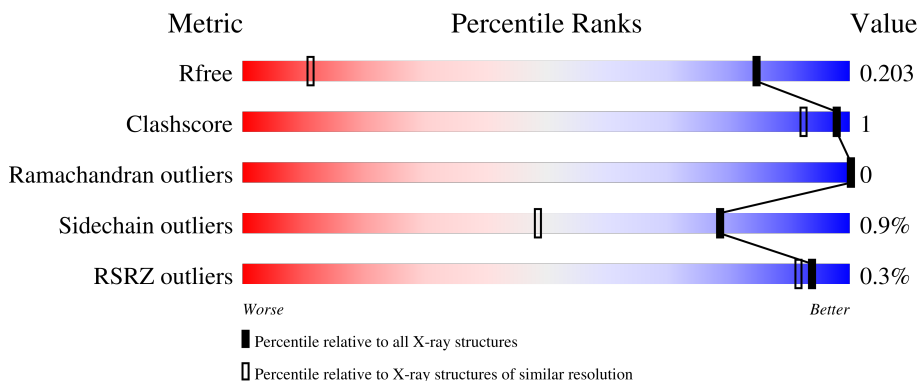
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1619 (1.14-1.06)
Clashscore	141614	1671 (1.14-1.06)
Ramachandran outliers	138981	1615 (1.14-1.06)
Sidechain outliers	138945	1613 (1.14-1.06)
RSRZ outliers	127900	1588 (1.14-1.06)

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 <=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	362	97%
1	B	362	96%

2 Entry composition ⓘ

There are 5 unique types of molecules in this entry. The entry contains 11776 atoms, of which 5432 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 Beta-lactamase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	359	Total	C	H	N	O	S	0	5	0
			5437	1748	2702	459	518	10			
1	B	359	Total	C	H	N	O	S	0	8	0
			5479	1760	2724	462	523	10			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	expression tag	UNP Q9L387
A	1	HIS	-	expression tag	UNP Q9L387
A	2	MET	-	expression tag	UNP Q9L387
B	0	GLY	-	expression tag	UNP Q9L387
B	1	HIS	-	expression tag	UNP Q9L387
B	2	MET	-	expression tag	UNP Q9L387

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	4	Total	Zn	0	0
			4	4		
2	B	4	Total	Zn	0	0
			4	4		

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O_4S).

SO4

O1

O3 O⁻ — S — O⁻ O4

S

O2

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total O S 5 4 1			0	0
3	B	1	Total O S 5 4 1			0	0
3	B	1	Total O S 5 4 1			0	0
3	B	1	Total O S 5 4 1			0	0

- Molecule 4 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).

ACT

OXT

O⁻

C

CH3 — O

O
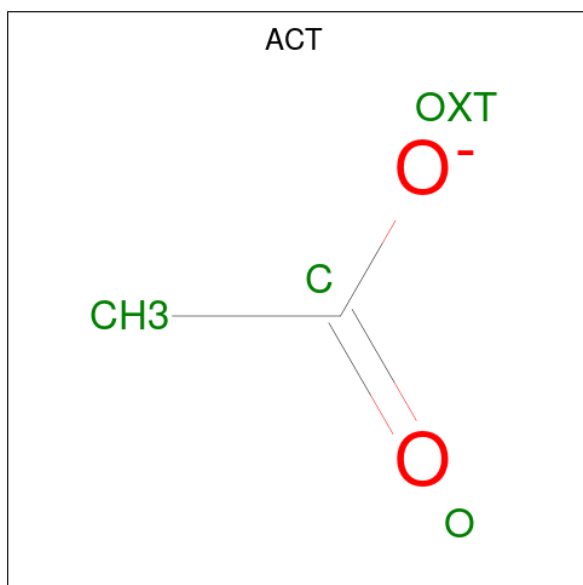
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	B	1	Total 7	C 2	H 3	O 2	0	0
4	B	1	Total 7	C 2	H 3	O 2	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	404	Total 410	O 410	0	6
5	B	406	Total 408	O 408	0	2

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 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: Beta-lactamase

Chain A:

• Molecule 1: Beta-lactamase

Chain B:

4 Data and refinement statistics ⓘ

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	53.42Å 53.92Å 116.17Å 90.00° 89.87° 90.00°	Depositor
Resolution (Å)	19.83 – 1.10 19.83 – 1.07	Depositor EDS
% Data completeness (in resolution range)	99.6 (19.83-1.10) 96.2 (19.83-1.07)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$< I/\sigma(I) >$ [1]	1.59 (at 1.07Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.188 , 0.204 0.189 , 0.203	Depositor DCC
R_{free} test set	13986 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å^2)	10.7	Xtriage
Anisotropy	0.724	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.43 , 44.5	EDS
L-test for twinning[2]	$< |L| > = 0.51, < L^2 > = 0.35$	Xtriage
Estimated twinning fraction	0.007 for k,h,-l 0.007 for -k,-h,-l 0.469 for h,-k,-l	Xtriage
F_o,F_c correlation	0.98	EDS
Total number of atoms	11776	wwPDB-VP
Average B, all atoms (Å^2)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.44% of the height of the origin peak. No significant pseudotranslation is detected.*

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[1]Intensities estimated from amplitudes.
[2]Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5   Model quality ⓘ

5.1   Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, ACT, ZN

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	0.42	0/2817	0.61	0/3828
1	B	0.42	0/2840	0.61	0/3858
All	All	0.42	0/5657	0.61	0/7686

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 ⓘ

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	2735	2702	2702	3	0
1	B	2755	2724	2723	5	1
2	A	4	0	0	0	0
2	B	4	0	0	0	1
3	A	5	0	0	0	0
3	B	15	0	0	0	0
4	B	8	6	6	1	0
5	A	410	0	0	1	4
5	B	408	0	0	2	3
All	All	6344	5432	5431	8	5

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 1.

The worst 5 of 8 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:160:HIS:NE2	5:A:768[B]:HOH:O	2.21	0.72
1:A:8:LEU:HD11	1:A:359:LEU:HD21	1.95	0.48
1:B:343:TYR:CZ	1:B:348:ARG:HG2	2.49	0.47
1:A:343:TYR:CZ	1:A:348:ARG:HG2	2.49	0.47
1:B:8:LEU:HD11	1:B:359:LEU:HD21	1.98	0.44

All (5) 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:B:186:HIS:HE2	2:B:404:ZN:ZN[2_656]	1.18	0.42
5:A:841:HOH:O	5:B:880:HOH:O[1_565]	2.11	0.09
5:A:841:HOH:O	5:B:602:HOH:O[1_565]	2.11	0.09
5:A:591:HOH:O	5:B:846:HOH:O[1_665]	2.12	0.08
5:A:743:HOH:O	5:A:891:HOH:O[2_755]	2.17	0.03

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	362/362 (100%)	350 (97%)	12 (3%)	0	100	100
1	B	365/362 (101%)	353 (97%)	12 (3%)	0	100	100
All	All	727/724 (100%)	703 (97%)	24 (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	281/278 (101%)	279 (99%)	2 (1%)	84	57
1	B	284/278 (102%)	281 (99%)	3 (1%)	73	40
All	All	565/556 (102%)	560 (99%)	5 (1%)	78	48

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	199	TYR
1	A	313	ASN
1	B	199	TYR
1	B	207	ASP
1	B	313	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 monosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 14 ligands modelled in this entry, 8 are monoatomic - leaving 6 for Mogul analysis.

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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	$\#|Z| > 2$
4	ACT	B	407	2	3,3,3	0.69	0	3,3,3	0.76	0
3	SO4	B	409	-	4,4,4	0.28	0	6,6,6	0.54	0
3	SO4	B	406	-	4,4,4	0.21	0	6,6,6	0.31	0
3	SO4	B	405	-	4,4,4	0.19	0	6,6,6	0.28	0
3	SO4	A	405	-	4,4,4	0.12	0	6,6,6	0.51	0
4	ACT	B	408	2	3,3,3	0.79	0	3,3,3	0.80	0

There are no bond length outliers.

There are no bond angle outliers.

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
4	B	407	ACT	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 ⓘ

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(Å²)	Q<0.9
1	A	359/362 (99%)	-0.43	1 (0%)	94	91	9, 15, 30, 52	0
1	B	359/362 (99%)	-0.41	1 (0%)	94	91	9, 15, 30, 55	0
All	All	718/724 (99%)	-0.42	2 (0%)	94	91	9, 15, 30, 55	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	213	THR	2.5
1	A	213	THR	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 ⓘ

There are no monosaccharides 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. 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	RSCC	RSR	B-factors(Å²)	Q<0.9
3	SO4	A	405	5/5	0.97	0.07	15,15,19,19	0
2	ZN	B	402	1/1	0.98	0.07	15,15,15,15	0
2	ZN	A	401	1/1	0.98	0.06	18,18,18,18	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	SO4	B	405	5/5	0.98	0.09	20,21,24,29	0
3	SO4	B	406	5/5	0.98	0.10	21,23,26,32	0
3	SO4	B	409	5/5	0.98	0.07	15,16,19,20	0
4	ACT	B	407	4/4	0.98	0.06	13,15,18,18	0
4	ACT	B	408	4/4	0.98	0.05	13,15,19,19	0
2	ZN	A	402	1/1	0.99	0.07	15,15,15,15	0
2	ZN	B	404	1/1	0.99	0.04	19,19,19,19	0
2	ZN	A	404	1/1	0.99	0.04	17,17,17,17	0
2	ZN	B	401	1/1	0.99	0.05	19,19,19,19	0
2	ZN	B	403	1/1	1.00	0.04	14,14,14,14	0
2	ZN	A	403	1/1	1.00	0.05	14,14,14,14	0

6.5 Other polymers ⓘ

There are no such residues in this entry.