Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 04:03 PM GMT

PDB ID	:	4E57
Title	:	Crystal Structure of spacer 6aa-shortened cephalosporin acylase mutant
Authors	:	Yin, J.; Zhang, Z.; Wu, G.; Huang, X.
Deposited on	:	2012-03-14
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 (1)) were used in the production of this report:

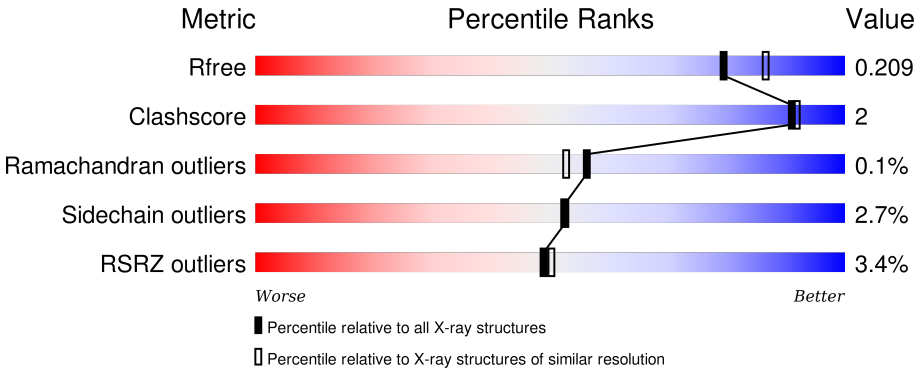
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)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

1 Overall quality at a glance (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))
R_{free}	91344	6249 (2.00-2.00)
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (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.

Mol	Chain	Length	Quality of chain
1	A	698	3% 90% 6%
1	B	698	3% 91% 5%

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	CD	B	701	-	-	-	X

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 12015 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 Cephalosporin acylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	677	Total	C	N	O	S	0	0	0
			5329	3373	940	1004	12			
1	B	677	Total	C	N	O	S	0	0	0
			5329	3373	940	1004	12			

There are 46 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	EXPRESSION TAG	UNP O86089
A	2	GLY	-	EXPRESSION TAG	UNP O86089
A	3	ILE	-	EXPRESSION TAG	UNP O86089
A	127	GLU	ASP	SEE REMARK 999	UNP O86089
A	?	-	PRO	DELETION	UNP O86089
A	?	-	ASP	DELETION	UNP O86089
A	?	-	LEU	DELETION	UNP O86089
A	?	-	ALA	DELETION	UNP O86089
A	?	-	ASP	DELETION	UNP O86089
A	?	-	GLN	DELETION	UNP O86089
A	166	ALA	SER	ENGINEERED MUTATION	UNP O86089
A	687	LEU	-	EXPRESSION TAG	UNP O86089
A	688	ALA	-	EXPRESSION TAG	UNP O86089
A	689	ALA	-	EXPRESSION TAG	UNP O86089
A	690	ALA	-	EXPRESSION TAG	UNP O86089
A	691	LEU	-	EXPRESSION TAG	UNP O86089
A	692	GLU	-	EXPRESSION TAG	UNP O86089
A	693	HIS	-	EXPRESSION TAG	UNP O86089
A	694	HIS	-	EXPRESSION TAG	UNP O86089
A	695	HIS	-	EXPRESSION TAG	UNP O86089
A	696	HIS	-	EXPRESSION TAG	UNP O86089
A	697	HIS	-	EXPRESSION TAG	UNP O86089
A	698	HIS	-	EXPRESSION TAG	UNP O86089
B	1	MET	-	EXPRESSION TAG	UNP O86089
B	2	GLY	-	EXPRESSION TAG	UNP O86089

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Chain	Residue	Modelled	Actual	Comment	Reference
B	3	ILE	-	EXPRESSION TAG	UNP O86089
B	127	GLU	ASP	SEE REMARK 999	UNP O86089
B	?	-	PRO	DELETION	UNP O86089
B	?	-	ASP	DELETION	UNP O86089
B	?	-	LEU	DELETION	UNP O86089
B	?	-	ALA	DELETION	UNP O86089
B	?	-	ASP	DELETION	UNP O86089
B	?	-	GLN	DELETION	UNP O86089
B	166	ALA	SER	ENGINEERED MUTATION	UNP O86089
B	687	LEU	-	EXPRESSION TAG	UNP O86089
B	688	ALA	-	EXPRESSION TAG	UNP O86089
B	689	ALA	-	EXPRESSION TAG	UNP O86089
B	690	ALA	-	EXPRESSION TAG	UNP O86089
B	691	LEU	-	EXPRESSION TAG	UNP O86089
B	692	GLU	-	EXPRESSION TAG	UNP O86089
B	693	HIS	-	EXPRESSION TAG	UNP O86089
B	694	HIS	-	EXPRESSION TAG	UNP O86089
B	695	HIS	-	EXPRESSION TAG	UNP O86089
B	696	HIS	-	EXPRESSION TAG	UNP O86089
B	697	HIS	-	EXPRESSION TAG	UNP O86089
B	698	HIS	-	EXPRESSION TAG	UNP O86089

- Molecule 2 is CADMIUM ION (three-letter code: CD) (formula: Cd).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	14	Total Cd 14 14	0	0
2	A	19	Total Cd 19 19	0	0

- Molecule 3 is water.

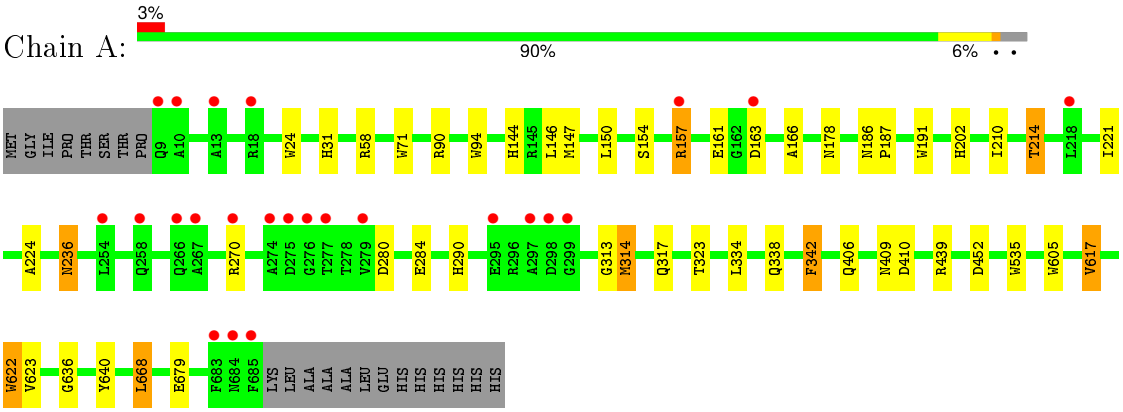
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	681	Total O 681 681	0	0
3	B	643	Total O 643 643	0	0

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.

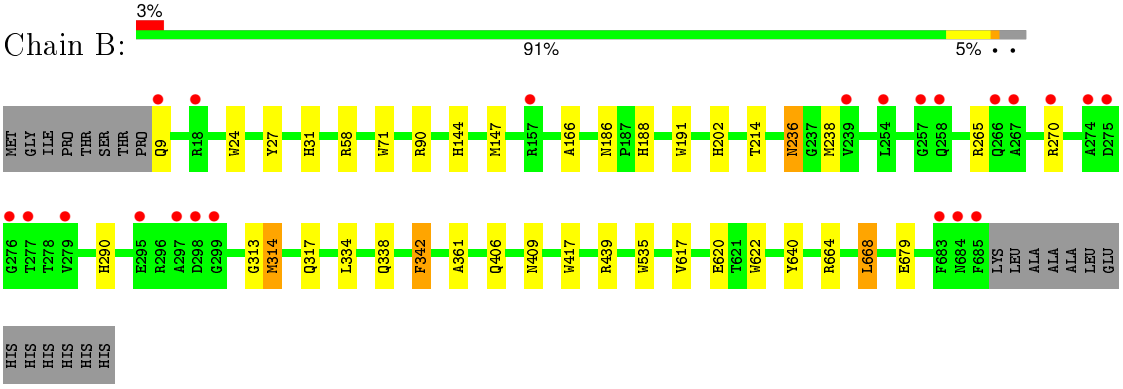
- Molecule 1: Cephalosporin acylase

Chain A:



- Molecule 1: Cephalosporin acylase

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 41	Depositor
Cell constants a, b, c, α, β, γ	73.60Å 73.60Å 383.83Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.28 – 2.00 48.28 – 2.00	Depositor EDS
% Data completeness (in resolution range)	100.0 (48.28-2.00) 100.0 (48.28-2.00)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.65 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.182 , 0.207 0.185 , 0.209	Depositor DCC
R_{free} test set	6866 reflections (5.29%)	DCC
Wilson B-factor (Å^2)	26.2	Xtriage
Anisotropy	0.046	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.38 , 39.4	EDS
Estimated twinning fraction	0.470 for h,-k,-l	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.51$, $< L^2 > = 0.35$	Xtriage
Outliers	1 of 136599 reflections (0.001%)	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	12015	wwPDB-VP
Average B, all atoms (Å^2)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.81% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: CD

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	0.50	6/5484 (0.1%)	0.54	2/7493 (0.0%)
1	B	0.50	5/5484 (0.1%)	0.54	1/7493 (0.0%)
All	All	0.50	11/10968 (0.1%)	0.54	3/14986 (0.0%)

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	71	TRP	CD2-CE2	5.21	1.47	1.41
1	A	191	TRP	CD2-CE2	5.17	1.47	1.41
1	B	417	TRP	CD2-CE2	5.13	1.47	1.41
1	A	605	TRP	CD2-CE2	5.12	1.47	1.41
1	B	191	TRP	CD2-CE2	5.10	1.47	1.41
1	B	535	TRP	CD2-CE2	5.10	1.47	1.41
1	B	622	TRP	CD2-CE2	5.09	1.47	1.41
1	B	71	TRP	CD2-CE2	5.08	1.47	1.41
1	A	535	TRP	CD2-CE2	5.07	1.47	1.41
1	A	622	TRP	CD2-CE2	5.02	1.47	1.41
1	A	94	TRP	CD2-CE2	5.01	1.47	1.41

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	668	LEU	CA-CB-CG	5.96	129.00	115.30
1	B	668	LEU	CA-CB-CG	5.79	128.60	115.30
1	A	163	ASP	CB-CG-OD2	5.21	122.99	118.30

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	5329	0	5070	28	0
1	B	5329	0	5071	17	0
2	A	19	0	0	0	0
2	B	14	0	0	0	0
3	A	681	0	0	2	0
3	B	643	0	0	1	0
All	All	12015	0	10141	45	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 2.

All (45) 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:B:317:GLN:HE22	1:B:338:GLN:H	1.25	0.84
1:A:317:GLN:HE22	1:A:338:GLN:H	1.27	0.82
1:B:166:ALA:H	1:B:409:ASN:HD21	1.26	0.81
1:A:166:ALA:H	1:A:409:ASN:HD21	1.25	0.79
1:A:409:ASN:HD22	1:A:439:ARG:HH22	1.38	0.71
1:B:313:GLY:H	1:B:338:GLN:HE21	1.38	0.70
1:A:313:GLY:H	1:A:338:GLN:HE21	1.38	0.69
1:A:90:ARG:HH11	1:A:144:HIS:HE2	1.39	0.69
1:B:409:ASN:HD22	1:B:439:ARG:HH22	1.38	0.68
1:B:90:ARG:HH11	1:B:144:HIS:HE2	1.38	0.68
1:B:317:GLN:NE2	1:B:338:GLN:H	1.94	0.63
1:A:161:GLU:HA	1:A:617:VAL:HG13	1.83	0.59
1:A:317:GLN:NE2	1:A:338:GLN:H	1.96	0.59
1:A:290:HIS:HE1	3:A:1080:HOH:O	1.87	0.58
1:B:147:MET:CE	1:B:342:PHE:HE1	2.18	0.57
1:B:290:HIS:HE1	3:B:1045:HOH:O	1.88	0.56
1:A:178:ASN:ND2	3:A:1392:HOH:O	2.41	0.54
1:A:31:HIS:ND1	1:A:202:HIS:HD2	2.08	0.51
1:A:147:MET:CE	1:A:342:PHE:CE1	2.94	0.50
1:A:147:MET:CE	1:A:342:PHE:HE1	2.25	0.49
1:A:147:MET:HE1	1:A:342:PHE:HE1	1.78	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:210:ILE:HD11	1:A:323:THR:CG2	2.44	0.47
1:B:31:HIS:ND1	1:B:202:HIS:HD2	2.12	0.46
1:A:314:MET:HB2	1:A:338:GLN:O	2.14	0.46
1:B:313:GLY:H	1:B:338:GLN:NE2	2.12	0.45
1:B:314:MET:HB2	1:B:338:GLN:O	2.17	0.45
1:A:187:PRO:HD2	1:A:622:TRP:O	2.16	0.44
1:B:188:HIS:HE1	1:B:620:GLU:OE1	2.00	0.43
1:A:157:ARG:NE	1:A:157:ARG:HA	2.33	0.43
1:A:31:HIS:ND1	1:A:202:HIS:CD2	2.87	0.43
1:B:147:MET:HE1	1:B:342:PHE:HE1	1.84	0.43
1:A:24:TRP:HB2	1:A:679:GLU:HB2	2.00	0.43
1:B:24:TRP:HB2	1:B:679:GLU:HB2	2.01	0.43
1:A:210:ILE:HG23	1:A:224:ALA:HB1	2.00	0.42
1:B:236:ASN:ND2	1:B:342:PHE:H	2.18	0.42
1:A:157:ARG:HE	1:A:157:ARG:HA	1.85	0.42
1:B:27:TYR:CD1	1:B:664:ARG:HD3	2.55	0.41
1:A:623:VAL:O	1:A:636:GLY:HA2	2.20	0.41
1:A:236:ASN:ND2	1:A:342:PHE:H	2.19	0.41
1:A:166:ALA:HB2	1:A:187:PRO:HA	2.03	0.41
1:B:238:MET:SD	1:B:361:ALA:HB2	2.61	0.41
1:A:147:MET:HE1	1:A:342:PHE:CE1	2.54	0.40
1:A:410:ASP:OD1	1:A:410:ASP:N	2.52	0.40
1:A:146:LEU:HA	1:A:150:LEU:HD12	2.03	0.40
1:A:214:THR:HB	1:A:221:ILE:HA	2.03	0.40

There are no symmetry-related clashes.

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	675/698 (97%)	658 (98%)	16 (2%)	1 (0%)	56 53

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	B	675/698 (97%)	660 (98%)	14 (2%)	1 (0%)	56 53
All	All	1350/1396 (97%)	1318 (98%)	30 (2%)	2 (0%)	56 53

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	342	PHE
1	A	342	PHE

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	547/564 (97%)	531 (97%)	16 (3%)	50 49
1	B	547/564 (97%)	534 (98%)	13 (2%)	57 58
All	All	1094/1128 (97%)	1065 (97%)	29 (3%)	52 52

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	58	ARG
1	A	154	SER
1	A	157	ARG
1	A	186	ASN
1	A	214	THR
1	A	236	ASN
1	A	270	ARG
1	A	280	ASP
1	A	284	GLU
1	A	314	MET
1	A	334	LEU
1	A	406	GLN
1	A	452	ASP
1	A	617	VAL

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Mol	Chain	Res	Type
1	A	640	TYR
1	A	668	LEU
1	B	9	GLN
1	B	58	ARG
1	B	186	ASN
1	B	214	THR
1	B	236	ASN
1	B	265	ARG
1	B	270	ARG
1	B	314	MET
1	B	334	LEU
1	B	406	GLN
1	B	617	VAL
1	B	640	TYR
1	B	668	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	148	ASN
1	A	188	HIS
1	A	202	HIS
1	A	236	ASN
1	A	249	GLN
1	A	290	HIS
1	A	317	GLN
1	A	338	GLN
1	A	354	ASN
1	A	406	GLN
1	A	409	ASN
1	A	442	GLN
1	B	9	GLN
1	B	148	ASN
1	B	188	HIS
1	B	202	HIS
1	B	236	ASN
1	B	249	GLN
1	B	290	HIS
1	B	317	GLN
1	B	338	GLN
1	B	354	ASN
1	B	406	GLN

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Mol	Chain	Res	Type
1	B	409	ASN
1	B	523	GLN

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

Of 33 ligands modelled in this entry, 33 are monoatomic - leaving 0 for Mogul analysis.

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.

No monomer is involved in short contacts.

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, 95^{th} 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	677/698 (96%)	0.08	24 (3%) 48 49	18, 26, 44, 55	0
1	B	677/698 (96%)	0.09	22 (3%) 51 52	18, 26, 45, 56	1 (0%)
All	All	1354/1396 (96%)	0.09	46 (3%) 49 50	18, 26, 44, 56	1 (0%)

All (46) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	239	VAL	7.2
1	B	684	ASN	4.0
1	B	279	VAL	3.8
1	B	18	ARG	3.7
1	A	684	ASN	3.7
1	A	276	GLY	3.5
1	B	274	ALA	3.4
1	A	18	ARG	3.3
1	A	258	GLN	3.3
1	A	685	PHE	3.2
1	B	297	ALA	3.1
1	B	257	GLY	3.1
1	A	279	VAL	3.0
1	B	685	PHE	3.0
1	B	9	GLN	3.0
1	A	275	ASP	2.9
1	B	276	GLY	2.8
1	A	297	ALA	2.8
1	A	9	GLN	2.7
1	A	270	ARG	2.7
1	A	299	GLY	2.7
1	B	298	ASP	2.6
1	B	258	GLN	2.6
1	A	277	THR	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	267	ALA	2.5
1	B	270	ARG	2.5
1	B	275	ASP	2.5
1	A	274	ALA	2.5
1	B	277	THR	2.5
1	A	298	ASP	2.4
1	A	254	LEU	2.4
1	B	683	PHE	2.4
1	A	13	ALA	2.4
1	A	267	ALA	2.4
1	B	295	GLU	2.4
1	A	157	ARG	2.3
1	B	266	GLN	2.2
1	A	295	GLU	2.2
1	A	218	LEU	2.2
1	B	299	GLY	2.2
1	A	10	ALA	2.2
1	A	266	GLN	2.1
1	A	163	ASP	2.1
1	B	254	LEU	2.1
1	A	683	PHE	2.1
1	B	157	ARG	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, 95^{th} 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(Å^2)	Q<0.9
2	CD	B	701	1/1	0.97	0.22	2.99	14,14,14,14	1
2	CD	A	706	1/1	0.97	0.07	-1.63	30,30,30,30	0
2	CD	B	704	1/1	0.97	0.05	-2.00	36,36,36,36	0
2	CD	A	707	1/1	0.98	0.05	-2.12	36,36,36,36	0
2	CD	B	702	1/1	0.88	0.11	-	75,75,75,75	0
2	CD	A	719	1/1	0.85	0.14	-	129,129,129,129	0
2	CD	B	705	1/1	0.72	0.10	-	45,45,45,45	0
2	CD	A	703	1/1	0.85	0.13	-	98,98,98,98	0
2	CD	A	717	1/1	0.89	0.11	-	73,73,73,73	0
2	CD	A	705	1/1	0.80	0.28	-	123,123,123,123	0
2	CD	B	707	1/1	0.95	0.09	-	50,50,50,50	1
2	CD	A	701	1/1	0.85	0.12	-	83,83,83,83	0
2	CD	B	710	1/1	0.88	0.08	-	63,63,63,63	1
2	CD	A	715	1/1	0.93	0.07	-	61,61,61,61	0
2	CD	A	708	1/1	0.89	0.07	-	45,45,45,45	0
2	CD	B	713	1/1	0.77	0.19	-	128,128,128,128	0
2	CD	B	708	1/1	0.92	0.06	-	52,52,52,52	1
2	CD	A	718	1/1	0.70	0.14	-	81,81,81,81	1
2	CD	A	711	1/1	0.95	0.05	-	52,52,52,52	1
2	CD	B	711	1/1	0.92	0.08	-	57,57,57,57	1
2	CD	A	714	1/1	0.93	0.07	-	63,63,63,63	1
2	CD	A	710	1/1	0.97	0.10	-	51,51,51,51	1
2	CD	A	702	1/1	0.86	0.17	-	109,109,109,109	0
2	CD	A	704	1/1	0.79	0.23	-	105,105,105,105	0
2	CD	A	709	1/1	0.93	0.04	-	52,52,52,52	1
2	CD	A	713	1/1	0.93	0.09	-	55,55,55,55	1
2	CD	A	716	1/1	0.75	0.16	-	129,129,129,129	0
2	CD	B	714	1/1	0.82	0.17	-	80,80,80,80	1
2	CD	A	712	1/1	0.98	0.06	-	52,52,52,52	1
2	CD	B	703	1/1	0.79	0.24	-	117,117,117,117	0
2	CD	B	706	1/1	0.95	0.05	-	54,54,54,54	1
2	CD	B	712	1/1	0.94	0.08	-	62,62,62,62	0
2	CD	B	709	1/1	0.94	0.06	-	53,53,53,53	1

6.5 Other polymers

There are no such residues in this entry.