Full wwPDB X-ray Structure Validation Report (i)

Jun 19, 2024 – 10:03 AM EDT

PDB ID	:	4EPA
Title	:	The crystal structure of the ferric yersiniabactin uptake receptor FyuA from Yersinia pestis
Authors	:	Lukacik, P.; Barnard, T.J.; Buchanan, S.K.
Deposited on	:	2012-04-17
Resolution	:	3.20 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
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

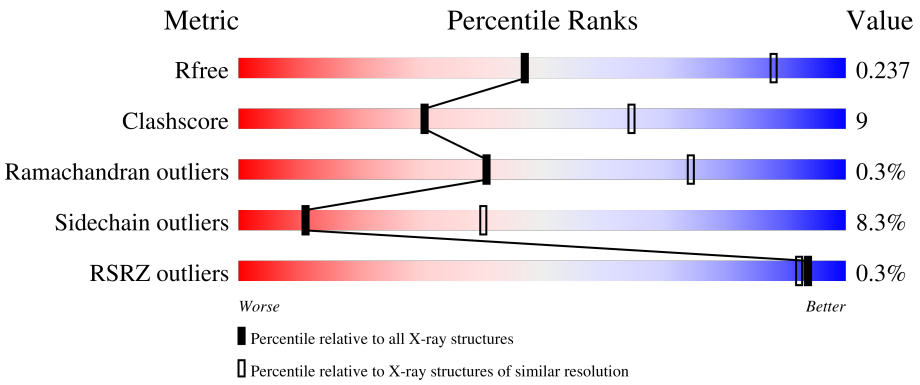
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

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 3.20 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	1133 (3.20-3.20)
Clashscore	141614	1253 (3.20-3.20)
Ramachandran outliers	138981	1234 (3.20-3.20)
Sidechain outliers	138945	1233 (3.20-3.20)
RSRZ outliers	127900	1095 (3.20-3.20)

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	655	75% 20% • •

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 4884 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.

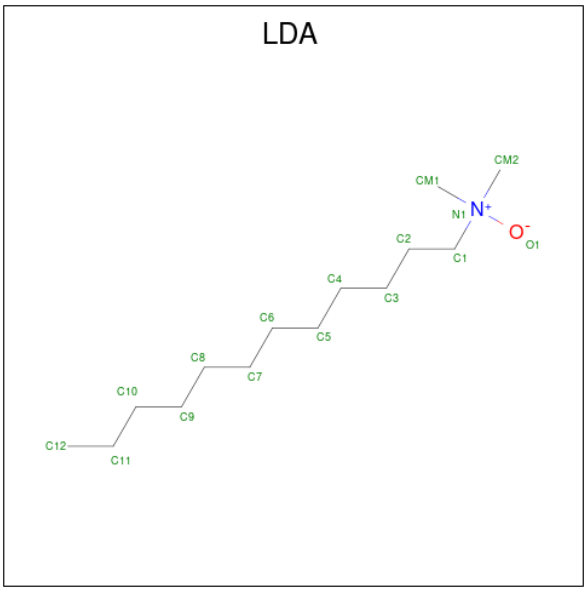
- Molecule 1 is a protein called Pesticin receptor.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	A	632	4863	3029	855	962	2	15	0	1	0

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	GLY	-	EXPRESSION TAG	UNP P46359
A	-2	ALA	-	EXPRESSION TAG	UNP P46359
A	-1	MSE	-	EXPRESSION TAG	UNP P46359
A	0	GLY	-	EXPRESSION TAG	UNP P46359

- Molecule 2 is LAURYL DIMETHYLAMINE-N-OXIDE (three-letter code: LDA) (formula: $C_{14}H_{31}NO$).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	C		
2	A	1	11	11	0	0

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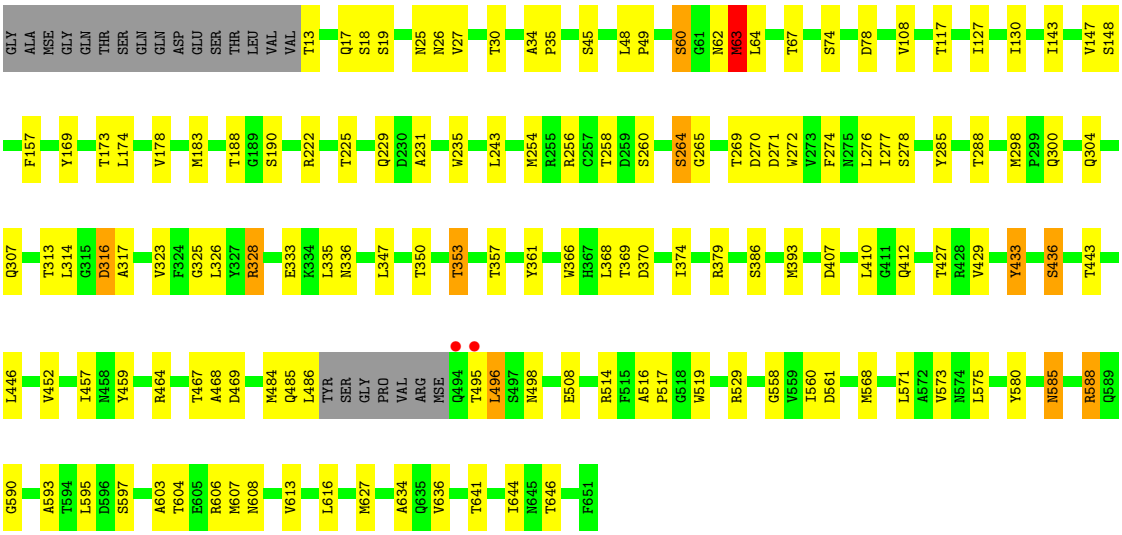
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	C		
2	A	1	10	10	0	0

3 Residue-property plots (i)

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: Pesticin receptor

Chain A:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	I 41 2 2	Depositor
Cell constants a, b, c, α, β, γ	181.18Å 181.18Å 319.86Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 3.20 49.99 – 3.20	Depositor EDS
% Data completeness (in resolution range)	97.7 (50.00-3.20) 97.7 (49.99-3.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.25	Depositor
$< I/\sigma(I) >$ [1]	2.39 (at 3.19Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.227 , 0.244 0.222 , 0.237	Depositor DCC
R_{free} test set	2179 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å^2)	65.2	Xtriage
Anisotropy	0.013	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.28 , 14.8	EDS
L-test for twinning[2]	$< \|L\| > = 0.47$, $< L^2 > = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.91	EDS
Total number of atoms	4884	wwPDB-VP
Average B, all atoms (Å^2)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.48% 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, 0.333 respectively for untwinned datasets, and 0.375, 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: LDA

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	0.62	0/4953	0.71	0/6713

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	0	2

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	590	GLY	Peptide
1	A	74	SER	Peptide

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	4863	0	4641	88	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	21	0	40	0	0
All	All	4884	0	4681	88	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 9.

All (88) 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:333:GLU:OE1	1:A:353:THR:HG23	1.52	1.10
1:A:63:MSE:HE1	1:A:225:THR:HG21	1.36	1.02
1:A:347:LEU:HD22	1:A:393:MSE:HE2	1.55	0.89
1:A:60:SER:HB3	1:A:67:THR:HG23	1.61	0.82
1:A:64:LEU:HD21	1:A:254:MSE:HE1	1.74	0.69
1:A:410:LEU:HD11	1:A:433:TYR:CD1	2.28	0.69
1:A:63:MSE:CE	1:A:225:THR:HG21	2.20	0.67
1:A:264:SER:OG	1:A:265:GLY:N	2.29	0.65
1:A:467:THR:HG22	1:A:469:ASP:H	1.61	0.65
1:A:298:MSE:HE3	1:A:335:LEU:HD21	1.78	0.64
1:A:269:THR:HG22	1:A:271:ASP:H	1.62	0.63
1:A:369:THR:HG22	1:A:370:ASP:N	2.14	0.63
1:A:127:ILE:C	1:A:127:ILE:HD12	2.18	0.63
1:A:269:THR:HG22	1:A:270:ASP:N	2.14	0.62
1:A:467:THR:HG22	1:A:468:ALA:N	2.15	0.61
1:A:127:ILE:HD12	1:A:127:ILE:O	2.03	0.59
1:A:485:GLN:HB3	1:A:496:LEU:HD21	1.85	0.58
1:A:173:THR:O	1:A:174:LEU:HD23	2.03	0.58
1:A:231:ALA:HB1	1:A:243:LEU:HD22	1.85	0.58
1:A:369:THR:HG22	1:A:370:ASP:H	1.68	0.57
1:A:269:THR:CG2	1:A:270:ASP:N	2.68	0.56
1:A:613:VAL:O	1:A:613:VAL:HG13	2.07	0.55
1:A:143:ILE:HD11	1:A:157:PHE:HD1	1.72	0.54
1:A:366:TRP:CD2	1:A:368:LEU:HD11	2.42	0.54
1:A:386:SER:O	1:A:386:SER:OG	2.15	0.54
1:A:285:TYR:OH	1:A:300:GLN:OE1	2.23	0.53
1:A:467:THR:CG2	1:A:468:ALA:N	2.71	0.53
1:A:560:ILE:HG22	1:A:561:ASP:N	2.24	0.52
1:A:235:TRP:CD1	1:A:235:TRP:O	2.63	0.52
1:A:25:ASN:O	1:A:464[A]:ARG:NH2	2.42	0.52
1:A:78:ASP:HB2	1:A:496:LEU:HD11	1.92	0.52
1:A:62:ASN:ND2	1:A:254:MSE:SE	2.94	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:568:MSE:HE1	1:A:608:ASN:ND2	2.27	0.50
1:A:17:GLN:HB3	1:A:323:VAL:CG2	2.41	0.50
1:A:34:ALA:N	1:A:35:PRO:HD2	2.26	0.50
1:A:568:MSE:HE1	1:A:608:ASN:HD21	1.77	0.49
1:A:143:ILE:HD11	1:A:157:PHE:CD1	2.47	0.49
1:A:516:ALA:HB1	1:A:517:PRO:CD	2.42	0.48
1:A:585:ASN:HD22	1:A:585:ASN:N	2.11	0.48
1:A:235:TRP:O	1:A:235:TRP:CG	2.66	0.48
1:A:429:VAL:HG22	1:A:459:TYR:HD2	1.79	0.48
1:A:325:GLY:O	1:A:326:LEU:HD23	2.14	0.47
1:A:595:LEU:HD23	1:A:616:LEU:HD23	1.96	0.47
1:A:231:ALA:CB	1:A:243:LEU:HD22	2.45	0.47
1:A:62:ASN:HD22	1:A:254:MSE:SE	2.48	0.46
1:A:328:ARG:HA	1:A:357:THR:O	2.15	0.46
1:A:117:THR:HB	1:A:412:GLN:NE2	2.31	0.45
1:A:183:MSE:HE1	1:A:636:VAL:HG23	1.99	0.45
1:A:316:ASP:CB	1:A:317:ALA:HA	2.47	0.45
1:A:108:VAL:HG12	1:A:130:ILE:HG12	1.98	0.45
1:A:60:SER:O	1:A:60:SER:OG	2.30	0.45
1:A:26:ASN:O	1:A:27:VAL:HG23	2.16	0.45
1:A:366:TRP:CD1	1:A:368:LEU:HD12	2.52	0.45
1:A:580:TYR:CE2	1:A:588:ARG:HB2	2.52	0.45
1:A:48:LEU:N	1:A:49:PRO:CD	2.80	0.45
1:A:484:MSE:HE2	1:A:486:LEU:HD11	1.99	0.44
1:A:178:VAL:O	1:A:178:VAL:HG12	2.17	0.44
1:A:183:MSE:HE1	1:A:636:VAL:CG2	2.48	0.44
1:A:269:THR:CG2	1:A:270:ASP:H	2.31	0.44
1:A:436:SER:HB2	1:A:452:VAL:HA	2.00	0.44
1:A:361:TYR:HB3	1:A:379:ARG:HG3	1.99	0.44
1:A:410:LEU:HD23	1:A:410:LEU:N	2.33	0.44
1:A:147:VAL:HG12	1:A:644:ILE:HG22	2.00	0.44
1:A:374:ILE:O	1:A:374:ILE:CG2	2.66	0.43
1:A:336:ASN:ND2	1:A:350:THR:HG23	2.33	0.43
1:A:467:THR:CG2	1:A:468:ALA:H	2.30	0.43
1:A:17:GLN:HB3	1:A:323:VAL:HG21	2.01	0.43
1:A:256:ARG:HB2	1:A:285:TYR:HB3	2.00	0.43
1:A:274:PHE:CD2	1:A:274:PHE:N	2.87	0.43
1:A:169:TYR:CD2	1:A:169:TYR:N	2.87	0.42
1:A:272:TRP:CE2	1:A:313:THR:HG22	2.54	0.42
1:A:603:ALA:HB3	1:A:607:MSE:HE2	2.02	0.42
1:A:277:ILE:CG2	1:A:278:SER:N	2.81	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:519:TRP:CD2	1:A:558:GLY:HA3	2.55	0.42
1:A:235:TRP:CE3	1:A:634:ALA:HB2	2.55	0.42
1:A:316:ASP:HB2	1:A:317:ALA:HA	2.01	0.42
1:A:457:ILE:O	1:A:457:ILE:HG23	2.19	0.42
1:A:571:LEU:HD12	1:A:571:LEU:HA	1.91	0.42
1:A:60:SER:HB3	1:A:67:THR:CG2	2.40	0.41
1:A:17:GLN:HB3	1:A:323:VAL:HG22	2.02	0.41
1:A:366:TRP:CG	1:A:368:LEU:HD12	2.56	0.41
1:A:298:MSE:CE	1:A:335:LEU:HD21	2.47	0.41
1:A:369:THR:CG2	1:A:370:ASP:N	2.81	0.41
1:A:608:ASN:OD1	1:A:608:ASN:C	2.59	0.41
1:A:516:ALA:HB1	1:A:517:PRO:HD2	2.03	0.41
1:A:429:VAL:HG22	1:A:459:TYR:CD2	2.57	0.40
1:A:62:ASN:O	1:A:64:LEU:N	2.54	0.40
1:A:575:LEU:HD12	1:A:593:ALA:HB2	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	629/655 (96%)	584 (93%)	43 (7%)	2 (0%)	41 74

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	63	MSE
1	A	229	GLN

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	518/529 (98%)	475 (92%)	43 (8%)	11 40

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

Mol	Chain	Res	Type
1	A	13	THR
1	A	18	SER
1	A	19	SER
1	A	30	THR
1	A	45	SER
1	A	60	SER
1	A	63	MSE
1	A	148	SER
1	A	188	THR
1	A	190	SER
1	A	222	ARG
1	A	258	THR
1	A	260	SER
1	A	264	SER
1	A	276	LEU
1	A	288	THR
1	A	304	GLN
1	A	307	GLN
1	A	314	LEU
1	A	316	ASP
1	A	328	ARG
1	A	353	THR
1	A	407	ASP
1	A	427	THR
1	A	433	TYR
1	A	436	SER
1	A	443	THR
1	A	446	LEU
1	A	495	THR
1	A	496	LEU

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Mol	Chain	Res	Type
1	A	498	ASN
1	A	508	GLU
1	A	514	ARG
1	A	529	ARG
1	A	573	VAL
1	A	585	ASN
1	A	588	ARG
1	A	597	SER
1	A	604	THR
1	A	606	ARG
1	A	627	MSE
1	A	641	THR
1	A	646	THR

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

Mol	Chain	Res	Type
1	A	59	ASN
1	A	229	GLN

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 monosaccharides 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	LDA	A	701	-	10,10,15	0.46	0	9,9,17	0.35	0
2	LDA	A	702	-	9,9,15	0.50	0	8,8,17	0.28	0

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	LDA	A	701	-	-	3/8/8/13	-
2	LDA	A	702	-	-	1/7/7/13	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	701	LDA	C11-C10-C9-C8
2	A	701	LDA	C7-C8-C9-C10
2	A	701	LDA	C5-C6-C7-C8
2	A	702	LDA	C4-C5-C6-C7

There are no ring outliers.

No monomer is involved in short contacts.

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 (i)

6.1 Protein, DNA and RNA chains (i)

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	617/655 (94%)	-0.36	2 (0%) 94 92	20, 40, 66, 91	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	494	GLN	2.6
1	A	495	THR	2.6

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

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

6.3 Carbohydrates (i)

There are no monosaccharides 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. 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(Å^2)	Q<0.9
2	LDA	A	702	10/16	0.93	0.27	48,57,60,61	0
2	LDA	A	701	11/16	0.94	0.23	36,50,68,69	0

6.5 Other polymers (i)

There are no such residues in this entry.