



wwPDB X-ray Structure Validation Summary Report ⓘ

Nov 8, 2017 – 11:31 PM EST

PDB ID : 4MB8
Title : Evolutionary history and metabolic insights of ancient mammalian uricases
Authors : Ortlund, E.O.; Murphy, M.N.
Deposited on : unknown
Resolution : 2.40 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20030345
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
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) : rb-20030345

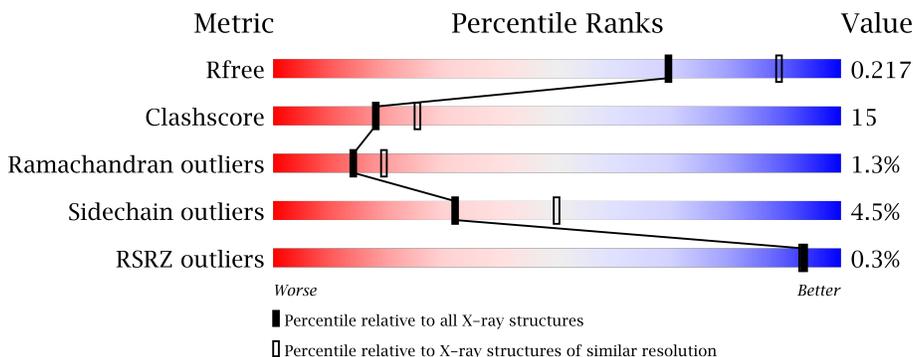
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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}	100719	3166 (2.40-2.40)
Clashscore	112137	3674 (2.40-2.40)
Ramachandran outliers	110173	3616 (2.40-2.40)
Sidechain outliers	110143	3617 (2.40-2.40)
RSRZ outliers	101464	3195 (2.40-2.40)

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 $\leq 5\%$. 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	304	 .% 67% 23% . . 6%
1	B	304	 70% 21% . 6%
1	C	304	 68% 24% . 6%
1	D	304	 .% 67% 25% . . 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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	ACT	B	401	-	-	-	X
2	ACT	D	402	-	-	-	X

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 9529 atoms, of which 6 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 Uricase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	285	2304	1480	391	423	10	0	0	0
1	B	286	2313	1486	393	424	10	0	0	0
1	C	285	2302	1482	389	421	10	0	0	0
1	D	290	2344	1504	400	430	10	0	0	0

There are 52 discrepancies between the modelled and reference sequences:

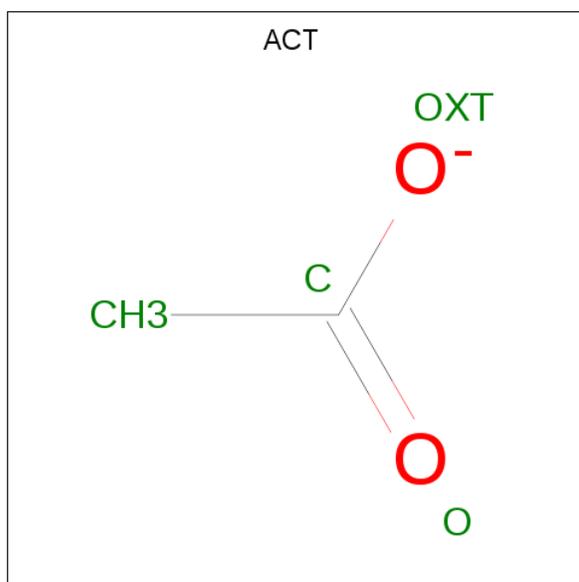
Chain	Residue	Modelled	Actual	Comment	Reference
A	58	LEU	VAL	ENGINEERED MUTATION	UNP Q5FZI9
A	59	HIS	TYR	ENGINEERED MUTATION	UNP Q5FZI9
A	89	ALA	THR	ENGINEERED MUTATION	UNP Q5FZI9
A	134	THR	ASN	ENGINEERED MUTATION	UNP Q5FZI9
A	147	LYS	ARG	ENGINEERED MUTATION	UNP Q5FZI9
A	192	GLN	LYS	ENGINEERED MUTATION	UNP Q5FZI9
A	220	LYS	GLU	ENGINEERED MUTATION	UNP Q5FZI9
A	245	LEU	HIS	ENGINEERED MUTATION	UNP Q5FZI9
A	253	ILE	MET	ENGINEERED MUTATION	UNP Q5FZI9
A	291	LYS	ARG	ENGINEERED MUTATION	UNP Q5FZI9
A	296	VAL	ALA	ENGINEERED MUTATION	UNP Q5FZI9
A	301	SER	ALA	ENGINEERED MUTATION	UNP Q5FZI9
A	303	ARG	LYS	ENGINEERED MUTATION	UNP Q5FZI9
B	58	LEU	VAL	ENGINEERED MUTATION	UNP Q5FZI9
B	59	HIS	TYR	ENGINEERED MUTATION	UNP Q5FZI9
B	89	ALA	THR	ENGINEERED MUTATION	UNP Q5FZI9
B	134	THR	ASN	ENGINEERED MUTATION	UNP Q5FZI9
B	147	LYS	ARG	ENGINEERED MUTATION	UNP Q5FZI9
B	192	GLN	LYS	ENGINEERED MUTATION	UNP Q5FZI9
B	220	LYS	GLU	ENGINEERED MUTATION	UNP Q5FZI9
B	245	LEU	HIS	ENGINEERED MUTATION	UNP Q5FZI9

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Chain	Residue	Modelled	Actual	Comment	Reference
B	253	ILE	MET	ENGINEERED MUTATION	UNP Q5FZ19
B	291	LYS	ARG	ENGINEERED MUTATION	UNP Q5FZ19
B	296	VAL	ALA	ENGINEERED MUTATION	UNP Q5FZ19
B	301	SER	ALA	ENGINEERED MUTATION	UNP Q5FZ19
B	303	ARG	LYS	ENGINEERED MUTATION	UNP Q5FZ19
C	58	LEU	VAL	ENGINEERED MUTATION	UNP Q5FZ19
C	59	HIS	TYR	ENGINEERED MUTATION	UNP Q5FZ19
C	89	ALA	THR	ENGINEERED MUTATION	UNP Q5FZ19
C	134	THR	ASN	ENGINEERED MUTATION	UNP Q5FZ19
C	147	LYS	ARG	ENGINEERED MUTATION	UNP Q5FZ19
C	192	GLN	LYS	ENGINEERED MUTATION	UNP Q5FZ19
C	220	LYS	GLU	ENGINEERED MUTATION	UNP Q5FZ19
C	245	LEU	HIS	ENGINEERED MUTATION	UNP Q5FZ19
C	253	ILE	MET	ENGINEERED MUTATION	UNP Q5FZ19
C	291	LYS	ARG	ENGINEERED MUTATION	UNP Q5FZ19
C	296	VAL	ALA	ENGINEERED MUTATION	UNP Q5FZ19
C	301	SER	ALA	ENGINEERED MUTATION	UNP Q5FZ19
C	303	ARG	LYS	ENGINEERED MUTATION	UNP Q5FZ19
D	58	LEU	VAL	ENGINEERED MUTATION	UNP Q5FZ19
D	59	HIS	TYR	ENGINEERED MUTATION	UNP Q5FZ19
D	89	ALA	THR	ENGINEERED MUTATION	UNP Q5FZ19
D	134	THR	ASN	ENGINEERED MUTATION	UNP Q5FZ19
D	147	LYS	ARG	ENGINEERED MUTATION	UNP Q5FZ19
D	192	GLN	LYS	ENGINEERED MUTATION	UNP Q5FZ19
D	220	LYS	GLU	ENGINEERED MUTATION	UNP Q5FZ19
D	245	LEU	HIS	ENGINEERED MUTATION	UNP Q5FZ19
D	253	ILE	MET	ENGINEERED MUTATION	UNP Q5FZ19
D	291	LYS	ARG	ENGINEERED MUTATION	UNP Q5FZ19
D	296	VAL	ALA	ENGINEERED MUTATION	UNP Q5FZ19
D	301	SER	ALA	ENGINEERED MUTATION	UNP Q5FZ19
D	303	ARG	LYS	ENGINEERED MUTATION	UNP Q5FZ19

- Molecule 2 is ACETATE ION (three-letter code: ACT) (formula: C₂H₃O₂).



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

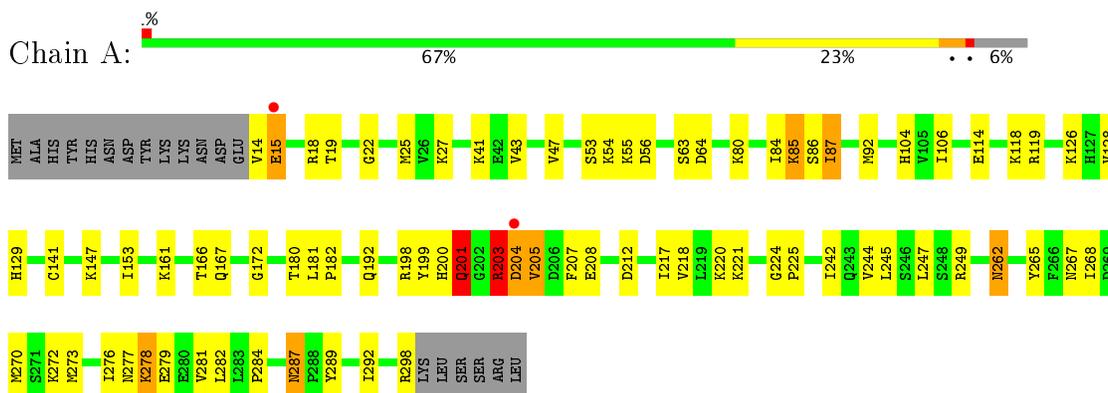
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	57	Total O 57 57	0	0
3	B	60	Total O 60 60	0	0
3	C	72	Total O 72 72	0	0
3	D	59	Total O 59 59	0	0

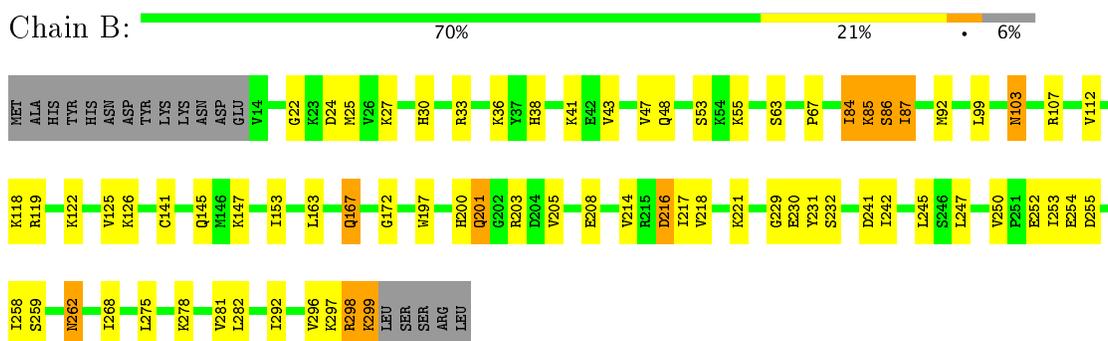
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

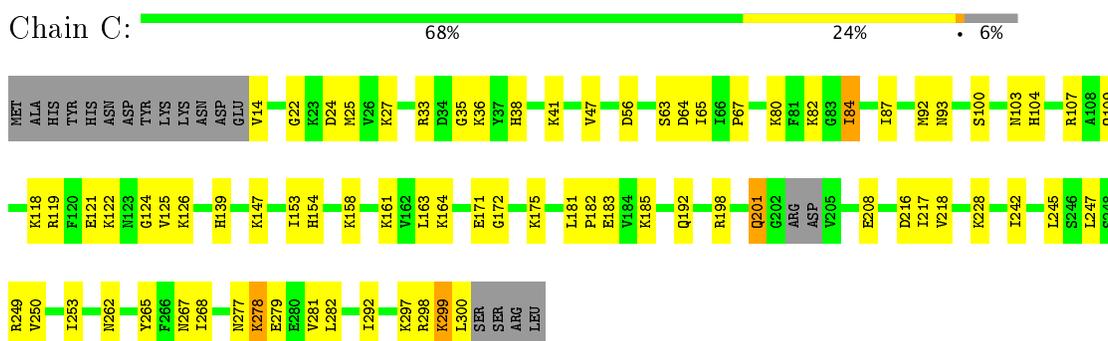
- Molecule 1: Uricase



- Molecule 1: Uricase

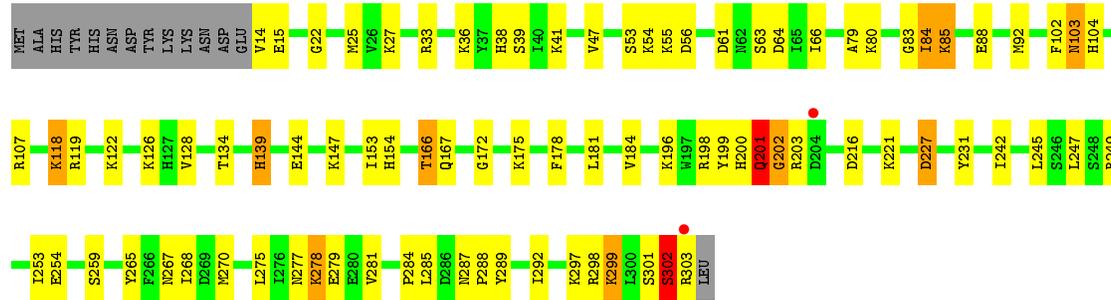


- Molecule 1: Uricase



- Molecule 1: Uricase

Chain D: %



4 Data and refinement statistics

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, α , β , γ	143.78Å 143.78Å 138.87Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	39.77 – 2.40 39.77 – 2.40	Depositor EDS
% Data completeness (in resolution range)	100.0 (39.77-2.40) 97.0 (39.77-2.40)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.11 (at 2.39Å)	Xtrriage
Refinement program	PHENIX dev_1458	Depositor
R, R_{free}	0.174 , 0.220 0.175 , 0.217	Depositor DCC
R_{free} test set	3103 reflections (5.04%)	DCC
Wilson B-factor (Å ²)	44.3	Xtrriage
Anisotropy	0.137	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	(Not available) , (Not available)	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.487 for h,-h-k,-l	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	9529	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.81% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ 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:
ACT

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	0.48	0/2360	0.62	0/3189
1	B	0.46	0/2369	0.64	1/3200 (0.0%)
1	C	0.46	0/2357	0.61	0/3183
1	D	0.46	0/2400	0.61	0/3241
All	All	0.46	0/9486	0.62	1/12813 (0.0%)

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	B	0	1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	B	229	GLY	N-CA-C	5.71	127.37	113.10

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	298	ARG	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	2304	0	2306	89	0
1	B	2313	0	2319	69	0
1	C	2302	0	2312	71	0
1	D	2344	0	2353	85	0
2	B	4	3	3	0	0
2	D	8	3	6	0	0
3	A	57	0	0	7	0
3	B	60	0	0	2	0
3	C	72	0	0	3	0
3	D	59	0	0	4	0
All	All	9523	6	9299	281	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 15.

The worst 5 of 281 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:270:MET:HE3	1:D:277:ASN:HB2	1.42	1.01
1:A:270:MET:HE2	1:A:277:ASN:HB2	1.39	1.00
1:D:41:LYS:HD2	1:D:84:ILE:HD13	1.41	0.99
1:A:270:MET:CE	1:A:277:ASN:HB2	1.99	0.92
1:B:221:LYS:HE2	1:B:241:ASP:HB3	1.51	0.91

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	283/304 (93%)	264 (93%)	14 (5%)	5 (2%)	10	12
1	B	284/304 (93%)	270 (95%)	11 (4%)	3 (1%)	17	23
1	C	281/304 (92%)	267 (95%)	12 (4%)	2 (1%)	25	37
1	D	288/304 (95%)	265 (92%)	18 (6%)	5 (2%)	11	13
All	All	1136/1216 (93%)	1066 (94%)	55 (5%)	15 (1%)	14	19

5 of 15 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	201	GLN
1	A	203	ARG
1	B	201	GLN
1	B	230	GLU
1	C	201	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	259/277 (94%)	249 (96%)	10 (4%)	37	56
1	B	260/277 (94%)	248 (95%)	12 (5%)	31	49
1	C	259/277 (94%)	250 (96%)	9 (4%)	41	61
1	D	264/277 (95%)	248 (94%)	16 (6%)	22	34
All	All	1042/1108 (94%)	995 (96%)	47 (4%)	32	50

5 of 47 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	299	LYS
1	C	201	GLN
1	D	285	LEU
1	C	100	SER

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Mol	Chain	Res	Type
1	C	208	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 21 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	167	GLN
1	B	267	ASN
1	D	76	HIS
1	B	103	ASN
1	D	93	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](#)

3 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	ACT	B	401	-	1,3,3	1.55	0	0,3,3	0.00	-
2	ACT	D	401	-	1,3,3	2.19	1 (100%)	0,3,3	0.00	-
2	ACT	D	402	-	1,3,3	1.49	0	0,3,3	0.00	-

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	ACT	B	401	-	-	0/0/0/0	0/0/0/0
2	ACT	D	401	-	-	0/0/0/0	0/0/0/0
2	ACT	D	402	-	-	0/0/0/0	0/0/0/0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	401	ACT	CH3-C	2.19	1.51	1.48

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 [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	285/304 (93%)	-0.47	2 (0%) 87 86	31, 54, 82, 110	0
1	B	286/304 (94%)	-0.51	0 100 100	31, 52, 82, 102	0
1	C	285/304 (93%)	-0.48	0 100 100	30, 52, 84, 109	0
1	D	290/304 (95%)	-0.47	2 (0%) 87 86	31, 55, 84, 119	0
All	All	1146/1216 (94%)	-0.48	4 (0%) 93 93	30, 53, 84, 119	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	204	ASP	3.3
1	D	303	ARG	2.3
1	A	204	ASP	2.3
1	A	15	GLU	2.2

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 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, 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	LLDF	B-factors(\AA^2)	Q<0.9
2	ACT	D	402	4/4	0.97	0.19	7.39	66,73,79,79	0
2	ACT	B	401	4/4	0.90	0.20	6.34	59,71,79,79	0
2	ACT	D	401	4/4	0.91	0.16	1.60	56,61,75,78	0

6.5 Other polymers [i](#)

There are no such residues in this entry.