wwPDB X-ray Structure Validation Summary Report (i)

Feb 1, 2016 – 09:05 AM GMT

PDB ID	:	3H3F
Title	:	Rabbit muscle L-lactate dehydrogenase in complex with NADH and oxamate
Authors	:	Bujacz, A.; Bujacz, G.; Swiderek, K.; Paneth, P.
Deposited on	:	2009-04-16
Resolution	:	2.38 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

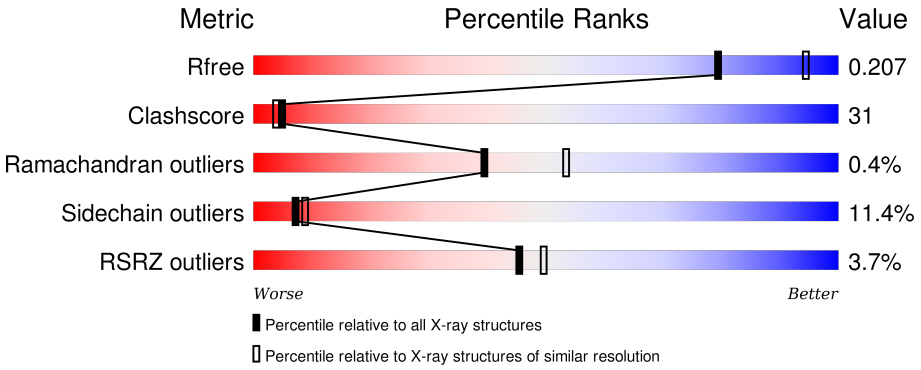
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	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 2.38 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	4019 (2.40-2.36)
Clashscore	102246	4595 (2.40-2.36)
Ramachandran outliers	100387	4520 (2.40-2.36)
Sidechain outliers	100360	4522 (2.40-2.36)
RSRZ outliers	91569	4034 (2.40-2.36)

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	331	4% 58% 34% 8%
1	B	331	2% 66% 30% 5%
1	C	331	3% 56% 36% 8%
1	D	331	3% 57% 37% 6%
1	E	331	7% 50% 42% 8%

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Mol	Chain	Length	Quality of chain
1	F	331	2% 61% 32% 7%
1	G	331	% 62% 31% 7%
1	H	331	7% 52% 40% 8%

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	NAI	D	332	-	-	-	X
3	OXM	A	333	-	-	-	X
3	OXM	D	333	-	-	-	X
3	OXM	E	333	-	-	-	X
3	OXM	F	333	-	-	X	-
4	ACT	A	334	-	-	-	X
4	ACT	B	334	-	-	-	X
4	ACT	C	334	-	-	X	X
4	ACT	D	336	-	-	X	-
4	ACT	E	335	-	-	-	X
4	ACT	E	338	-	-	-	X
4	ACT	F	334	-	-	-	X
4	ACT	H	337	-	-	-	X

2 Entry composition (i)

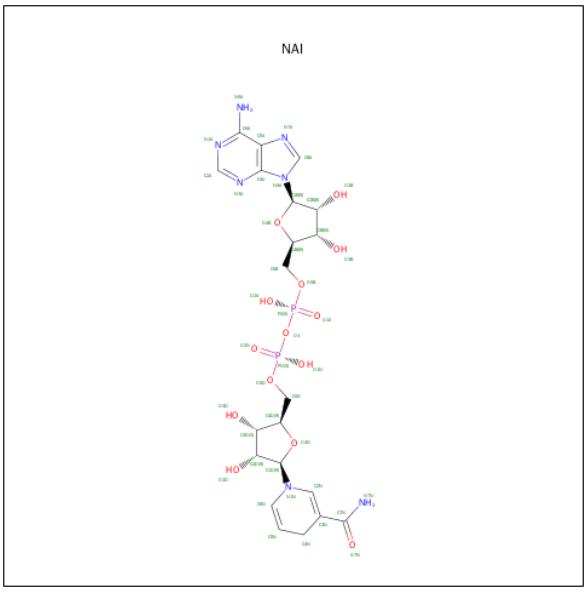
There are 5 unique types of molecules in this entry. The entry contains 22218 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 L-lactate dehydrogenase A chain.

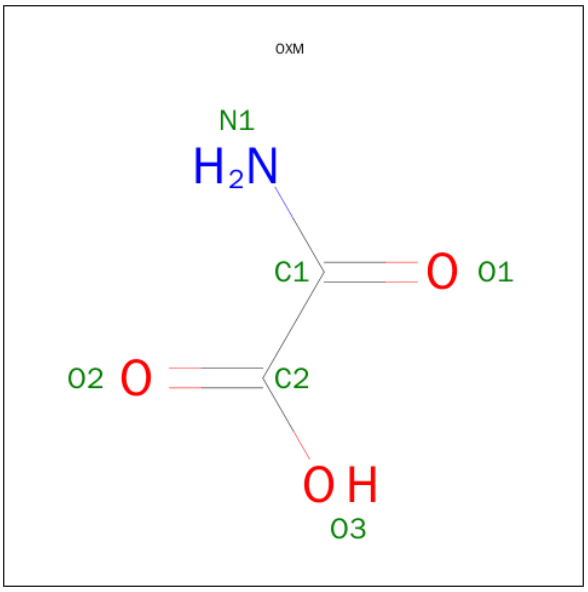
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	331	Total	C	N	O	S	0	1	0
			2562	1635	442	471	14			
1	B	331	Total	C	N	O	S	0	2	0
			2568	1641	442	471	14			
1	C	331	Total	C	N	O	S	0	0	0
			2559	1633	441	471	14			
1	D	331	Total	C	N	O	S	0	4	0
			2575	1644	442	475	14			
1	E	331	Total	C	N	O	S	0	1	0
			2564	1636	441	473	14			
1	F	331	Total	C	N	O	S	0	1	0
			2563	1636	441	471	15			
1	G	331	Total	C	N	O	S	0	3	0
			2571	1639	442	476	14			
1	H	331	Total	C	N	O	S	0	2	0
			2568	1640	441	473	14			

- Molecule 2 is 1,4-DIHYDRONICOTINAMIDE ADENINE DINUCLEOTIDE (three-letter code: NAI) (formula: $C_{21}H_{29}N_7O_{14}P_2$).



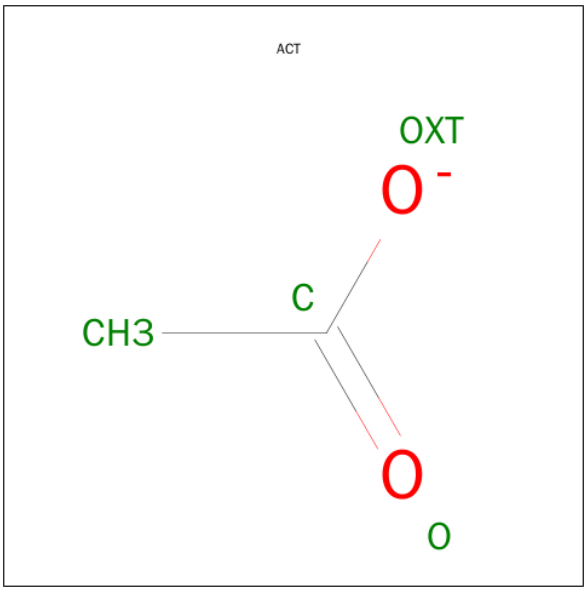
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	B	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	C	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	D	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	E	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	F	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	G	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	H	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

- Molecule 3 is OXAMIC ACID (three-letter code: OXM) (formula: $C_2H_3NO_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			6	2	1	3		
3	B	1	Total	C	N	O	0	0
			6	2	1	3		
3	C	1	Total	C	N	O	0	0
			6	2	1	3		
3	D	1	Total	C	N	O	0	0
			6	2	1	3		
3	E	1	Total	C	N	O	0	0
			6	2	1	3		
3	F	1	Total	C	N	O	0	0
			6	2	1	3		
3	G	1	Total	C	N	O	0	0
			6	2	1	3		
3	H	1	Total	C	N	O	0	0
			6	2	1	3		

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



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O 4 2 2	0	0
4	B	1	Total C O 4 2 2	0	0
4	C	1	Total C O 4 2 2	0	0
4	C	1	Total C O 4 2 2	0	0
4	D	1	Total C O 4 2 2	0	0
4	D	1	Total C O 4 2 2	0	0
4	D	1	Total C O 4 2 2	0	0
4	D	1	Total C O 4 2 2	0	0
4	D	1	Total C O 4 2 2	0	0
4	E	1	Total C O 4 2 2	0	0
4	E	1	Total C O 4 2 2	0	0
4	E	1	Total C O 4 2 2	0	0
4	E	1	Total C O 4 2 2	0	0
4	E	1	Total C O 4 2 2	0	0

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	F	1	Total C O 4 2 2	0	0
4	F	1	Total C O 4 2 2	0	0
4	G	1	Total C O 4 2 2	0	0
4	G	1	Total C O 4 2 2	0	0
4	G	1	Total C O 4 2 2	0	0
4	H	1	Total C O 4 2 2	0	0
4	H	1	Total C O 4 2 2	0	0
4	H	1	Total C O 4 2 2	0	0
4	H	1	Total C O 4 2 2	0	0

- Molecule 5 is water.

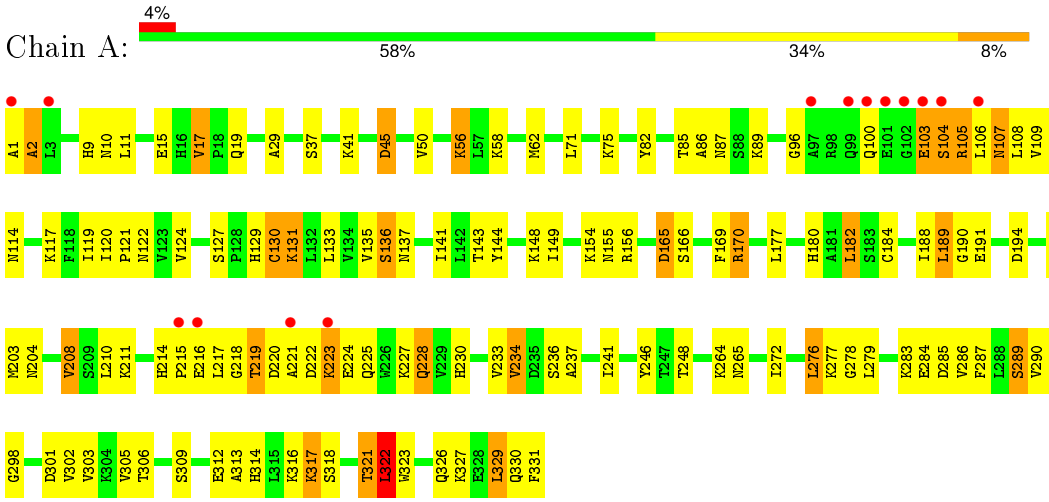
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	146	Total O 146 146	0	0
5	B	202	Total O 202 202	0	0
5	C	143	Total O 143 143	0	0
5	D	139	Total O 139 139	0	0
5	E	116	Total O 116 116	0	0
5	F	153	Total O 153 153	0	0
5	G	158	Total O 158 158	0	0
5	H	139	Total O 139 139	0	0

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.

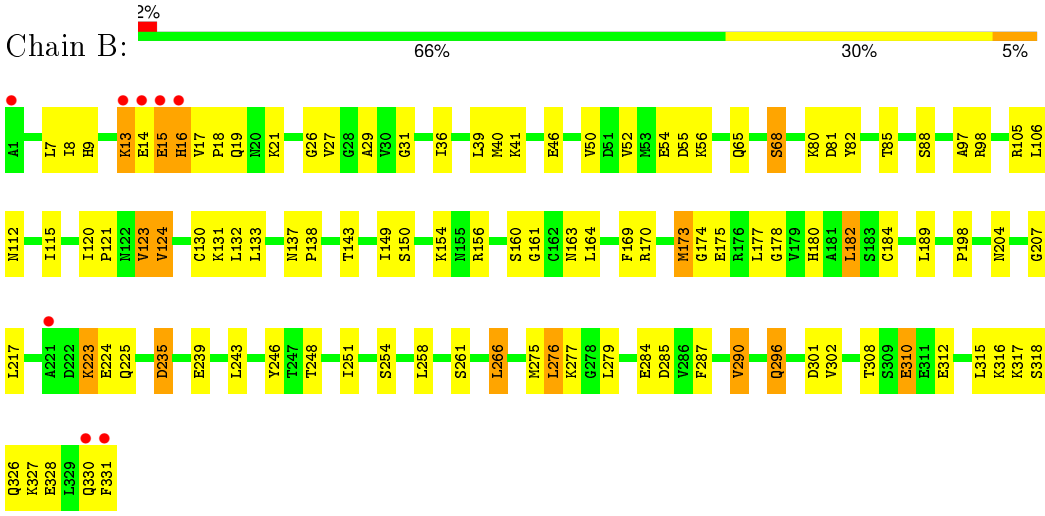
- Molecule 1: L-lactate dehydrogenase A chain

Chain A:



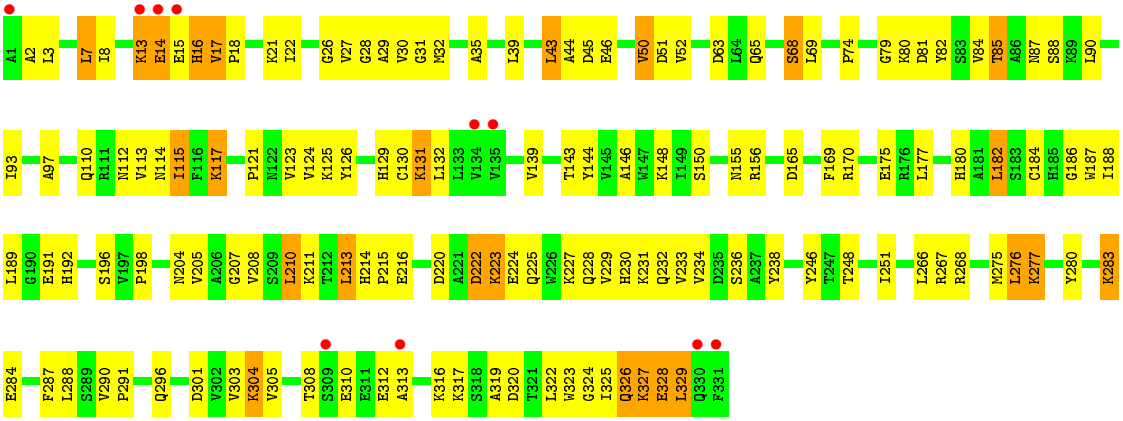
- Molecule 1: L-lactate dehydrogenase A chain

Chain B:



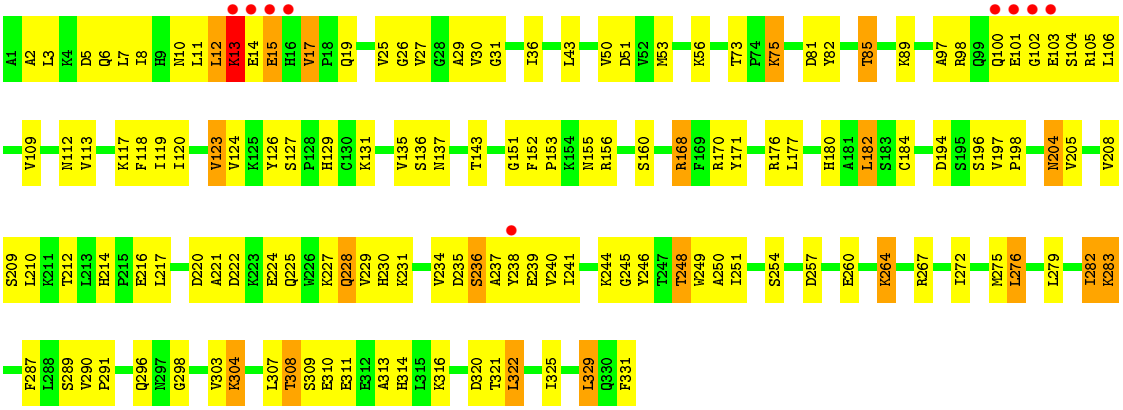
- Molecule 1: L-lactate dehydrogenase A chain

Chain C:



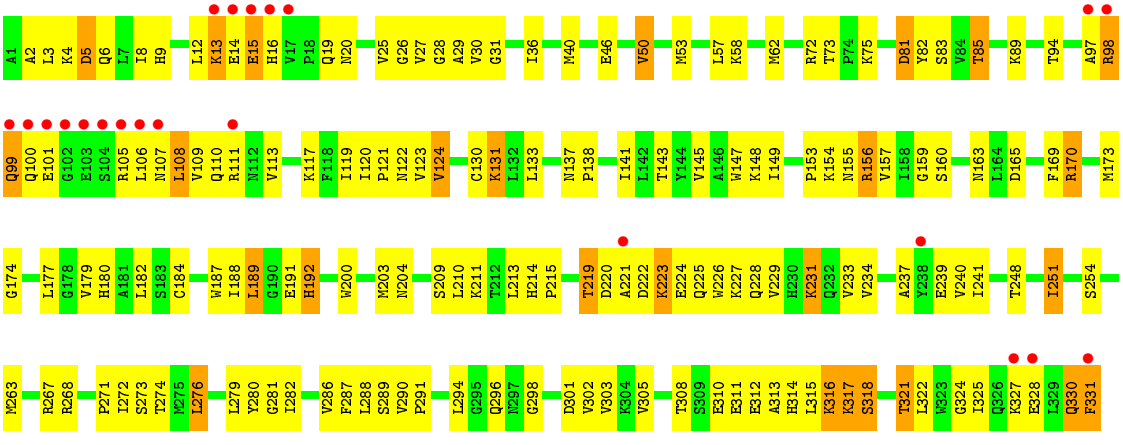
• Molecule 1: L-lactate dehydrogenase A chain

Chain D:



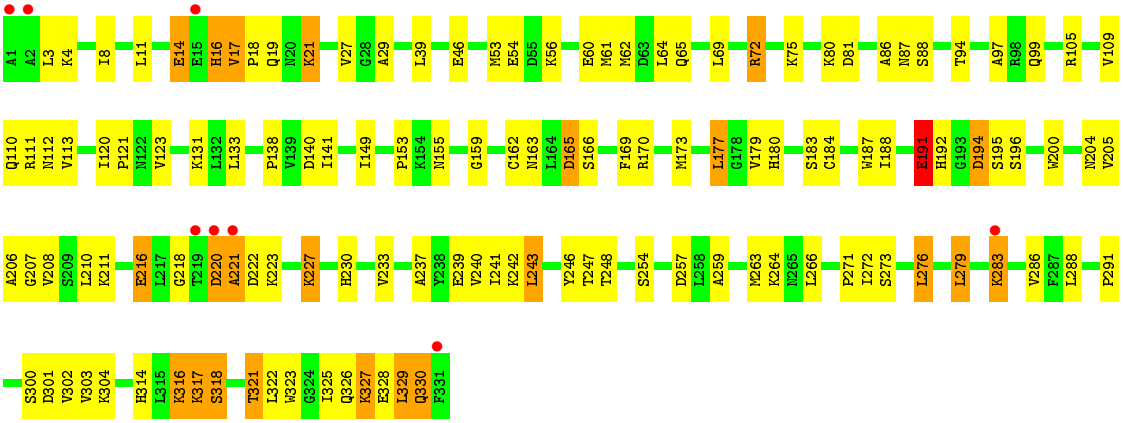
• Molecule 1: L-lactate dehydrogenase A chain

Chain E:



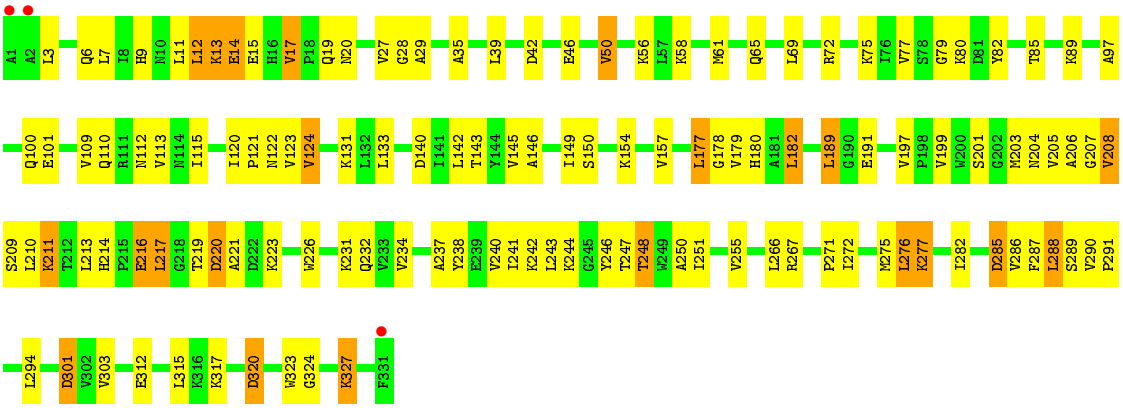
• Molecule 1: L-lactate dehydrogenase A chain

Chain F:



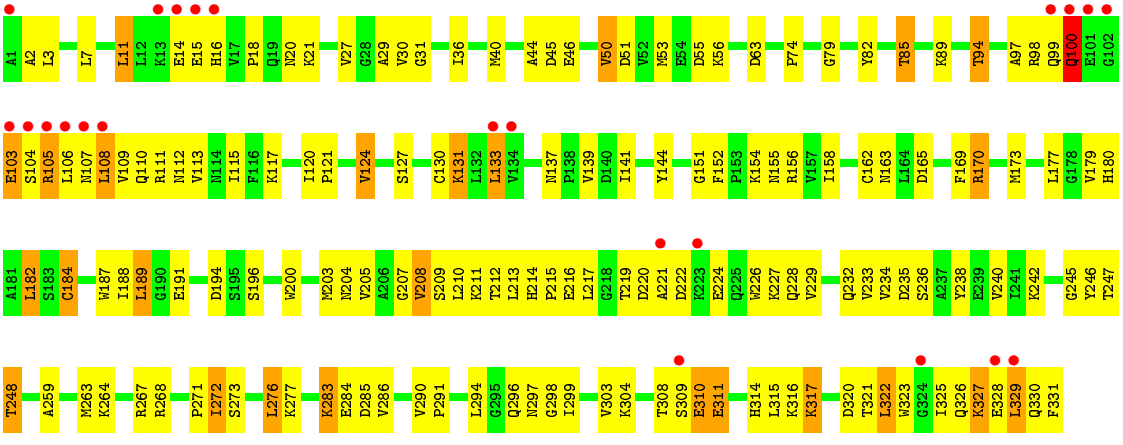
• Molecule 1: L-lactate dehydrogenase A chain

Chain G:



• Molecule 1: L-lactate dehydrogenase A chain

Chain H:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α, β, γ	65.50Å 85.28Å 138.53Å 98.49° 91.67° 111.59°	Depositor
Resolution (Å)	60.00 – 2.38 30.96 – 2.38	Depositor EDS
% Data completeness (in resolution range)	89.7 (60.00-2.38) 86.4 (30.96-2.38)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.13 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.5.0066	Depositor
R, R_{free}	0.160 , 0.207 0.164 , 0.207	Depositor DCC
R_{free} test set	4887 reflections (5.21%)	DCC
Wilson B-factor (Å^2)	28.8	Xtriage
Anisotropy	0.092	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.31 , 52.6	EDS
Estimated twinning fraction	0.034 for h,-h-k,-l	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.47$, $< L^2 > = 0.30$	Xtriage
Outliers	0 of 98703 reflections	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	22218	wwPDB-VP
Average B, all atoms (Å^2)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.46% 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: OXM, NAI, 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	0.89	0/2614	0.96	5/3534 (0.1%)
1	B	0.92	0/2622	1.00	3/3545 (0.1%)
1	C	0.88	0/2605	0.96	3/3523 (0.1%)
1	D	0.87	1/2639 (0.0%)	0.97	5/3568 (0.1%)
1	E	0.86	0/2613	0.96	2/3534 (0.1%)
1	F	0.91	1/2613 (0.0%)	0.96	0/3533
1	G	0.88	0/2632	0.97	3/3559 (0.1%)
1	H	0.83	1/2622 (0.0%)	0.95	4/3546 (0.1%)
All	All	0.88	3/20960 (0.0%)	0.97	25/28342 (0.1%)

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	D	0	2

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	H	184	CYS	CB-SG	-5.16	1.73	1.81
1	F	191	GLU	CG-CD	5.16	1.59	1.51
1	D	126	TYR	CB-CG	-5.04	1.44	1.51

The worst 5 of 25 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	268	ARG	NE-CZ-NH2	-8.94	115.83	120.30
1	B	235	ASP	CB-CG-OD1	6.82	124.44	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	G	42	ASP	CB-CG-OD1	6.26	123.94	118.30
1	H	170	ARG	NE-CZ-NH2	-6.17	117.21	120.30
1	E	156	ARG	NE-CZ-NH1	5.87	123.23	120.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	13	LYS	Peptide
1	D	282	ILE	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	2562	0	2644	199	0
1	B	2568	0	2655	124	0
1	C	2559	0	2639	180	0
1	D	2575	0	2653	175	0
1	E	2564	0	2643	204	0
1	F	2563	0	2644	161	0
1	G	2571	0	2647	146	0
1	H	2568	0	2648	200	0
2	A	44	0	27	4	0
2	B	44	0	27	2	0
2	C	44	0	27	4	0
2	D	44	0	27	13	0
2	E	44	0	27	10	0
2	F	44	0	27	4	0
2	G	44	0	27	2	0
2	H	44	0	27	1	0
3	A	6	0	2	1	0
3	B	6	0	2	0	0
3	C	6	0	2	1	0
3	D	6	0	2	0	0
3	E	6	0	2	1	0
3	F	6	0	2	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	G	6	0	2	1	0
3	H	6	0	2	0	0
4	A	4	0	3	0	0
4	B	4	0	3	0	0
4	C	8	0	6	2	0
4	D	20	0	15	2	0
4	E	20	0	15	1	0
4	F	8	0	6	1	0
4	G	12	0	9	0	0
4	H	16	0	12	0	0
5	A	146	0	0	24	0
5	B	202	0	0	18	0
5	C	143	0	0	13	0
5	D	139	0	0	21	0
5	E	116	0	0	30	0
5	F	153	0	0	16	0
5	G	158	0	0	23	0
5	H	139	0	0	21	0
All	All	22218	0	21474	1332	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 31.

The worst 5 of 1332 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:C:223:LYS:CD	1:C:223:LYS:H	1.25	1.45
1:E:317:LYS:CB	1:E:317:LYS:HZ2	0.88	1.41
1:A:106:LEU:HD23	1:A:106:LEU:O	1.26	1.28
1:H:308:THR:CG2	1:H:310:GLU:HG2	1.62	1.28
1:E:100:GLN:CG	1:E:101:GLU:H	1.42	1.27

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	330/331 (100%)	306 (93%)	22 (7%)	2 (1%)	30	40
1	B	331/331 (100%)	314 (95%)	17 (5%)	0	100	100
1	C	329/331 (99%)	307 (93%)	22 (7%)	0	100	100
1	D	333/331 (101%)	307 (92%)	25 (8%)	1 (0%)	46	61
1	E	330/331 (100%)	307 (93%)	21 (6%)	2 (1%)	30	40
1	F	330/331 (100%)	303 (92%)	25 (8%)	2 (1%)	30	40
1	G	332/331 (100%)	311 (94%)	20 (6%)	1 (0%)	46	61
1	H	331/331 (100%)	306 (92%)	23 (7%)	2 (1%)	30	40
All	All	2646/2648 (100%)	2461 (93%)	175 (7%)	10 (0%)	39	53

5 of 10 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	2	ALA
1	E	192	HIS
1	D	221	ALA
1	E	312	GLU
1	A	104	SER

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	285/284 (100%)	252 (88%)	33 (12%)	7	8
1	B	286/284 (101%)	261 (91%)	25 (9%)	13	17
1	C	284/284 (100%)	252 (89%)	32 (11%)	7	9
1	D	288/284 (101%)	254 (88%)	34 (12%)	6	8
1	E	285/284 (100%)	249 (87%)	36 (13%)	5	6
1	F	285/284 (100%)	255 (90%)	30 (10%)	8	11

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	G	287/284 (101%)	254 (88%)	33 (12%)	7 8
1	H	286/284 (101%)	246 (86%)	40 (14%)	4 5
All	All	2286/2272 (101%)	2023 (88%)	263 (12%)	7 8

5 of 263 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	290	VAL
1	E	223	LYS
1	H	208	VAL
1	D	322	LEU
1	E	99	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 55 such sidechains are listed below:

Mol	Chain	Res	Type
1	D	326	GLN
1	E	204	ASN
1	H	16	HIS
1	E	6	GLN
1	E	107	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)

39 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	NAI	A	332	-	38,48,48	1.30	5 (13%)	48,73,73	2.21	12 (25%)
3	OXM	A	333	-	2,5,5	1.05	0	2,6,6	1.85	1 (50%)
4	ACT	A	334	-	1,3,3	3.79	1 (100%)	0,3,3	0.00	-
2	NAI	B	332	-	38,48,48	1.48	5 (13%)	48,73,73	2.84	13 (27%)
3	OXM	B	333	-	2,5,5	1.18	0	2,6,6	0.18	0
4	ACT	B	334	-	1,3,3	3.49	1 (100%)	0,3,3	0.00	-
2	NAI	C	332	-	38,48,48	1.35	5 (13%)	48,73,73	2.51	10 (20%)
3	OXM	C	333	-	2,5,5	0.60	0	2,6,6	0.69	0
4	ACT	C	334	-	1,3,3	4.01	1 (100%)	0,3,3	0.00	-
4	ACT	C	335	-	1,3,3	2.36	1 (100%)	0,3,3	0.00	-
2	NAI	D	332	-	38,48,48	1.68	7 (18%)	48,73,73	2.23	11 (22%)
3	OXM	D	333	-	2,5,5	0.24	0	2,6,6	0.68	0
4	ACT	D	334	-	1,3,3	2.09	1 (100%)	0,3,3	0.00	-
4	ACT	D	335	-	1,3,3	2.93	1 (100%)	0,3,3	0.00	-
4	ACT	D	336	-	1,3,3	1.70	0	0,3,3	0.00	-
4	ACT	D	337	-	1,3,3	1.49	0	0,3,3	0.00	-
4	ACT	D	338	-	1,3,3	1.56	0	0,3,3	0.00	-
2	NAI	E	332	-	38,48,48	1.43	5 (13%)	48,73,73	2.20	8 (16%)
3	OXM	E	333	-	2,5,5	0.89	0	2,6,6	1.15	0
4	ACT	E	334	-	1,3,3	1.78	0	0,3,3	0.00	-
4	ACT	E	335	-	1,3,3	3.30	1 (100%)	0,3,3	0.00	-
4	ACT	E	336	-	1,3,3	2.67	1 (100%)	0,3,3	0.00	-
4	ACT	E	337	-	1,3,3	2.95	1 (100%)	0,3,3	0.00	-
4	ACT	E	338	-	1,3,3	2.88	1 (100%)	0,3,3	0.00	-
2	NAI	F	332	-	38,48,48	1.42	4 (10%)	48,73,73	2.59	14 (29%)
3	OXM	F	333	-	2,5,5	0.58	0	2,6,6	2.07	1 (50%)
4	ACT	F	334	-	1,3,3	1.79	0	0,3,3	0.00	-
4	ACT	F	335	-	1,3,3	2.07	1 (100%)	0,3,3	0.00	-
2	NAI	G	332	-	38,48,48	1.18	3 (7%)	48,73,73	2.59	16 (33%)
3	OXM	G	333	-	2,5,5	0.13	0	2,6,6	2.10	1 (50%)
4	ACT	G	334	-	1,3,3	1.32	0	0,3,3	0.00	-

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
4	ACT	G	335	-	1,3,3	2.21	1 (100%)	0,3,3	0.00	-
4	ACT	G	336	-	1,3,3	3.15	1 (100%)	0,3,3	0.00	-
2	NAI	H	332	-	38,48,48	1.43	5 (13%)	48,73,73	2.72	14 (29%)
3	OXM	H	333	-	2,5,5	1.04	0	2,6,6	2.21	1 (50%)
4	ACT	H	334	-	1,3,3	2.09	1 (100%)	0,3,3	0.00	-
4	ACT	H	335	-	1,3,3	2.08	1 (100%)	0,3,3	0.00	-
4	ACT	H	336	-	1,3,3	2.21	1 (100%)	0,3,3	0.00	-
4	ACT	H	337	-	1,3,3	1.34	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	NAI	A	332	-	-	0/25/72/72	0/5/5/5
3	OXM	A	333	-	-	0/0/4/4	0/0/0/0
4	ACT	A	334	-	-	0/0/0/0	0/0/0/0
2	NAI	B	332	-	-	0/25/72/72	0/5/5/5
3	OXM	B	333	-	-	0/0/4/4	0/0/0/0
4	ACT	B	334	-	-	0/0/0/0	0/0/0/0
2	NAI	C	332	-	-	0/25/72/72	0/5/5/5
3	OXM	C	333	-	-	0/0/4/4	0/0/0/0
4	ACT	C	334	-	-	0/0/0/0	0/0/0/0
4	ACT	C	335	-	-	0/0/0/0	0/0/0/0
2	NAI	D	332	-	-	0/25/72/72	0/5/5/5
3	OXM	D	333	-	-	0/0/4/4	0/0/0/0
4	ACT	D	334	-	-	0/0/0/0	0/0/0/0
4	ACT	D	335	-	-	0/0/0/0	0/0/0/0
4	ACT	D	336	-	-	0/0/0/0	0/0/0/0
4	ACT	D	337	-	-	0/0/0/0	0/0/0/0
4	ACT	D	338	-	-	0/0/0/0	0/0/0/0
2	NAI	E	332	-	-	0/25/72/72	0/5/5/5
3	OXM	E	333	-	-	0/0/4/4	0/0/0/0
4	ACT	E	334	-	-	0/0/0/0	0/0/0/0
4	ACT	E	335	-	-	0/0/0/0	0/0/0/0
4	ACT	E	336	-	-	0/0/0/0	0/0/0/0
4	ACT	E	337	-	-	0/0/0/0	0/0/0/0
4	ACT	E	338	-	-	0/0/0/0	0/0/0/0
2	NAI	F	332	-	-	0/25/72/72	0/5/5/5
3	OXM	F	333	-	-	0/0/4/4	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	ACT	F	334	-	-	0/0/0/0	0/0/0/0
4	ACT	F	335	-	-	0/0/0/0	0/0/0/0
2	NAI	G	332	-	-	0/25/72/72	0/5/5/5
3	OXM	G	333	-	-	0/0/4/4	0/0/0/0
4	ACT	G	334	-	-	0/0/0/0	0/0/0/0
4	ACT	G	335	-	-	0/0/0/0	0/0/0/0
4	ACT	G	336	-	-	0/0/0/0	0/0/0/0
2	NAI	H	332	-	-	0/25/72/72	0/5/5/5
3	OXM	H	333	-	-	0/0/4/4	0/0/0/0
4	ACT	H	334	-	-	0/0/0/0	0/0/0/0
4	ACT	H	335	-	-	0/0/0/0	0/0/0/0
4	ACT	H	336	-	-	0/0/0/0	0/0/0/0
4	ACT	H	337	-	-	0/0/0/0	0/0/0/0

The worst 5 of 55 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	332	NAI	C4N-C5N	-5.09	1.38	1.49
2	C	332	NAI	C4N-C5N	-4.12	1.40	1.49
2	H	332	NAI	C4N-C5N	-4.08	1.40	1.49
2	B	332	NAI	C4N-C5N	-3.93	1.40	1.49
2	E	332	NAI	C4N-C5N	-3.64	1.41	1.49

The worst 5 of 102 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	332	NAI	N3A-C2A-N1A	-13.63	118.45	128.89
2	F	332	NAI	N3A-C2A-N1A	-13.11	118.85	128.89
2	C	332	NAI	N3A-C2A-N1A	-11.69	119.94	128.89
2	G	332	NAI	N3A-C2A-N1A	-11.52	120.08	128.89
2	H	332	NAI	N3A-C2A-N1A	-10.85	120.58	128.89

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

17 monomers are involved in 48 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	332	NAI	4	0
3	A	333	OXM	1	0
2	B	332	NAI	2	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	332	NAI	4	0
3	C	333	OXM	1	0
4	C	334	ACT	2	0
2	D	332	NAI	13	0
4	D	336	ACT	2	0
2	E	332	NAI	10	0
3	E	333	OXM	1	0
4	E	334	ACT	1	0
2	F	332	NAI	4	0
3	F	333	OXM	3	0
4	F	335	ACT	1	0
2	G	332	NAI	2	0
3	G	333	OXM	1	0
2	H	332	NAI	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^{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	331/331 (100%)	-0.11	14 (4%)	40	45	8, 18, 34, 52	0
1	B	331/331 (100%)	-0.27	8 (2%)	62	65	7, 16, 29, 43	0
1	C	331/331 (100%)	-0.14	10 (3%)	54	57	9, 18, 33, 46	0
1	D	331/331 (100%)	-0.18	9 (2%)	58	61	8, 19, 33, 46	0
1	E	331/331 (100%)	-0.00	22 (6%)	22	25	7, 20, 38, 53	0
1	F	331/331 (100%)	-0.24	8 (2%)	62	65	8, 17, 33, 43	0
1	G	331/331 (100%)	-0.25	3 (0%)	85	87	8, 19, 29, 43	0
1	H	331/331 (100%)	0.04	23 (6%)	20	22	10, 20, 35, 53	0
All	All	2648/2648 (100%)	-0.14	97 (3%)	45	49	7, 19, 33, 53	0

The worst 5 of 97 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	100	GLN	6.1
1	A	102	GLY	5.8
1	H	102	GLY	5.0
1	E	99	GLN	4.9
1	H	106	LEU	4.8

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 (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^{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
4	ACT	H	337	4/4	0.95	0.34	19.05	46,46,46,47	0
4	ACT	C	334	4/4	0.95	0.25	9.29	27,28,28,28	0
4	ACT	E	338	4/4	0.92	0.25	8.56	39,40,40,40	0
4	ACT	E	335	4/4	0.86	0.17	5.66	32,32,33,33	0
2	NAI	D	332	44/44	0.85	0.27	5.42	25,32,44,45	0
3	OXM	D	333	6/6	0.72	0.36	5.31	60,62,63,64	0
3	OXM	E	333	6/6	0.90	0.27	5.10	34,36,37,37	0
4	ACT	A	334	4/4	0.87	0.16	4.32	31,32,33,33	0
4	ACT	B	334	4/4	0.89	0.16	4.32	30,30,32,33	0
3	OXM	A	333	6/6	0.86	0.23	4.30	34,34,35,36	0
4	ACT	F	334	4/4	0.87	0.13	3.28	38,38,40,40	0
4	ACT	D	336	4/4	0.93	0.28	1.95	39,40,40,40	0
4	ACT	E	334	4/4	0.94	0.15	1.87	37,38,38,38	0
3	OXM	H	333	6/6	0.88	0.17	1.76	11,16,20,22	0
4	ACT	D	338	4/4	0.94	0.11	1.25	32,34,34,34	0
4	ACT	G	335	4/4	0.96	0.20	0.53	39,39,39,40	0
4	ACT	E	336	4/4	0.90	0.18	0.25	41,42,43,43	0
2	NAI	E	332	44/44	0.93	0.15	0.23	20,29,33,36	0
4	ACT	H	335	4/4	0.96	0.10	0.08	33,33,34,34	0
4	ACT	G	334	4/4	0.97	0.10	0.03	30,31,31,33	0
2	NAI	A	332	44/44	0.95	0.12	-0.29	20,25,29,31	0
3	OXM	B	333	6/6	0.98	0.10	-0.46	9,14,16,17	0
2	NAI	B	332	44/44	0.98	0.10	-0.46	7,14,17,18	0
3	OXM	C	333	6/6	0.97	0.10	-0.57	12,15,17,17	0
2	NAI	C	332	44/44	0.99	0.11	-0.65	6,12,15,17	0
2	NAI	F	332	44/44	0.98	0.09	-0.70	6,13,17,19	0
2	NAI	G	332	44/44	0.98	0.10	-0.71	6,13,18,22	0
2	NAI	H	332	44/44	0.98	0.07	-1.23	10,19,23,27	0
3	OXM	G	333	6/6	0.99	0.09	-1.24	5,6,7,9	0
3	OXM	F	333	6/6	0.98	0.07	-1.43	14,16,17,18	0
4	ACT	H	334	4/4	0.97	0.07	-1.59	31,32,32,32	0
4	ACT	C	335	4/4	0.90	0.18	-	42,43,43,43	0
4	ACT	D	334	4/4	0.90	0.40	-	40,41,41,41	0
4	ACT	E	337	4/4	0.82	0.24	-	37,38,38,38	0
4	ACT	D	337	4/4	0.94	0.32	-	32,33,33,34	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å^2)	Q<0.9
4	ACT	F	335	4/4	0.95	0.19	-	37,38,38,39	0
4	ACT	H	336	4/4	0.76	0.37	-	45,46,46,47	0
4	ACT	D	335	4/4	0.92	0.18	-	41,41,42,42	0
4	ACT	G	336	4/4	0.88	0.22	-	42,44,44,44	0

6.5 Other polymers (i)

There are no such residues in this entry.