Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 08:44 AM GMT

PDB ID	:	3FVD
Title	:	Crystal structure of a member of enolase superfamily from ROSEOVARIUS NUBINHIBENS ISM complexed with magnesium
Authors	:	Malashkevich, V.N.; Rutter, M.; Bain, K.T.; Lau, C.; Ozyurt, S.; Smith, D.; Wasserman, S.; Sauder, J.M.; Burley, S.K.; Almo, S.C.; New York SGX Research Center for Structural Genomics (NYSGXRC)
Deposited on	:	2009-01-15
Resolution	:	2.30 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 ⓘ symbol.

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

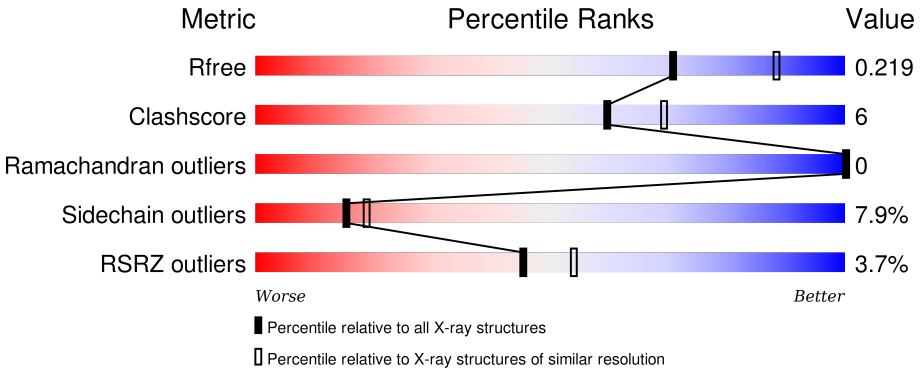
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	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 2.30 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	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (2.30-2.30)

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	378	3% 79% 15%
1	B	378	4% 82% 12%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 5859 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 Mandelate racemase/muconate lactonizing enzyme.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	367	Total	C	N	O	S	0	3	0
			2783	1751	505	516	11			
1	A	361	Total	C	N	O	S	0	4	0
			2732	1712	500	509	11			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-1	MET	-	EXPRESSION TAG	UNP A3SNF7
B	0	SER	-	EXPRESSION TAG	UNP A3SNF7
B	1	LEU	-	EXPRESSION TAG	UNP A3SNF7
B	369	GLU	-	EXPRESSION TAG	UNP A3SNF7
B	370	GLY	-	EXPRESSION TAG	UNP A3SNF7
B	371	HIS	-	EXPRESSION TAG	UNP A3SNF7
B	372	HIS	-	EXPRESSION TAG	UNP A3SNF7
B	373	HIS	-	EXPRESSION TAG	UNP A3SNF7
B	374	HIS	-	EXPRESSION TAG	UNP A3SNF7
B	375	HIS	-	EXPRESSION TAG	UNP A3SNF7
B	376	HIS	-	EXPRESSION TAG	UNP A3SNF7
A	-1	MET	-	EXPRESSION TAG	UNP A3SNF7
A	0	SER	-	EXPRESSION TAG	UNP A3SNF7
A	1	LEU	-	EXPRESSION TAG	UNP A3SNF7
A	369	GLU	-	EXPRESSION TAG	UNP A3SNF7
A	370	GLY	-	EXPRESSION TAG	UNP A3SNF7
A	371	HIS	-	EXPRESSION TAG	UNP A3SNF7
A	372	HIS	-	EXPRESSION TAG	UNP A3SNF7
A	373	HIS	-	EXPRESSION TAG	UNP A3SNF7
A	374	HIS	-	EXPRESSION TAG	UNP A3SNF7
A	375	HIS	-	EXPRESSION TAG	UNP A3SNF7
A	376	HIS	-	EXPRESSION TAG	UNP A3SNF7

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total Mg 1 1	0	0
2	A	1	Total Mg 1 1	0	0

- Molecule 3 is water.

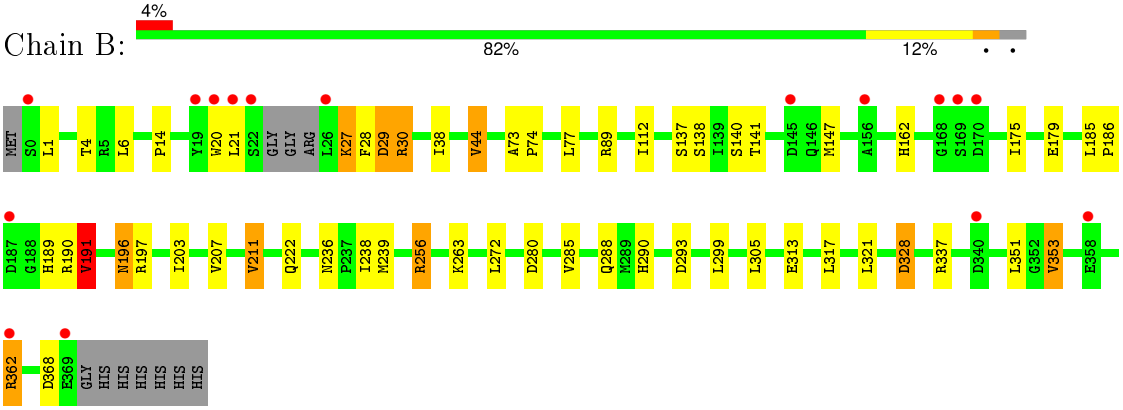
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	342	Total O 342 342	0	0

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.

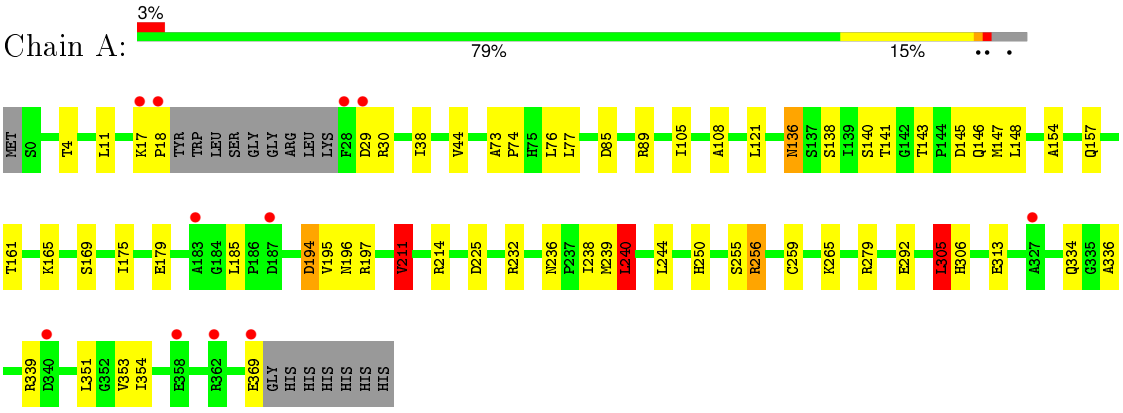
- Molecule 1: Mandelate racemase/muconate lactonizing enzyme

Chain B:



- Molecule 1: Mandelate racemase/muconate lactonizing enzyme

Chain A:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 4 21 2	Depositor
Cell constants a, b, c, α, β, γ	126.52Å 126.52Å 97.29Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.99 – 2.30 19.99 – 2.30	Depositor EDS
% Data completeness (in resolution range)	100.0 (19.99-2.30) 100.0 (19.99-2.30)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	0.17	Depositor
$< I/\sigma(I) >$ [1]	4.69 (at 2.30Å)	Xtriage
Refinement program	REFMAC 5.5.0070	Depositor
R, R_{free}	0.164 , 0.225 0.161 , 0.219	Depositor DCC
R_{free} test set	1789 reflections (5.29%)	DCC
Wilson B-factor (Å^2)	30.9	Xtriage
Anisotropy	0.116	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 46.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Outliers	0 of 35576 reflections	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	5859	wwPDB-VP
Average B, all atoms (Å^2)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.43% 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: MG

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	0.96	1/2813 (0.0%)	0.93	6/3843 (0.2%)
1	B	0.92	3/2864 (0.1%)	0.94	9/3914 (0.2%)
All	All	0.94	4/5677 (0.1%)	0.94	15/7757 (0.2%)

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

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	211	VAL	CB-CG2	-6.56	1.39	1.52
1	B	211	VAL	CB-CG2	-5.25	1.41	1.52
1	B	179	GLU	CG-CD	5.12	1.59	1.51
1	B	285	VAL	CB-CG1	5.04	1.63	1.52

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	239	MET	CA-CB-CG	7.08	125.34	113.30
1	B	191	VAL	CB-CA-C	-6.61	98.85	111.40
1	B	256	ARG	NE-CZ-NH1	-6.58	117.01	120.30
1	B	190	ARG	NE-CZ-NH2	-6.43	117.09	120.30
1	B	280	ASP	CB-CG-OD1	6.32	123.99	118.30
1	A	89	ARG	NE-CZ-NH1	5.99	123.29	120.30
1	B	256	ARG	NE-CZ-NH2	5.92	123.26	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	239	MET	CA-CB-CG	5.84	123.23	113.30
1	B	293	ASP	CB-CG-OD1	5.69	123.42	118.30
1	A	89	ARG	NE-CZ-NH2	-5.56	117.52	120.30
1	A	240	LEU	CB-CG-CD1	5.40	120.19	111.00
1	A	305	LEU	CA-CB-CG	-5.26	103.20	115.30
1	B	89	ARG	NE-CZ-NH1	5.20	122.90	120.30
1	A	194	ASP	CB-CG-OD1	5.14	122.93	118.30
1	B	197	ARG	NE-CZ-NH2	-5.12	117.74	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	28	PHE	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	2732	0	2690	35	0
1	B	2783	0	2742	27	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	342	0	0	5	0
All	All	5859	0	5432	62	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 6.

All (62) 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:20:TRP:CD1	1:B:27:LYS:HD2	2.00	0.95
1:A:165:LYS:NZ	1:A:194:ASP:OD2	1.98	0.95
1:B:20:TRP:HD1	1:B:27:LYS:HD2	1.36	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:334:GLN:HE22	1:A:354:ILE:H	1.33	0.77
1:B:141:THR:HA	1:B:147:MET:CE	2.17	0.74
1:B:299:LEU:HD11	1:B:351:LEU:HD21	1.73	0.69
1:A:18:PRO:HB3	1:A:29:ASP:HB2	1.73	0.69
1:A:175:ILE:O	1:A:179:GLU:HG2	1.94	0.68
1:A:334:GLN:NE2	1:A:354:ILE:H	1.91	0.68
1:A:141:THR:HA	1:A:147:MET:HE3	1.76	0.67
1:B:256:ARG:NH2	3:A:714:HOH:O	2.24	0.66
1:B:185:LEU:HD11	1:B:191:VAL:HG22	1.78	0.65
1:B:141:THR:HA	1:B:147:MET:HE3	1.78	0.65
1:B:299:LEU:CD1	1:B:351:LEU:HD21	2.26	0.65
1:A:256[A]:ARG:HH11	1:A:256[A]:ARG:HG2	1.63	0.64
1:A:141:THR:HB	1:A:165:LYS:O	1.98	0.63
1:B:196:ASN:HA	3:A:675:HOH:O	1.99	0.61
1:B:141:THR:HA	1:B:147:MET:HE2	1.84	0.58
1:B:175:ILE:HG13	1:B:207:VAL:HB	1.85	0.58
1:A:141:THR:HA	1:A:147:MET:CE	2.33	0.57
1:A:225:ASP:OD2	1:A:256[A]:ARG:NH2	2.37	0.57
1:A:18:PRO:HB3	1:A:29:ASP:CB	2.35	0.56
1:A:256[A]:ARG:NH1	1:A:256[A]:ARG:HG2	2.21	0.55
1:B:44:VAL:HG22	1:B:112:ILE:HG23	1.88	0.54
1:B:20:TRP:HD1	1:B:27:LYS:CD	2.16	0.54
1:A:143:THR:HG23	1:A:146:GLN:OE1	2.08	0.54
1:B:186:PRO:HB2	1:B:189:HIS:CD2	2.44	0.52
1:B:299:LEU:HD11	1:B:351:LEU:CD2	2.37	0.52
1:B:288:GLN:HB3	1:B:317:LEU:HD11	1.91	0.52
1:A:244:LEU:HB3	1:A:250:HIS:CE1	2.46	0.51
1:B:29:ASP:HB3	1:B:30:ARG:HD2	1.93	0.50
1:A:136:ASN:HB2	1:A:161:THR:O	2.11	0.50
1:B:196:ASN:ND2	3:A:384:HOH:O	2.44	0.50
1:B:299:LEU:HD13	1:B:353:VAL:HG22	1.94	0.50
1:B:203:ILE:O	1:B:207:VAL:HG13	2.12	0.49
1:A:154:ALA:HA	1:A:157:GLN:NE2	2.28	0.49
1:A:11:LEU:HD12	1:A:11:LEU:N	2.28	0.49
1:A:279:ARG:C	1:A:279:ARG:HD2	2.33	0.49
1:A:121:LEU:HB2	1:A:306:HIS:CG	2.48	0.49
1:A:196:ASN:HD21	1:A:197:ARG:HH21	1.61	0.48
1:A:76:LEU:HD11	1:A:105:ILE:HD13	1.94	0.48
1:A:236:ASN:O	1:A:238:ILE:HD12	2.13	0.48
1:A:256[A]:ARG:HH11	1:A:256[A]:ARG:CG	2.27	0.47
1:A:38:ILE:HD12	1:A:108:ALA:HB3	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:197:ARG:NH1	3:A:532:HOH:O	2.44	0.47
1:A:232:ARG:NH2	3:A:632:HOH:O	2.38	0.46
1:B:6:LEU:HD23	1:B:38:ILE:HG12	1.97	0.46
1:A:240:LEU:HD22	1:A:259:CYS:HB3	1.97	0.45
1:A:306:HIS:HE1	1:A:351:LEU:O	2.00	0.45
1:A:211:VAL:O	1:A:214:ARG:NH1	2.51	0.44
1:B:137:SER:OG	1:B:162:HIS:ND1	2.42	0.44
1:A:147:MET:HE3	1:A:165:LYS:O	2.18	0.43
1:A:265:LYS:HA	1:A:292:GLU:O	2.17	0.43
1:A:73:ALA:HB3	1:A:74:PRO:HD3	2.01	0.43
1:B:236:ASN:O	1:B:238[A]:ILE:HD13	2.19	0.42
1:B:73:ALA:HB3	1:B:74:PRO:HD3	2.01	0.42
1:B:263:LYS:HG2	1:B:290:HIS:HB2	2.02	0.42
1:A:305:LEU:HD22	1:A:336:ALA:CB	2.50	0.41
1:B:362:ARG:HH11	1:B:362:ARG:HB3	1.85	0.41
1:A:148:LEU:HA	1:A:148:LEU:HD12	1.84	0.41
1:A:154:ALA:HA	1:A:157:GLN:HE21	1.86	0.40
1:B:14:PRO:HD2	1:B:328:ASP: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	361/378 (96%)	354 (98%)	7 (2%)	0	100	100
1	B	366/378 (97%)	359 (98%)	7 (2%)	0	100	100
All	All	727/756 (96%)	713 (98%)	14 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	285/295 (97%)	261 (92%)	24 (8%)	14	16
1	B	290/295 (98%)	267 (92%)	23 (8%)	15	19
All	All	575/590 (98%)	528 (92%)	47 (8%)	15	17

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	1	LEU
1	B	4	THR
1	B	21	LEU
1	B	27	LYS
1	B	29	ASP
1	B	30	ARG
1	B	44	VAL
1	B	77	LEU
1	B	138	SER
1	B	140	SER
1	B	191	VAL
1	B	196	ASN
1	B	211	VAL
1	B	222	GLN
1	B	272	LEU
1	B	305	LEU
1	B	313	GLU
1	B	321	LEU
1	B	328	ASP
1	B	337	ARG
1	B	353	VAL
1	B	362	ARG
1	B	368	ASP
1	A	4	THR
1	A	17	LYS
1	A	30	ARG
1	A	44	VAL

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Mol	Chain	Res	Type
1	A	77	LEU
1	A	85	ASP
1	A	136	ASN
1	A	138	SER
1	A	140[A]	SER
1	A	140[B]	SER
1	A	145	ASP
1	A	169	SER
1	A	185	LEU
1	A	195	VAL
1	A	211	VAL
1	A	240	LEU
1	A	255	SER
1	A	256[A]	ARG
1	A	256[B]	ARG
1	A	305	LEU
1	A	313	GLU
1	A	339	ARG
1	A	353	VAL
1	A	369	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	10	HIS
1	B	136	ASN
1	B	277	GLN
1	A	10	HIS
1	A	94	GLN
1	A	123	GLN
1	A	157	GLN
1	A	196	ASN
1	A	222	GLN
1	A	277	GLN
1	A	306	HIS
1	A	334	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 2 ligands modelled in this entry, 2 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	361/378 (95%)	-0.29	11 (3%) 54 63	5, 18, 38, 58	0
1	B	367/378 (97%)	-0.22	16 (4%) 38 47	7, 19, 45, 88	0
All	All	728/756 (96%)	-0.26	27 (3%) 45 54	5, 19, 43, 88	0

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	20	TRP	6.3
1	B	21	LEU	5.8
1	B	19	TYR	4.5
1	B	362	ARG	4.4
1	B	26	LEU	4.4
1	B	22	SER	4.1
1	B	169	SER	3.7
1	B	168	GLY	3.6
1	A	362	ARG	3.6
1	A	28	PHE	3.6
1	B	145	ASP	3.1
1	A	18	PRO	3.1
1	B	156	ALA	3.1
1	B	187	ASP	3.1
1	A	369	GLU	3.0
1	A	358	GLU	2.6
1	B	0	SER	2.6
1	A	340	ASP	2.5
1	B	369	GLU	2.5
1	B	358	GLU	2.5
1	A	327	ALA	2.4
1	B	340	ASP	2.3
1	A	187	ASP	2.3
1	A	29	ASP	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	170	ASP	2.1
1	A	17	LYS	2.1
1	A	183	ALA	2.0

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, 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	MG	A	501	1/1	0.87	0.17	1.76	17,17,17,17	1
2	MG	B	501	1/1	0.90	0.11	-0.05	16,16,16,16	1

6.5 Other polymers (i)

There are no such residues in this entry.