Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 09:07 PM GMT

PDB ID	:	4UZE
Title	:	R66A mutant of FAD synthetase from Corynebacterium ammoniagenes
Authors	:	Martinez-Julvez, M.; Herguedas, B.; Milagros, M.
Deposited on	:	2014-09-05
Resolution	:	2.34 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:

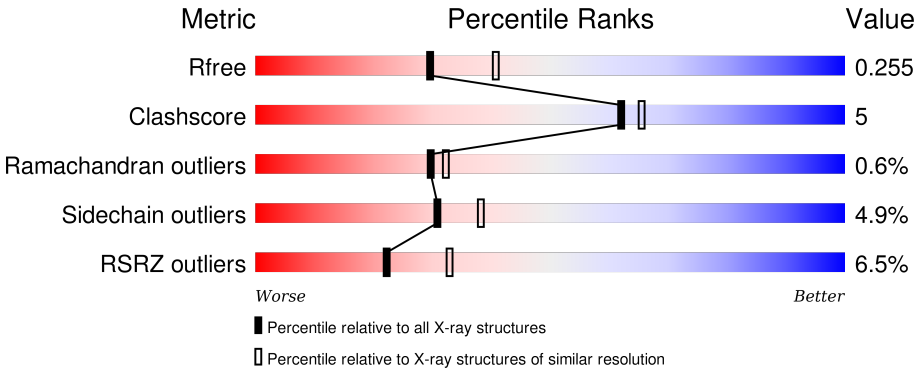
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	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 2.34 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	1406 (2.36-2.32)
Clashscore	102246	1509 (2.36-2.32)
Ramachandran outliers	100387	1490 (2.36-2.32)
Sidechain outliers	100360	1491 (2.36-2.32)
RSRZ outliers	91569	1412 (2.36-2.32)

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	338	5% 88% 10%
1	B	338	8% 87% 11%

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	SO4	A	1339	-	-	-	X
2	SO4	B	1339	-	-	-	X
3	PPV	A	1340	-	-	-	X
3	PPV	B	1340	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5346 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.

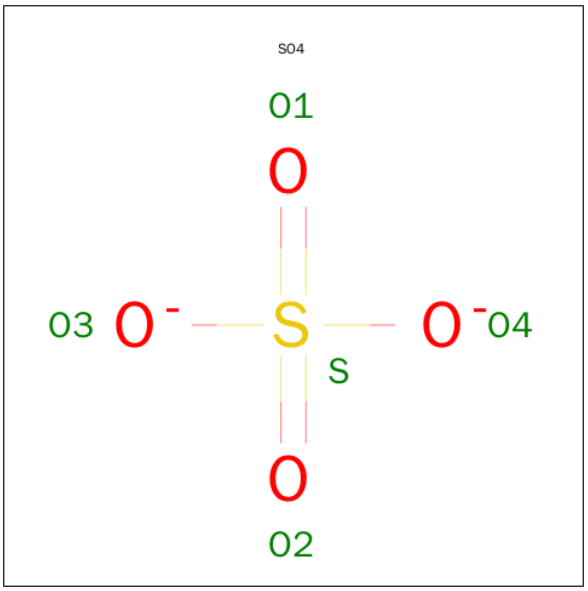
- Molecule 1 is a protein called RIBOFLAVIN BIOSYNTHESIS PROTEIN RIBF.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	338	Total	C	N	O	S	0	0	0
			2594	1642	441	504	7			
1	B	335	Total	C	N	O	S	0	0	0
			2571	1627	438	499	7			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	66	ALA	ARG	ENGINEERED MUTATION	UNP Q59263
B	66	ALA	ARG	ENGINEERED MUTATION	UNP Q59263

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		

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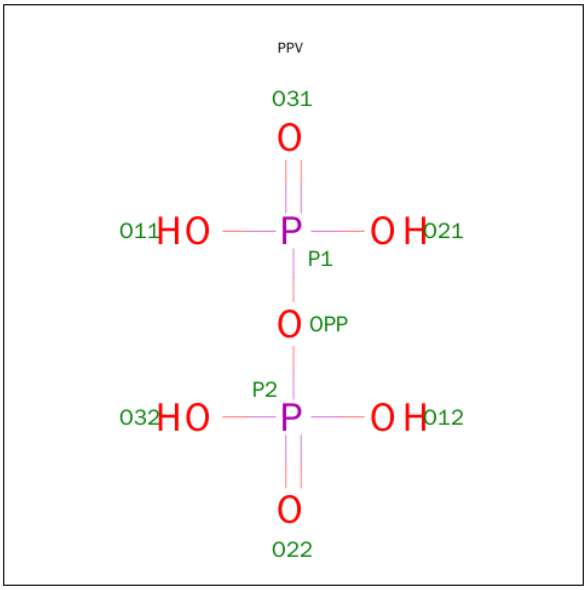
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total O S 5 4 1	0	0

- Molecule 3 is PYROPHOSPHATE (three-letter code: PPV) (formula: $H_4O_7P_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total O P 9 7 2	0	0
3	B	1	Total O P 9 7 2	0	0

- Molecule 4 is water.

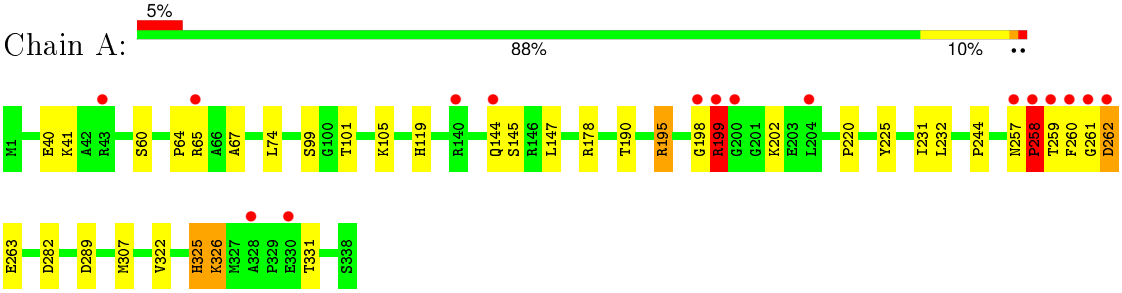
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	103	Total O 103 103	0	0
4	B	50	Total O 50 50	0	0

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.

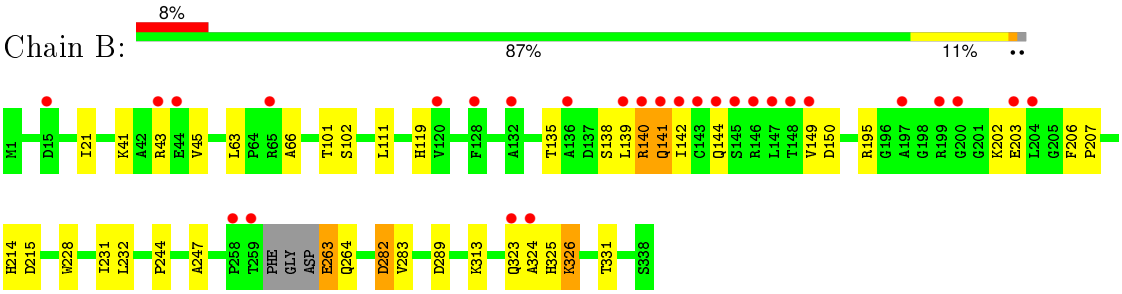
- Molecule 1: RIBOFLAVIN BIOSYNTHESIS PROTEIN RIBF

Chain A:



- Molecule 1: RIBOFLAVIN BIOSYNTHESIS PROTEIN RIBF

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 3	Depositor
Cell constants a, b, c, α, β, γ	133.61Å 133.61Å 133.61Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.00 – 2.34 44.54 – 2.34	Depositor EDS
% Data completeness (in resolution range)	99.5 (40.00-2.34) 99.6 (44.54-2.34)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.27 (at 2.34Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.218 , 0.256 0.220 , 0.255	Depositor DCC
R_{free} test set	1705 reflections (5.33%)	DCC
Wilson B-factor (Å^2)	35.2	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 28.9	EDS
Estimated twinning fraction	0.034 for l,-k,h	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Outliers	0 of 33738 reflections	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	5346	wwPDB-VP
Average B, all atoms (Å^2)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.28% 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: SO4, PPV

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	0.43	0/2649	0.62	2/3604 (0.1%)
1	B	0.39	0/2624	0.56	0/3569
All	All	0.41	0/5273	0.59	2/7173 (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	A	0	1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	258	PRO	CA-N-CD	-7.57	100.91	111.50
1	A	257	ASN	C-N-CD	6.22	141.47	128.40

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	258	PRO	Peptide

5.2 Too-close contacts

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	2594	0	2530	28	0
1	B	2571	0	2513	28	0
2	A	5	0	0	0	0
2	B	5	0	0	0	0
3	A	9	0	0	0	0
3	B	9	0	0	0	0
4	A	103	0	0	0	0
4	B	50	0	0	0	0
All	All	5346	0	5043	55	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 5.

All (55) 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:141:GLN:O	1:B:144:GLN:HG2	1.66	0.94
1:A:261:GLY:HA2	1:A:262:ASP:C	1.86	0.93
1:B:263:GLU:HB3	1:B:264:GLN:HA	1.59	0.85
1:A:258:PRO:HG3	1:A:263:GLU:HA	1.63	0.80
1:A:198:GLY:HA2	1:A:199:ARG:HB2	1.68	0.75
1:A:258:PRO:HG3	1:A:260:PHE:O	1.87	0.75
1:A:258:PRO:CG	1:A:263:GLU:HA	2.22	0.69
1:B:263:GLU:CB	1:B:264:GLN:HA	2.22	0.68
1:B:263:GLU:HB2	1:B:264:GLN:HB2	1.74	0.67
1:A:74:LEU:HD23	1:B:195:ARG:HH21	1.59	0.67
1:B:215:ASP:H	1:B:263:GLU:HG2	1.61	0.66
1:A:60:SER:HB3	1:A:67:ALA:HB2	1.77	0.65
1:A:258:PRO:HB3	1:A:260:PHE:O	2.00	0.62
1:B:41:LYS:HG2	1:B:119:HIS:CD2	2.36	0.61
1:B:63:LEU:HB3	1:B:66:ALA:HB3	1.82	0.61
1:B:263:GLU:CB	1:B:264:GLN:CA	2.80	0.59
1:A:41:LYS:HG2	1:A:119:HIS:CD2	2.38	0.58
1:A:198:GLY:CA	1:A:199:ARG:HB2	2.35	0.57
1:B:231:ILE:HB	1:B:244:PRO:HA	1.87	0.56
1:B:323:GLN:C	1:B:325:HIS:H	2.10	0.55
1:A:258:PRO:CG	1:A:260:PHE:O	2.55	0.54
1:A:202:LYS:HE2	1:A:202:LYS:HA	1.89	0.54
1:A:258:PRO:CB	1:A:260:PHE:O	2.56	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:322:VAL:O	1:A:325:HIS:HD2	1.91	0.54
1:A:261:GLY:HA2	1:A:262:ASP:O	2.08	0.54
1:B:232:LEU:HD12	1:B:282:ASP:HB3	1.90	0.53
1:A:64:PRO:O	1:A:65:ARG:HB2	2.09	0.52
1:B:215:ASP:H	1:B:263:GLU:CG	2.23	0.52
1:B:263:GLU:HB3	1:B:264:GLN:CA	2.35	0.51
1:B:215:ASP:N	1:B:263:GLU:HG2	2.25	0.51
1:B:323:GLN:C	1:B:325:HIS:N	2.64	0.50
1:B:323:GLN:O	1:B:324:ALA:HB3	2.12	0.50
1:B:135:THR:H	1:B:138:SER:HB2	1.78	0.48
1:B:215:ASP:CG	1:B:263:GLU:HG3	2.34	0.48
1:B:101:THR:HG22	1:B:102:SER:O	2.13	0.47
1:B:228:TRP:HB3	1:B:247:ALA:HB1	1.97	0.46
1:A:60:SER:HB3	1:A:67:ALA:CB	2.45	0.45
1:A:258:PRO:C	1:A:260:PHE:H	2.20	0.45
1:A:190:THR:HG21	1:A:232:LEU:HD11	1.97	0.45
1:B:140:ARG:HA	1:B:149:VAL:HG21	1.98	0.44
1:B:149:VAL:HG12	1:B:150:ASP:N	2.33	0.44
1:B:141:GLN:HG2	1:B:142:ILE:N	2.33	0.43
1:A:144:GLN:O	1:A:145:SER:HB3	2.18	0.43
1:A:258:PRO:CA	1:A:260:PHE:H	2.32	0.42
1:A:220:PRO:HB2	1:A:225:TYR:CZ	2.54	0.42
1:B:214:HIS:HA	1:B:263:GLU:HG2	2.00	0.42
1:A:326:LYS:HE2	1:A:326:LYS:H	1.84	0.42
1:B:324:ALA:O	1:B:326:LYS:N	2.42	0.42
1:A:261:GLY:CA	1:A:262:ASP:C	2.74	0.42
1:A:258:PRO:HA	1:A:260:PHE:H	1.85	0.41
1:A:101:THR:CG2	1:A:105:LYS:HB2	2.50	0.41
1:A:231:ILE:HB	1:A:244:PRO:HA	2.02	0.41
1:B:206:PHE:HA	1:B:207:PRO:HD3	1.94	0.40
1:B:21:ILE:HD11	1:B:111:LEU:HD21	2.03	0.40
1:A:195:ARG:HD2	1:A:195:ARG:HA	1.92	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	336/338 (99%)	324 (96%)	9 (3%)	3 (1%)	21	21
1	B	331/338 (98%)	308 (93%)	22 (7%)	1 (0%)	46	54
All	All	667/676 (99%)	632 (95%)	31 (5%)	4 (1%)	30	32

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	258	PRO
1	A	199	ARG
1	A	262	ASP
1	B	45	VAL

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	274/274 (100%)	260 (95%)	14 (5%)	29	36
1	B	272/274 (99%)	259 (95%)	13 (5%)	31	39
All	All	546/548 (100%)	519 (95%)	27 (5%)	31	38

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	40	GLU
1	A	99	SER
1	A	147	LEU
1	A	178	ARG
1	A	195	ARG
1	A	199	ARG
1	A	258	PRO
1	A	259	THR

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Mol	Chain	Res	Type
1	A	282	ASP
1	A	289	ASP
1	A	307	MET
1	A	325	HIS
1	A	326	LYS
1	A	331	THR
1	B	43	ARG
1	B	139	LEU
1	B	140	ARG
1	B	141	GLN
1	B	202	LYS
1	B	203	GLU
1	B	263	GLU
1	B	282	ASP
1	B	283	VAL
1	B	289	ASP
1	B	313	LYS
1	B	326	LYS
1	B	331	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	325	HIS
1	B	141	GLN
1	B	144	GLN
1	B	335	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 carbohydrates in this entry.

5.6 Ligand geometry (i)

4 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	SO4	A	1339	-	4,4,4	0.46	0	6,6,6	0.34	0
3	PPV	A	1340	-	6,8,8	0.67	0	11,13,13	0.99	0
2	SO4	B	1339	-	4,4,4	0.47	0	6,6,6	0.41	0
3	PPV	B	1340	-	6,8,8	0.67	0	11,13,13	0.97	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	SO4	A	1339	-	-	0/0/0/0	0/0/0/0
3	PPV	A	1340	-	-	0/6/6/6	0/0/0/0
2	SO4	B	1339	-	-	0/0/0/0	0/0/0/0
3	PPV	B	1340	-	-	0/6/6/6	0/0/0/0

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 (i)

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	338/338 (100%)	0.20	16 (4%)	35	48	19, 31, 64, 113	0
1	B	335/338 (99%)	0.51	28 (8%)	14	21	21, 39, 81, 107	0
All	All	673/676 (99%)	0.35	44 (6%)	22	33	19, 35, 78, 113	0

All (44) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	259	THR	7.8
1	A	260	PHE	7.2
1	A	65	ARG	6.9
1	A	261	GLY	5.4
1	B	204	LEU	4.7
1	B	324	ALA	4.6
1	A	258	PRO	4.6
1	B	43	ARG	4.6
1	B	259	THR	4.4
1	B	142	ILE	4.3
1	A	199	ARG	4.2
1	B	197	ALA	4.0
1	B	65	ARG	3.8
1	B	145	SER	3.7
1	B	149	VAL	3.6
1	B	200	GLY	3.6
1	B	323	GLN	3.6
1	B	258	PRO	3.6
1	A	257	ASN	3.5
1	A	144	GLN	3.2
1	A	330	GLU	3.1
1	B	136	ALA	3.0
1	A	200	GLY	2.9
1	B	146	ARG	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	144	GLN	2.9
1	B	147	LEU	2.8
1	B	140	ARG	2.8
1	B	128	PHE	2.6
1	B	199	ARG	2.6
1	A	198	GLY	2.5
1	B	132	ALA	2.5
1	B	143	CYS	2.3
1	A	328	ALA	2.3
1	A	140	ARG	2.3
1	A	43	ARG	2.3
1	B	139	LEU	2.3
1	B	203	GLU	2.3
1	B	141	GLN	2.2
1	B	44	GLU	2.2
1	A	204	LEU	2.2
1	B	15	ASP	2.1
1	A	262	ASP	2.1
1	B	148	THR	2.0
1	B	120	VAL	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
3	PPV	A	1340	9/9	0.93	0.19	4.47	46,47,63,67	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å^2)	Q<0.9
2	SO4	B	1339	5/5	0.86	0.18	3.57	69,69,74,74	0
2	SO4	A	1339	5/5	0.86	0.17	3.47	61,68,70,73	0
3	PPV	B	1340	9/9	0.90	0.21	2.53	49,55,69,74	0

6.5 Other polymers (i)

There are no such residues in this entry.