



# Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 03:27 PM GMT

PDB ID : 1R2F  
Title : RIBONUCLEOTIDE REDUCTASE R2F PROTEIN FROM SALMONELLA  
TYPHIMURIUM  
Authors : Eklund, H.; Eriksson, M.  
Deposited on : 1998-08-24  
Resolution : 2.10 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

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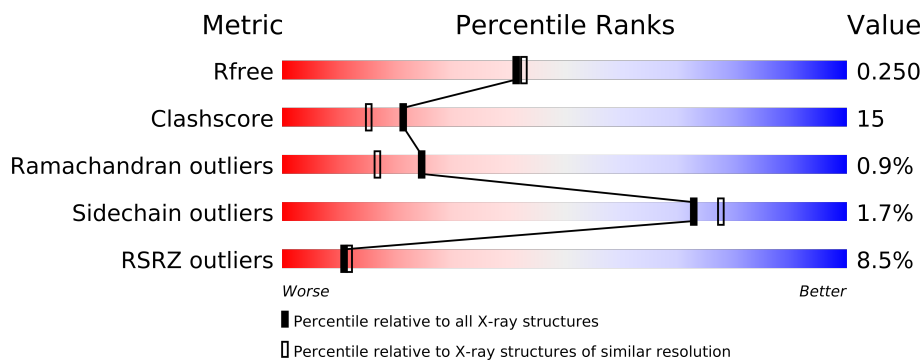
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 2.10 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))
$R_{free}$	66092	3012 (2.10-2.10)
Clashscore	79885	3649 (2.10-2.10)
Ramachandran outliers	78287	3610 (2.10-2.10)
Sidechain outliers	78261	3611 (2.10-2.10)
RSRZ outliers	66119	3013 (2.10-2.10)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	319	
1	B	319	

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4924 atoms, of which 0 are hydrogen and 0 are deuterium.

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 PROTEIN (RIBONUCLEOTIDE REDUCTASE R2).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	283	Total	C	N	O	S	0	0	0
			2268	1471	364	425	8			
1	B	281	Total	C	N	O	S	0	0	0
			2255	1463	362	422	8			

- Molecule 2 is FE (III) ION (three-letter code: FE) (formula: Fe).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	2	Total	Fe	0	0
			2	2		
2	A	2	Total	Fe	0	0
			2	2		

- Molecule 3 is water.

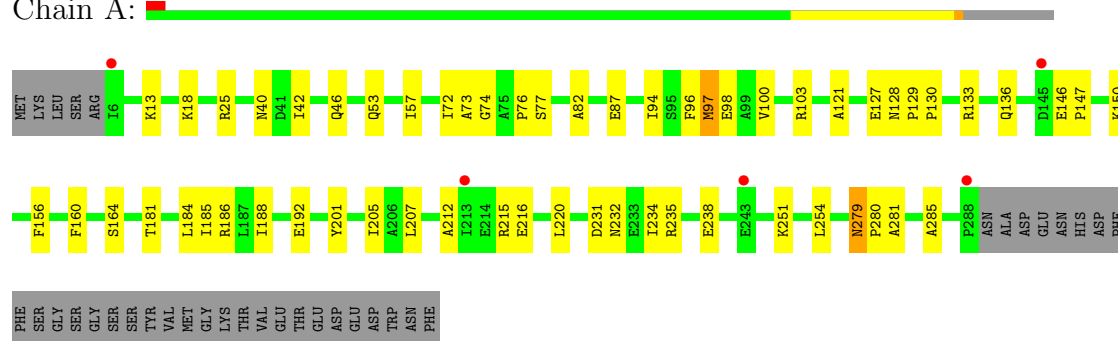
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	266	Total	O	0	0
			266	266		
3	B	131	Total	O	0	0
			131	131		

### 3 Residue-property plots

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.

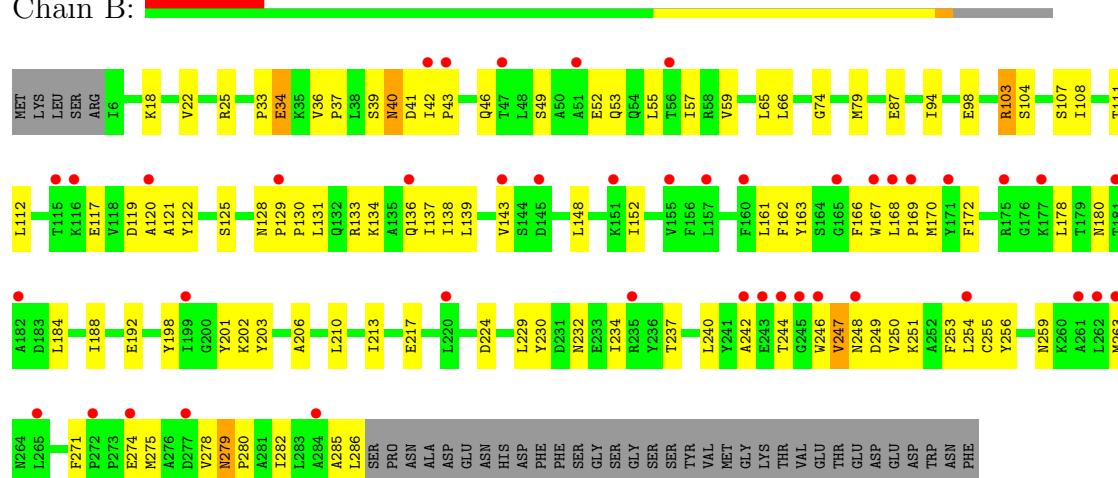
#### • Molecule 1: PROTEIN (RIBONUCLEOTIDE REDUCTASE R2)

Chain A:



#### • Molecule 1: PROTEIN (RIBONUCLEOTIDE REDUCTASE R2)

Chain B:



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	58.49Å 71.74Å 96.07Å 90.00° 95.01° 90.00°	Depositor
Resolution (Å)	20.00 – 2.10 19.42 – 2.08	Depositor EDS
% Data completeness (in resolution range)	93.1 (20.00-2.10) 91.2 (19.42-2.08)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.07	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.93 (at 2.07Å)	Xtriage
Refinement program	CNS	Depositor
R, $R_{free}$	0.226 , 0.247 0.223 , 0.250	Depositor DCC
$R_{free}$ test set	2175 reflections (5.04%)	DCC
Wilson B-factor (Å <sup>2</sup> )	31.9	Xtriage
Anisotropy	0.312	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.38 , 51.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 43452 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	4924	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.20% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: FE

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	0.37	0/2324	0.57	0/3166
1	B	0.30	0/2310	0.52	0/3146
All	All	0.33	0/4634	0.55	0/6312

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2268	0	2247	53	0
1	B	2255	0	2235	86	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	266	0	0	3	0
3	B	131	0	0	0	0
All	All	4924	0	4482	138	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 15.

All (138) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:131:LEU:HA	1:B:134:LYS:HZ3	1.42	0.82
1:A:97:MET:HA	1:A:97:MET:CE	2.11	0.81
1:B:108:ILE:HD11	1:B:184:LEU:HD23	1.66	0.75
1:B:167:TRP:NE1	1:B:282:ILE:HD13	2.04	0.73
1:B:138:ILE:HD11	1:B:161:LEU:CD1	2.21	0.71
1:B:188:ILE:O	1:B:192:GLU:HG2	1.90	0.71
1:A:216:GLU:O	1:A:220:LEU:HD13	1.91	0.69
1:B:254:LEU:HD12	1:B:255:CYS:N	2.08	0.68
1:A:97:MET:HA	1:A:97:MET:HE2	1.75	0.68
1:A:25:ARG:HH11	1:A:25:ARG:HB2	1.61	0.66
1:B:152:ILE:HD11	1:B:203:TYR:CZ	2.31	0.65
1:B:168:LEU:HB3	1:B:169:PRO:HD3	1.78	0.65
1:A:201:TYR:CZ	1:A:205:ILE:HD11	2.31	0.64
1:A:215:ARG:NH1	1:A:215:ARG:HB2	2.12	0.64
1:B:131:LEU:HA	1:B:134:LYS:NZ	2.13	0.64
1:B:129:PRO:HB2	1:B:130:PRO:HD3	1.79	0.63
1:A:279:ASN:ND2	1:A:281:ALA:H	1.96	0.63
1:B:247:VAL:HG12	1:B:251:LYS:HE3	1.79	0.63
1:A:279:ASN:HD22	1:A:280:PRO:N	1.97	0.62
1:A:279:ASN:HD22	1:A:279:ASN:C	2.00	0.62
1:B:42:ILE:N	1:B:43:PRO:HD2	2.15	0.61
1:B:138:ILE:HD11	1:B:161:LEU:HD12	1.83	0.61
1:B:49:SER:OG	1:B:52:GLU:HG3	2.00	0.61
1:A:215:ARG:HB2	1:A:215:ARG:HH11	1.65	0.60
1:A:212:ALA:HA	1:A:215:ARG:NH1	2.17	0.60
1:B:259:ASN:HD21	1:B:271:PHE:H	1.49	0.60
1:B:170:MET:HG3	1:B:286:LEU:HD21	1.84	0.59
1:B:133:ARG:HB3	1:B:232:ASN:HD21	1.68	0.58
1:A:207:LEU:HD22	1:A:215:ARG:HG2	1.86	0.57
1:B:39:SER:O	1:B:42:ILE:HG13	2.04	0.57
1:B:55:LEU:O	1:B:59:VAL:HG23	2.05	0.57
1:B:167:TRP:HE1	1:B:282:ILE:HD13	1.67	0.57
1:A:215:ARG:CB	1:A:215:ARG:HH11	2.19	0.56
1:B:148:LEU:HD12	1:B:210:LEU:HD11	1.87	0.56
1:A:77:SER:OG	1:A:150:LYS:HD3	2.06	0.56
1:A:73:ALA:O	1:A:77:SER:HB2	2.05	0.56
1:A:234:ILE:O	1:A:238:GLU:HG3	2.06	0.56
1:B:25:ARG:HH11	1:B:25:ARG:HG3	1.71	0.55
1:A:97:MET:HA	1:A:97:MET:HE3	1.87	0.55
3:A:559:HOH:O	1:B:79:MET:HE1	2.07	0.55
1:B:230:TYR:O	1:B:234:ILE:HG12	2.07	0.54
1:B:259:ASN:O	1:B:263:MET:HG2	2.08	0.54

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:279:ASN:HD22	1:B:280:PRO:CD	2.21	0.54
1:A:130:PRO:HA	1:A:232:ASN:HD21	1.74	0.53
1:B:46:GLN:OE1	1:B:46:GLN:HA	2.08	0.53
1:B:167:TRP:HE1	1:B:282:ILE:CD1	2.21	0.53
1:B:230:TYR:HE1	1:B:254:LEU:HD13	1.74	0.53
1:B:148:LEU:CD1	1:B:210:LEU:HD11	2.39	0.52
1:A:42:ILE:O	1:A:46:GLN:HG2	2.09	0.52
1:A:279:ASN:HD22	1:A:281:ALA:H	1.56	0.52
1:B:247:VAL:CG1	1:B:251:LYS:HE3	2.40	0.52
1:B:40:ASN:HB2	1:B:180:ASN:OD1	2.11	0.51
1:A:234:ILE:HG23	1:A:251:LYS:HE2	1.92	0.51
1:B:128:ASN:HD22	1:B:131:LEU:HB2	1.77	0.50
1:B:274:GLU:HG3	1:B:275:MET:HG3	1.94	0.50
1:B:57:ILE:HG22	1:B:121:ALA:CB	2.42	0.49
1:B:87:GLU:HG3	1:B:202:LYS:HE3	1.95	0.49
1:B:22:VAL:HG21	1:B:198:TYR:CD1	2.47	0.49
1:B:117:GLU:O	1:B:120:ALA:HB3	2.13	0.49
1:A:96:PHE:CD2	1:A:97:MET:HE3	2.48	0.48
1:A:25:ARG:NH1	1:A:25:ARG:HB2	2.26	0.48
1:B:253:PHE:O	1:B:256:TYR:HB3	2.13	0.48
1:B:128:ASN:HD22	1:B:131:LEU:CB	2.27	0.48
1:B:131:LEU:C	1:B:131:LEU:HD13	2.33	0.48
1:A:97:MET:CE	1:A:100:VAL:HG21	2.44	0.48
1:A:129:PRO:HB2	1:A:130:PRO:HD3	1.96	0.48
1:A:231:ASP:OD2	1:A:235:ARG:NH1	2.44	0.48
1:A:188:ILE:O	1:A:192:GLU:HG2	2.13	0.48
1:B:18:LYS:HD2	1:B:201:TYR:CD2	2.50	0.47
1:B:111:THR:HG22	1:B:111:THR:O	2.14	0.47
1:A:184:LEU:HD23	1:A:184:LEU:C	2.35	0.47
1:B:246:TRP:O	1:B:248:ASN:N	2.48	0.47
1:B:253:PHE:HA	1:B:278:VAL:CG2	2.45	0.47
1:B:213:ILE:O	1:B:217:GLU:HG3	2.14	0.47
1:B:53:GLN:O	1:B:57:ILE:HG13	2.15	0.47
1:B:279:ASN:HD22	1:B:280:PRO:HD2	1.78	0.47
1:B:133:ARG:NH1	1:B:136:GLN:NE2	2.63	0.46
1:B:36:VAL:HA	1:B:37:PRO:HD3	1.78	0.46
1:A:128:ASN:ND2	1:A:130:PRO:HD2	2.30	0.46
1:B:170:MET:SD	1:B:285:ALA:O	2.73	0.46
1:A:57:ILE:HG22	1:A:121:ALA:CB	2.45	0.46
1:A:156:PHE:O	1:A:160:PHE:HB3	2.16	0.46
1:A:72:ILE:O	1:A:76:PRO:HG3	2.15	0.46
1:A:133:ARG:NH1	1:A:136:GLN:OE1	2.48	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:53:GLN:O	1:A:57:ILE:HG13	2.16	0.46
1:B:163:TYR:HD1	1:B:163:TYR:H	1.65	0.45
1:B:94:ILE:O	1:B:98:GLU:HG2	2.16	0.45
1:B:55:LEU:CD1	1:B:240:LEU:HG	2.47	0.45
1:B:244:THR:HB	1:B:246:TRP:CE3	2.51	0.45
1:B:253:PHE:HA	1:B:278:VAL:HG21	1.97	0.45
1:B:172:PHE:HD2	1:B:178:LEU:HD12	1.82	0.45
1:B:163:TYR:HB3	1:B:253:PHE:CE2	2.52	0.44
1:A:94:ILE:O	1:A:98:GLU:HG2	2.16	0.44
1:B:163:TYR:HA	1:B:166:PHE:HB2	1.99	0.44
1:A:13:LYS:HE2	1:B:119:ASP:OD1	2.16	0.44
1:B:33:PRO:HG2	1:B:34:GLU:OE1	2.17	0.44
1:A:164:SER:HB3	1:A:254:LEU:HD21	1.99	0.44
1:A:97:MET:HE2	1:A:100:VAL:CG2	2.47	0.44
1:B:168:LEU:HD11	1:B:172:PHE:CZ	2.52	0.44
1:A:181:THR:O	1:A:185:ILE:HG12	2.18	0.44
1:A:96:PHE:CE2	1:A:97:MET:HE1	2.53	0.44
1:B:167:TRP:CH2	1:B:246:TRP:HB3	2.53	0.44
1:B:25:ARG:HG3	1:B:25:ARG:NH1	2.33	0.44
1:B:230:TYR:HE1	1:B:254:LEU:CD1	2.31	0.43
1:B:66:LEU:HB2	1:B:162:PHE:HE1	1.82	0.43
1:B:111:THR:O	1:B:112:LEU:HD23	2.19	0.43
1:B:103:ARG:C	1:B:103:ARG:HD3	2.39	0.43
1:B:18:LYS:HD2	1:B:201:TYR:CE2	2.54	0.43
1:B:163:TYR:N	1:B:163:TYR:CD1	2.87	0.43
1:A:129:PRO:HD2	1:A:130:PRO:HD3	2.01	0.42
1:B:279:ASN:HD22	1:B:280:PRO:N	2.17	0.42
1:A:82:ALA:HB1	1:A:87:GLU:HB3	2.01	0.42
1:A:207:LEU:O	1:A:215:ARG:HD3	2.19	0.42
1:B:246:TRP:O	1:B:249:ASP:N	2.50	0.42
1:B:65:LEU:HD23	1:B:131:LEU:HD11	2.01	0.42
1:B:166:PHE:O	1:B:169:PRO:HD2	2.19	0.42
1:A:212:ALA:HA	1:A:215:ARG:CZ	2.49	0.42
1:B:139:LEU:O	1:B:143:VAL:HG22	2.20	0.42
1:B:65:LEU:CD2	1:B:131:LEU:HD11	2.49	0.42
1:A:146:GLU:HA	1:A:147:PRO:HD2	1.81	0.42
1:A:186:ARG:NH1	1:A:285:ALA:O	2.53	0.42
1:B:237:THR:HG23	1:B:250:VAL:HG11	2.02	0.41
1:B:103:ARG:HD3	1:B:103:ARG:O	2.20	0.41
1:A:18:LYS:HB2	1:A:18:LYS:HE2	1.95	0.41
1:A:18:LYS:HD2	1:A:201:TYR:CD2	2.55	0.41
1:B:167:TRP:NE1	1:B:282:ILE:CD1	2.77	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:39:SER:C	1:B:41:ASP:H	2.24	0.41
1:B:104:SER:O	1:B:107:SER:HB2	2.21	0.41
1:A:96:PHE:CE2	1:A:97:MET:CE	3.04	0.41
1:B:246:TRP:O	1:B:247:VAL:C	2.59	0.41
1:B:172:PHE:CD2	1:B:178:LEU:HD12	2.56	0.41
1:B:148:LEU:HD21	1:B:206:ALA:CB	2.51	0.41
1:B:137:ILE:HG21	1:B:229:LEU:HD21	2.02	0.41
1:A:97:MET:HE2	1:A:100:VAL:HB	2.02	0.41
1:A:136:GLN:HG2	3:A:591:HOH:O	2.21	0.40
1:A:40:ASN:ND2	3:A:420:HOH:O	2.54	0.40
1:B:122:TYR:O	1:B:125:SER:HB3	2.21	0.40
1:A:18:LYS:HD2	1:A:201:TYR:CE2	2.57	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles

### 5.3.1 Protein backbone ⓘ

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	281/319 (88%)	276 (98%)	4 (1%)	1 (0%)	43	39
1	B	279/319 (88%)	265 (95%)	10 (4%)	4 (1%)	16	9
All	All	560/638 (88%)	541 (97%)	14 (2%)	5 (1%)	25	17

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	74	GLY
1	B	247	VAL
1	B	40	ASN
1	B	242	ALA
1	A	74	GLY

### 5.3.2 Protein sidechains ⓘ

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	243/275 (88%)	239 (98%)	4 (2%)	75	79
1	B	241/275 (88%)	237 (98%)	4 (2%)	73	78
All	All	484/550 (88%)	476 (98%)	8 (2%)	73	78

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	97	MET
1	A	103	ARG
1	A	127	GLU
1	A	279	ASN
1	B	34	GLU
1	B	103	ARG
1	B	224	ASP
1	B	279	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	15	GLN
1	A	29	ASN
1	A	40	ASN
1	A	46	GLN
1	A	53	GLN
1	A	71	ASN
1	A	128	ASN
1	A	132	GLN
1	A	141	HIS
1	A	208	GLN
1	A	232	ASN
1	A	264	ASN
1	A	279	ASN
1	B	29	ASN
1	B	40	ASN
1	B	53	GLN

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Mol	Chain	Res	Type
1	B	71	ASN
1	B	128	ASN
1	B	132	GLN
1	B	136	GLN
1	B	180	ASN
1	B	208	GLN
1	B	232	ASN
1	B	257	ASN
1	B	259	ASN
1	B	264	ASN
1	B	279	ASN

### 5.3.3 RNA ⓘ

There are no RNA chains 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 4 ligands modelled in this entry, 4 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.

## 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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	283/319 (88%)	0.03	5 (1%) 65 69	23, 30, 43, 45	0
1	B	281/319 (88%)	0.96	43 (15%) 3 3	27, 55, 73, 81	0
All	All	564/638 (88%)	0.50	48 (8%) 11 12	23, 37, 69, 81	0

All (48) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	243	GLU	5.6
1	B	274	GLU	4.5
1	B	175	ARG	4.5
1	B	254	LEU	4.2
1	B	116	LYS	4.2
1	B	245	GLY	3.9
1	B	145	ASP	3.8
1	B	246	TRP	3.7
1	A	145	ASP	3.7
1	B	284	ALA	3.5
1	B	220	LEU	3.2
1	B	248	ASN	3.1
1	B	263	MET	3.1
1	B	277	ASP	3.1
1	B	155	VAL	3.1
1	B	51	ALA	3.0
1	B	171	TYR	3.0
1	B	235	ARG	3.0
1	B	168	LEU	2.9
1	B	181	THR	2.8
1	B	242	ALA	2.8
1	B	265	LEU	2.8
1	B	47	THR	2.7
1	A	213	ILE	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	288	PRO	2.6
1	A	243	GLU	2.6
1	B	143	VAL	2.6
1	B	56	THR	2.6
1	A	6	ILE	2.5
1	B	261	ALA	2.5
1	B	120	ALA	2.5
1	B	43	PRO	2.5
1	B	244	THR	2.5
1	B	167	TRP	2.4
1	B	272	PRO	2.4
1	B	182	ALA	2.4
1	B	115	THR	2.4
1	B	42	ILE	2.3
1	B	129	PRO	2.3
1	B	165	GLY	2.2
1	B	169	PRO	2.2
1	B	160	PHE	2.1
1	B	157	LEU	2.1
1	B	199	ILE	2.1
1	B	136	GLN	2.1
1	B	262	LEU	2.0
1	B	151	LYS	2.0
1	B	177	LYS	2.0

## 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 ⓘ

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<sup>th</sup> 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	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	FE	B	403	1/1	0.04	-3.13	43,43,43,43	0
2	FE	B	402	1/1	0.03	-3.26	43,43,43,43	1
2	FE	A	400	1/1	0.02	-7.27	32,32,32,32	0
2	FE	A	401	1/1	0.01	-7.92	31,31,31,31	0

## 6.5 Other polymers

There are no such residues in this entry.