



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 19, 2016 – 06:39 PM GMT

PDB ID : 1SZE
Title : L230A mutant flavocytochrome b2 with benzoylformate
Authors : Mowat, C.G.; Wehenkel, A.; Green, A.J.; Walkinshaw, M.D.; Reid, G.A.; Chapman, S.K.
Deposited on : 2004-04-05
Resolution : 3.00 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 : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026982
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) : rb-20026982

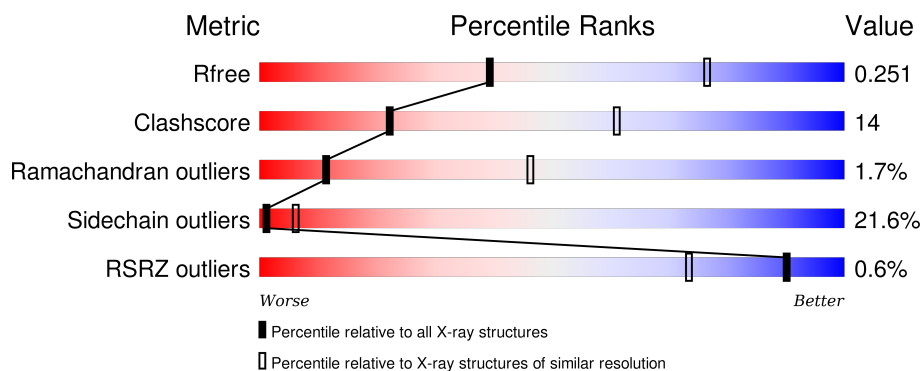
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	1578 (3.00-3.00)
Clashscore	102246	1912 (3.00-3.00)
Ramachandran outliers	100387	1853 (3.00-3.00)
Sidechain outliers	100360	1856 (3.00-3.00)
RSRZ outliers	91569	1592 (3.00-3.00)

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	511	 46% 25% 5% 23%
1	B	511	 42% 27% 7% 24%

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
3	173	A	5580	-	-	-	X
3	173	B	9580	-	-	-	X

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 6291 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 Cytochrome b2, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	392	Total	C	N	O	S	0	0	0
			3053	1936	521	585	11			
1	B	389	Total	C	N	O	S	0	0	0
			3030	1924	514	581	11			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	230	ALA	LEU	ENGINEERED	UNP P00175
B	230	ALA	LEU	ENGINEERED	UNP P00175

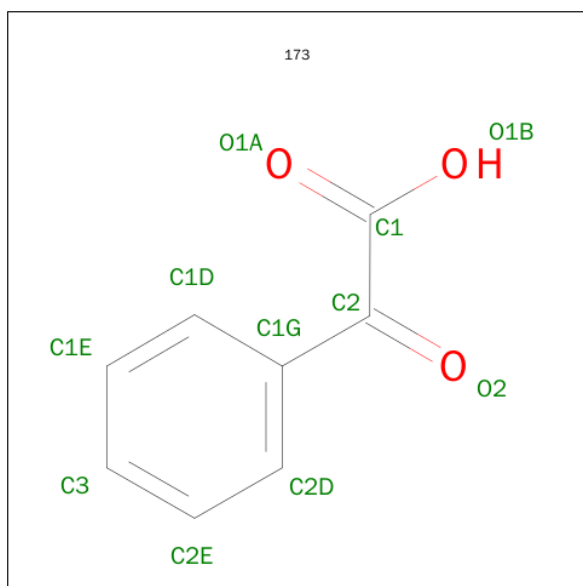
- Molecule 2 is FLAVIN MONONUCLEOTIDE (three-letter code: FMN) (formula: C₁₇H₂₁N₄O₉P).



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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	B	1	Total	C	N	O	P	0	0
			31	17	4	9	1		

- Molecule 3 is BENZOYL-FORMIC ACID (three-letter code: 173) (formula: $C_8H_6O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			11	8	3		
3	B	1	Total	C	O	0	0
			11	8	3		

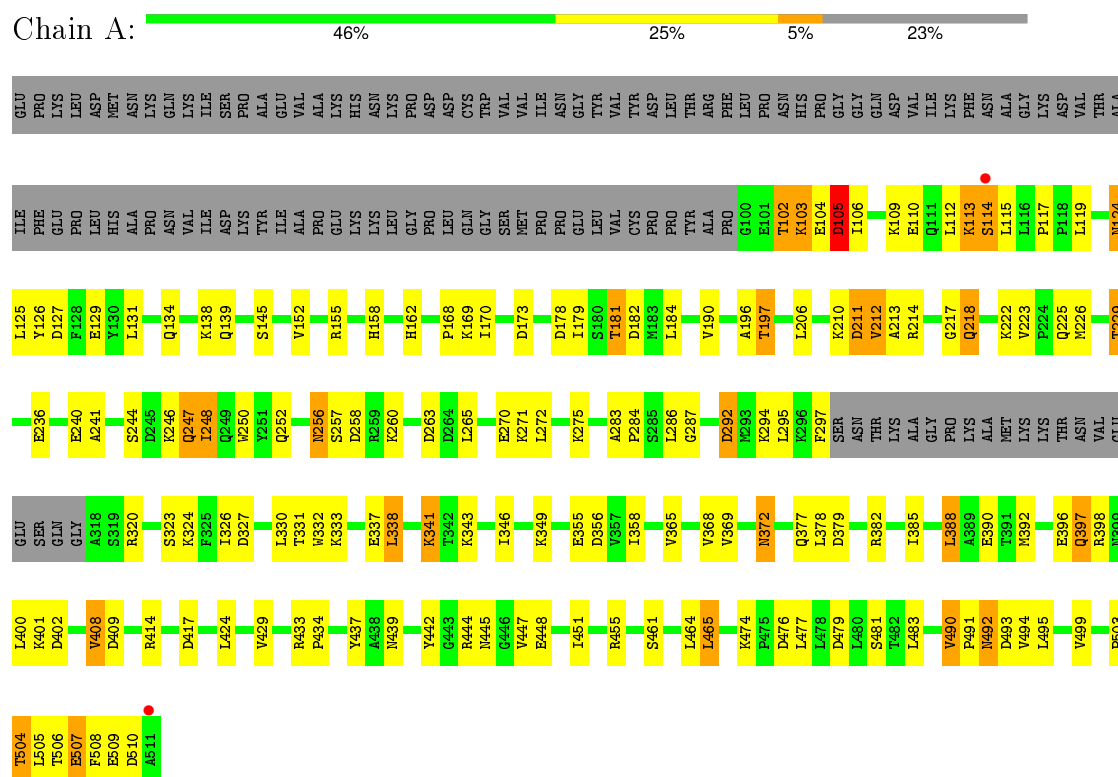
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	69	Total	O	0	0
			69	69		
4	B	55	Total	O	0	0
			55	55		

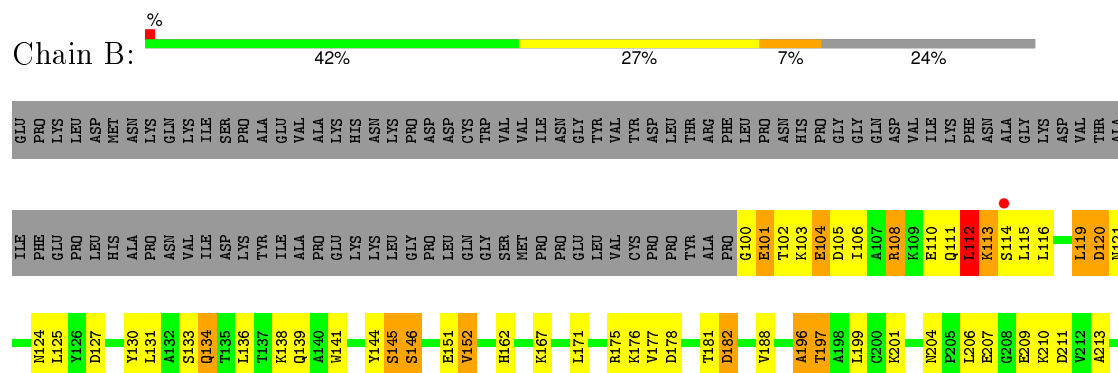
3 Residue-property plots

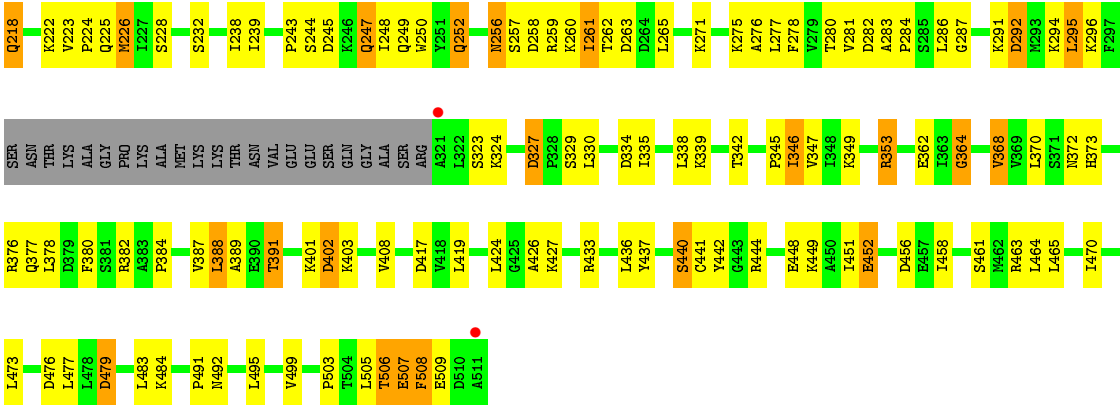
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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: Cytochrome b2, mitochondrial



- Molecule 1: Cytochrome b2, mitochondrial





4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	163.03Å 163.03Å 112.47Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	24.00 – 3.00 23.80 – 3.00	Depositor EDS
% Data completeness (in resolution range)	90.6 (24.00-3.00) 90.6 (23.80-3.00)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.47 (at 2.99Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.200 , 0.249 0.206 , 0.251	Depositor DCC
R_{free} test set	1577 reflections (5.27%)	DCC
Wilson B-factor (Å ²)	57.5	Xtriage
Anisotropy	0.010	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 48.2	EDS
Estimated twinning fraction	0.024 for -h,-k,l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 31502 reflections	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	6291	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.77% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: FMN, 173

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	0.77	1/3102 (0.0%)	0.99	9/4191 (0.2%)
1	B	0.74	0/3079	1.00	18/4161 (0.4%)
All	All	0.76	1/6181 (0.0%)	0.99	27/8352 (0.3%)

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	2
1	B	0	2
All	All	0	4

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	492	ASN	CB-CG	5.63	1.64	1.51

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	120	ASP	CB-CG-OD2	8.35	125.81	118.30
1	A	327	ASP	CB-CG-OD2	8.25	125.72	118.30
1	B	245	ASP	CB-CG-OD2	7.17	124.76	118.30
1	B	456	ASP	CB-CG-OD2	7.15	124.74	118.30
1	A	402	ASP	CB-CG-OD2	6.86	124.47	118.30
1	B	182	ASP	CB-CG-OD2	6.86	124.47	118.30
1	B	211	ASP	CB-CG-OD2	6.64	124.28	118.30
1	B	327	ASP	CB-CG-OD2	6.59	124.23	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	479	ASP	CB-CG-OD2	6.51	124.16	118.30
1	B	112	LEU	CA-CB-CG	6.28	129.74	115.30
1	A	493	ASP	CB-CG-OD2	6.17	123.85	118.30
1	A	178	ASP	CB-CG-OD2	6.12	123.81	118.30
1	B	509	GLU	N-CA-C	5.92	126.99	111.00
1	A	127	ASP	CB-CG-OD2	5.90	123.61	118.30
1	B	507	GLU	N-CA-C	5.83	126.74	111.00
1	A	356	ASP	CB-CG-OD2	5.75	123.47	118.30
1	B	417	ASP	CB-CG-OD2	5.68	123.42	118.30
1	A	292	ASP	CB-CG-OD2	5.45	123.20	118.30
1	B	334	ASP	CB-CG-OD2	5.37	123.14	118.30
1	B	127	ASP	CB-CG-OD2	5.37	123.13	118.30
1	B	463	ARG	NE-CZ-NH2	-5.30	117.65	120.30
1	A	355	GLU	OE1-CD-OE2	-5.18	117.09	123.30
1	B	292	ASP	CB-CG-OD2	5.16	122.95	118.30
1	B	108	ARG	NE-CZ-NH1	5.14	122.87	120.30
1	A	263	ASP	CB-CG-OD2	5.06	122.86	118.30
1	B	178	ASP	CB-CG-OD2	5.03	122.83	118.30
1	B	402	ASP	CB-CG-OD2	5.02	122.81	118.30

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	492	ASN	Peptide
1	A	506	THR	Peptide
1	B	104	GLU	Peptide
1	B	506	THR	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	3053	0	3113	83	0
1	B	3030	0	3088	97	0
2	A	31	0	19	4	0
2	B	31	0	18	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	11	0	5	1	0
3	B	11	0	5	0	0
4	A	69	0	0	3	0
4	B	55	0	0	4	0
All	All	6291	0	6248	171	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 14.

All (171) 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:A:256:ASN:HD22	1:A:258:ASP:H	1.07	0.94
1:A:113:LYS:O	1:A:115:LEU:N	2.02	0.91
1:A:218:GLN:HE22	1:A:444:ARG:HD2	1.39	0.88
1:A:331:THR:HB	4:A:5623:HOH:O	1.80	0.80
1:A:211:ASP:HB3	1:A:439:ASN:HD21	1.47	0.79
1:B:197:THR:HG21	1:B:436:LEU:CD1	2.11	0.79
1:A:129:GLU:HG3	1:A:437:TYR:CE2	2.19	0.78
1:B:197:THR:HG21	1:B:436:LEU:HD11	1.64	0.78
1:A:256:ASN:ND2	1:A:258:ASP:H	1.84	0.75
1:B:105:ASP:HA	1:B:108:ARG:HB2	1.66	0.75
1:B:102:THR:O	1:B:102:THR:HG22	1.87	0.74
1:B:339:LYS:NZ	1:B:364:GLY:O	2.22	0.73
1:B:113:LYS:O	1:B:115:LEU:N	2.22	0.73
1:A:124:ASN:C	1:A:124:ASN:HD22	1.95	0.70
1:A:113:LYS:HG3	1:A:113:LYS:O	1.93	0.68
1:A:479:ASP:OD2	1:B:382:ARG:HD3	1.94	0.67
1:B:507:GLU:O	1:B:508:PHE:HB2	1.96	0.66
1:A:119:LEU:HD21	1:A:445:ASN:O	1.96	0.65
1:A:499:VAL:HG13	1:B:505:LEU:HD11	1.80	0.64
1:A:113:LYS:C	1:A:115:LEU:H	2.00	0.64
1:A:223:VAL:HG21	1:A:451:ILE:HG12	1.79	0.64
1:B:136:LEU:HD21	1:B:440:SER:HB3	1.81	0.63
1:A:495:LEU:HD13	1:B:503:PRO:HB2	1.81	0.62
1:B:108:ARG:O	1:B:112:LEU:HD23	1.99	0.62
1:B:111:GLN:O	1:B:115:LEU:HB2	1.99	0.62
1:B:104:GLU:C	1:B:106:ILE:H	2.02	0.62
1:A:155:ARG:NH2	1:B:492:ASN:O	2.32	0.61
1:B:256:ASN:HD22	1:B:258:ASP:H	1.49	0.61
1:B:243:PRO:HD2	1:B:247:GLN:HE22	1.65	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:213:ALA:CA	1:B:225:GLN:HE22	2.15	0.60
1:B:124:ASN:ND2	4:B:9623:HOH:O	2.32	0.60
1:B:113:LYS:C	1:B:115:LEU:H	2.06	0.59
1:B:119:LEU:HD23	1:B:442:TYR:CD1	2.38	0.59
1:B:349:LYS:NZ	2:B:9570:FMN:O2'	2.35	0.59
1:A:213:ALA:HB2	1:A:225:GLN:HE22	1.67	0.59
1:A:408:VAL:HG12	1:A:429:VAL:HB	1.84	0.58
1:B:197:THR:CG2	1:B:436:LEU:HD11	2.32	0.58
1:A:382:ARG:HD2	4:A:5583:HOH:O	2.03	0.58
1:B:389:ALA:HB2	1:B:424:LEU:HD13	1.86	0.57
1:B:144:TYR:CD2	1:B:436:LEU:HD13	2.40	0.57
1:B:278:PHE:CE2	1:B:347:VAL:HG11	2.40	0.56
1:B:119:LEU:HD23	1:B:442:TYR:CG	2.41	0.56
1:A:210:LYS:O	1:A:214:ARG:HG3	2.06	0.56
1:B:120:ASP:C	1:B:121:ASN:HD22	2.10	0.56
1:A:331:THR:HG22	1:A:332:TRP:N	2.21	0.55
1:A:129:GLU:HG3	1:A:437:TYR:CD2	2.42	0.55
1:A:173:ASP:OD1	1:B:329:SER:HA	2.05	0.55
1:B:218:GLN:HE22	1:B:444:ARG:HG3	1.72	0.55
1:B:275:LYS:O	1:B:345:PRO:HD2	2.07	0.55
1:A:385:ILE:HD11	1:A:424:LEU:HD12	1.89	0.54
1:B:388:LEU:HD11	1:B:426:ALA:HB2	1.89	0.53
1:B:448:GLU:O	1:B:452:GLU:HG2	2.09	0.53
1:B:223:VAL:HG21	1:B:451:ILE:HG12	1.91	0.53
1:A:337:GLU:HB3	1:A:341:LYS:HZ3	1.72	0.53
1:B:213:ALA:HB2	1:B:225:GLN:HE22	1.74	0.52
1:B:327:ASP:OD1	1:B:329:SER:OG	2.27	0.52
1:A:155:ARG:CZ	1:B:491:PRO:HB2	2.39	0.52
1:B:144:TYR:CG	1:B:436:LEU:HD13	2.45	0.52
1:A:124:ASN:C	1:A:124:ASN:ND2	2.63	0.52
1:B:389:ALA:HB2	1:B:424:LEU:CD1	2.39	0.52
1:B:209:GLU:HG3	1:B:238:ILE:HD11	1.92	0.52
1:B:207:GLU:O	1:B:210:LYS:HB2	2.10	0.52
1:A:119:LEU:HD13	1:A:442:TYR:CD1	2.45	0.51
1:A:349:LYS:NZ	2:A:5570:FMN:N1	2.58	0.51
1:B:287:GLY:H	1:B:377:GLN:HE22	1.58	0.51
1:A:113:LYS:C	1:A:115:LEU:N	2.62	0.51
1:A:499:VAL:HG13	1:B:505:LEU:CD1	2.40	0.51
1:B:196:ALA:HB2	1:B:226:MET:HG2	1.93	0.51
1:B:213:ALA:HA	1:B:225:GLN:HE22	1.76	0.50
1:B:152:VAL:HG21	1:B:380:PHE:CE1	2.47	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:197:THR:HG21	1:B:436:LEU:CG	2.42	0.50
1:A:181:THR:OG1	1:A:182:ASP:N	2.45	0.49
1:A:283:ALA:N	1:A:284:PRO:CD	2.76	0.49
1:A:490:VAL:HG23	1:A:491:PRO:HD2	1.95	0.49
1:A:409:ASP:OD2	2:A:5570:FMN:O3'	2.24	0.48
1:A:256:ASN:HD22	1:A:257:SER:N	2.11	0.48
1:A:117:PRO:HD2	1:A:442:TYR:OH	2.13	0.48
1:A:349:LYS:HG3	1:A:369:VAL:HG21	1.94	0.48
1:A:326:ILE:HD11	3:A:5580:173:H1E	1.95	0.48
1:B:347:VAL:HG13	1:B:368:VAL:N	2.29	0.48
1:A:504:THR:CG2	4:A:5581:HOH:O	2.61	0.48
1:A:170:ILE:HG12	1:B:281:VAL:HG23	1.94	0.48
1:B:283:ALA:N	1:B:284:PRO:CD	2.77	0.48
1:B:102:THR:CG2	1:B:102:THR:O	2.57	0.48
1:B:138:LYS:HB3	1:B:295:LEU:CD1	2.44	0.48
1:A:236:GLU:HG3	1:A:272:LEU:HD21	1.96	0.48
1:B:213:ALA:CB	1:B:225:GLN:HE22	2.26	0.47
1:A:505:LEU:HD22	1:B:130:TYR:HB2	1.96	0.47
1:A:102:THR:C	1:A:104:GLU:N	2.67	0.47
1:A:102:THR:C	1:A:104:GLU:H	2.16	0.47
1:A:217:GLY:HA3	1:A:247:GLN:OE1	2.15	0.47
1:B:141:TRP:O	1:B:145:SER:OG	2.31	0.47
1:A:287:GLY:H	1:A:377:GLN:HE22	1.62	0.47
1:B:121:ASN:HD22	1:B:121:ASN:N	2.11	0.47
1:A:179:ILE:HD11	1:A:455:ARG:HG3	1.95	0.47
1:B:465:LEU:HD13	1:B:477:LEU:HD23	1.97	0.47
1:B:226:MET:CE	1:B:349:LYS:HD2	2.44	0.47
1:B:282:ASP:O	1:B:377:GLN:HG3	2.15	0.47
1:A:507:GLU:O	1:A:508:PHE:HB2	2.15	0.47
1:B:295:LEU:HD23	1:B:296:LYS:N	2.30	0.47
1:B:209:GLU:OE2	1:B:232:SER:HB2	2.15	0.47
1:B:470:ILE:HA	1:B:473:LEU:HD12	1.97	0.47
1:A:109:LYS:HA	1:A:112:LEU:HD13	1.96	0.46
1:A:372:ASN:C	1:A:372:ASN:HD22	2.18	0.46
1:A:248:ILE:HD11	1:A:250:TRP:CE2	2.50	0.46
1:A:168:PRO:O	1:B:353:ARG:HD2	2.15	0.46
1:B:507:GLU:O	1:B:508:PHE:CB	2.62	0.46
1:A:126:TYR:OH	1:A:158:HIS:CE1	2.69	0.46
1:B:335:ILE:O	1:B:338:LEU:N	2.49	0.46
1:A:213:ALA:CB	1:A:225:GLN:HE22	2.28	0.45
1:A:287:GLY:H	1:A:377:GLN:NE2	2.15	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:252:GLN:HA	1:B:278:PHE:O	2.16	0.45
1:B:258:ASP:HB3	1:B:261:ILE:HG13	1.99	0.45
1:A:211:ASP:HB3	1:A:439:ASN:ND2	2.24	0.45
1:A:433:ARG:N	1:A:434:PRO:CD	2.79	0.45
1:B:197:THR:HG22	2:B:9570:FMN:HM73	1.98	0.45
1:A:388:LEU:HD22	1:A:392:MET:HG2	1.98	0.45
1:A:337:GLU:OE1	1:A:341:LYS:NZ	2.49	0.45
1:B:104:GLU:C	1:B:106:ILE:N	2.68	0.44
1:A:461:SER:O	1:A:465:LEU:HB2	2.18	0.44
1:B:171:LEU:HA	1:B:171:LEU:HD23	1.92	0.44
1:A:229:THR:H	1:A:252:GLN:NE2	2.16	0.44
1:B:338:LEU:HG	1:B:346:ILE:HD11	2.00	0.44
1:A:414:ARG:O	1:A:417:ASP:HB2	2.18	0.44
1:A:105:ASP:O	1:A:109:LYS:HB2	2.18	0.44
1:B:100:GLY:O	1:B:138:LYS:HD2	2.18	0.44
1:B:213:ALA:HA	1:B:225:GLN:NE2	2.33	0.44
1:B:495:LEU:O	1:B:499:VAL:HG22	2.18	0.44
1:B:479:ASP:C	1:B:479:ASP:OD1	2.56	0.44
1:B:108:ARG:HH22	1:B:138:LYS:N	2.16	0.43
1:A:124:ASN:HD22	1:A:126:TYR:N	2.17	0.43
1:B:372:ASN:ND2	1:B:384:PRO:HD3	2.34	0.43
1:A:256:ASN:HD22	1:A:258:ASP:N	1.92	0.43
1:A:197:THR:O	1:A:197:THR:CG2	2.67	0.43
1:A:210:LYS:HG2	1:A:241:ALA:HB2	2.01	0.43
2:B:9570:FMN:H9	2:B:9570:FMN:H1'2	1.82	0.43
1:B:209:GLU:CG	1:B:238:ILE:HD11	2.49	0.43
1:B:387:VAL:O	1:B:391:THR:HG23	2.19	0.43
1:B:226:MET:HE3	1:B:349:LYS:HD2	2.00	0.42
1:B:250:TRP:CD1	1:B:276:ALA:HB3	2.54	0.42
1:A:170:ILE:HA	1:A:170:ILE:HD12	1.86	0.42
1:A:244:SER:HB2	1:A:247:GLN:HB3	2.01	0.42
1:B:133:SER:OG	1:B:134:GLN:NE2	2.52	0.42
1:A:252:GLN:NE2	2:A:5570:FMN:HN3	2.16	0.42
1:A:349:LYS:HZ2	2:A:5570:FMN:C2	2.32	0.42
1:A:196:ALA:HB2	1:A:226:MET:HE2	2.01	0.42
1:B:238:ILE:HG22	1:B:239:ILE:N	2.35	0.42
1:B:101:GLU:H	1:B:101:GLU:HG2	1.60	0.42
1:B:181:THR:HG21	4:B:9606:HOH:O	2.18	0.42
1:B:151:GLU:OE2	1:B:291:LYS:NZ	2.53	0.42
1:B:201:LYS:HA	1:B:204:ASN:O	2.20	0.41
1:B:252:GLN:NE2	1:B:280:THR:OG1	2.53	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:449:LYS:HA	1:B:452:GLU:CG	2.50	0.41
1:A:170:ILE:HB	4:B:9594:HOH:O	2.20	0.41
1:B:419:LEU:HD21	1:B:458:ILE:HG23	2.02	0.41
1:A:409:ASP:OD1	1:A:409:ASP:C	2.58	0.41
1:B:373:HIS:O	1:B:376:ARG:HB2	2.20	0.41
1:B:108:ARG:NH1	4:B:9589:HOH:O	2.46	0.41
1:A:448:GLU:HA	1:A:451:ILE:HD12	2.03	0.41
1:B:433:ARG:HB3	1:B:437:TYR:CE2	2.56	0.41
1:B:188:VAL:HG11	1:B:224:PRO:HG3	2.03	0.41
1:B:449:LYS:HA	1:B:452:GLU:HG3	2.03	0.41
1:A:212:VAL:HA	1:A:447:VAL:HG21	2.02	0.41
1:A:331:THR:CG2	1:A:332:TRP:N	2.84	0.40
1:A:102:THR:O	1:A:104:GLU:N	2.54	0.40
1:A:214:ARG:O	1:A:218:GLN:HB3	2.20	0.40
1:B:197:THR:HG21	1:B:436:LEU:HG	2.02	0.40
1:A:337:GLU:O	1:A:341:LYS:HD3	2.21	0.40
1:B:461:SER:O	1:B:465:LEU:HB2	2.22	0.40
1:A:338:LEU:HD13	1:A:346:ILE:HD13	2.04	0.40
1:A:338:LEU:HD13	1:A:346:ILE:CD1	2.52	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	388/511 (76%)	355 (92%)	26 (7%)	7 (2%)	11	45
1	B	385/511 (75%)	334 (87%)	45 (12%)	6 (2%)	12	48
All	All	773/1022 (76%)	689 (89%)	71 (9%)	13 (2%)	11	46

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	103	LYS
1	A	114	SER
1	A	397	GLN
1	B	114	SER
1	B	146	SER
1	A	105	ASP
1	A	138	LYS
1	B	362	GLU
1	B	364	GLY
1	B	508	PHE
1	A	510	ASP
1	B	196	ALA
1	A	503	PRO

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	336/439 (76%)	261 (78%)	75 (22%)	1	5
1	B	334/439 (76%)	264 (79%)	70 (21%)	1	7
All	All	670/878 (76%)	525 (78%)	145 (22%)	1	6

All (145) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	102	THR
1	A	103	LYS
1	A	105	ASP
1	A	106	ILE
1	A	110	GLU
1	A	113	LYS
1	A	114	SER
1	A	124	ASN
1	A	125	LEU
1	A	131	LEU
1	A	134	GLN

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Mol	Chain	Res	Type
1	A	139	GLN
1	A	145	SER
1	A	152	VAL
1	A	162	HIS
1	A	169	LYS
1	A	181	THR
1	A	184	LEU
1	A	190	VAL
1	A	197	THR
1	A	206	LEU
1	A	211	ASP
1	A	212	VAL
1	A	218	GLN
1	A	222	LYS
1	A	229	THR
1	A	240	GLU
1	A	246	LYS
1	A	247	GLN
1	A	248	ILE
1	A	256	ASN
1	A	260	LYS
1	A	265	LEU
1	A	270	GLU
1	A	271	LYS
1	A	275	LYS
1	A	286	LEU
1	A	292	ASP
1	A	294	LYS
1	A	295	LEU
1	A	297	PHE
1	A	320	ARG
1	A	323	SER
1	A	324	LYS
1	A	330	LEU
1	A	333	LYS
1	A	338	LEU
1	A	341	LYS
1	A	343	LYS
1	A	358	ILE
1	A	365	VAL
1	A	368	VAL
1	A	372	ASN

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Mol	Chain	Res	Type
1	A	378	LEU
1	A	379	ASP
1	A	388	LEU
1	A	390	GLU
1	A	396	GLU
1	A	397	GLN
1	A	398	ARG
1	A	400	LEU
1	A	401	LYS
1	A	408	VAL
1	A	464	LEU
1	A	465	LEU
1	A	474	LYS
1	A	476	ASP
1	A	477	LEU
1	A	481	SER
1	A	483	LEU
1	A	490	VAL
1	A	494	VAL
1	A	504	THR
1	A	507	GLU
1	A	509	GLU
1	B	101	GLU
1	B	103	LYS
1	B	110	GLU
1	B	112	LEU
1	B	113	LYS
1	B	116	LEU
1	B	119	LEU
1	B	125	LEU
1	B	131	LEU
1	B	134	GLN
1	B	139	GLN
1	B	145	SER
1	B	146	SER
1	B	152	VAL
1	B	162	HIS
1	B	167	LYS
1	B	175	ARG
1	B	176	LYS
1	B	177	VAL
1	B	182	ASP

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Mol	Chain	Res	Type
1	B	197	THR
1	B	199	LEU
1	B	206	LEU
1	B	218	GLN
1	B	222	LYS
1	B	226	MET
1	B	228	SER
1	B	244	SER
1	B	247	GLN
1	B	248	ILE
1	B	249	GLN
1	B	252	GLN
1	B	256	ASN
1	B	257	SER
1	B	259	ARG
1	B	260	LYS
1	B	261	ILE
1	B	262	THR
1	B	263	ASP
1	B	265	LEU
1	B	271	LYS
1	B	277	LEU
1	B	286	LEU
1	B	292	ASP
1	B	294	LYS
1	B	295	LEU
1	B	323	SER
1	B	324	LYS
1	B	330	LEU
1	B	342	THR
1	B	346	ILE
1	B	353	ARG
1	B	368	VAL
1	B	370	LEU
1	B	378	LEU
1	B	388	LEU
1	B	391	THR
1	B	401	LYS
1	B	402	ASP
1	B	403	LYS
1	B	408	VAL
1	B	427	LYS

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Mol	Chain	Res	Type
1	B	440	SER
1	B	441	CYS
1	B	452	GLU
1	B	464	LEU
1	B	476	ASP
1	B	483	LEU
1	B	484	LYS
1	B	506	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	124	ASN
1	A	134	GLN
1	A	139	GLN
1	A	157	ASN
1	A	218	GLN
1	A	225	GLN
1	A	252	GLN
1	A	256	ASN
1	A	372	ASN
1	A	377	GLN
1	A	439	ASN
1	A	497	ASN
1	B	121	ASN
1	B	134	GLN
1	B	139	GLN
1	B	159	ASN
1	B	162	HIS
1	B	218	GLN
1	B	225	GLN
1	B	247	GLN
1	B	252	GLN
1	B	256	ASN
1	B	377	GLN
1	B	397	GLN
1	B	497	ASN

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 ⓘ

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	FMN	A	5570	-	32,33,33	1.12	2 (6%)	34,50,50	3.63	15 (44%)
3	173	A	5580	-	8,11,11	1.72	1 (12%)	10,14,14	1.84	4 (40%)
2	FMN	B	9570	-	32,33,33	1.38	5 (15%)	34,50,50	4.26	17 (50%)
3	173	B	9580	-	8,11,11	1.51	1 (12%)	10,14,14	1.98	4 (40%)

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	FMN	A	5570	-	-	0/18/18/18	0/3/3/3
3	173	A	5580	-	-	0/4/8/8	0/1/1/1
2	FMN	B	9570	-	-	0/18/18/18	0/3/3/3
3	173	B	9580	-	-	0/4/8/8	0/1/1/1

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	5580	173	C1G-C2	-4.62	1.41	1.49
2	B	9570	FMN	C10-N1	-3.58	1.29	1.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	9580	173	C1G-C2	-3.06	1.43	1.49
2	A	5570	FMN	C10-N1	-2.68	1.31	1.35
2	B	9570	FMN	C2-N1	-2.15	1.33	1.38
2	B	9570	FMN	C9A-N10	-2.14	1.35	1.38
2	B	9570	FMN	C5A-N5	2.32	1.39	1.35
2	A	5570	FMN	C4-C4A	2.71	1.46	1.41
2	B	9570	FMN	C4-C4A	3.76	1.48	1.41

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	9570	FMN	C4-C4A-C10	-9.03	114.16	119.94
2	A	5570	FMN	C4A-C10-N10	-8.04	114.68	120.52
2	B	9570	FMN	N3-C2-N1	-7.75	114.64	127.69
2	A	5570	FMN	N3-C2-N1	-6.96	115.97	127.69
2	B	9570	FMN	C4A-C10-N10	-6.52	115.78	120.52
2	A	5570	FMN	C4A-C4-N3	-6.45	115.09	123.52
2	B	9570	FMN	C4A-C4-N3	-6.12	115.52	123.52
2	A	5570	FMN	C4-C4A-C10	-4.48	117.08	119.94
2	B	9570	FMN	O3'-C3'-C2'	-3.89	98.63	108.73
2	B	9570	FMN	O2P-P-O5'	-3.16	97.50	106.72
3	B	9580	173	C1D-C1G-C2D	-2.66	114.97	118.61
2	B	9570	FMN	C7M-C7-C8	-2.48	115.39	120.73
2	B	9570	FMN	C6-C5A-C9A	-2.43	116.42	119.11
3	A	5580	173	C1D-C1G-C2D	-2.38	115.35	118.61
2	A	5570	FMN	C6-C7-C8	-2.30	115.48	119.97
2	A	5570	FMN	C8M-C8-C9	-2.20	114.11	120.33
3	A	5580	173	C1E-C1D-C1G	2.19	122.96	120.35
2	A	5570	FMN	C6-C5A-N5	2.20	121.66	118.92
3	B	9580	173	C2D-C1G-C2	2.24	125.61	120.58
3	A	5580	173	C2D-C1G-C2	2.26	125.64	120.58
2	A	5570	FMN	C7-C6-C5A	2.30	124.65	120.90
2	B	9570	FMN	C7-C6-C5A	2.35	124.74	120.90
2	A	5570	FMN	O4'-C4'-C5'	2.38	115.29	110.09
2	A	5570	FMN	O3P-P-O1P	2.56	118.99	110.63
2	A	5570	FMN	C1'-C2'-C3'	2.67	117.45	109.82
3	B	9580	173	C1E-C1D-C1G	2.68	123.54	120.35
2	A	5570	FMN	O3'-C3'-C4'	2.77	115.92	108.73
2	A	5570	FMN	O2'-C2'-C3'	3.22	117.25	108.96
3	A	5580	173	C1G-C2-C1	3.35	125.43	119.45
2	A	5570	FMN	C4A-N5-C5A	3.45	120.79	116.72
2	B	9570	FMN	O4'-C4'-C5'	3.47	117.65	110.09

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	9570	FMN	C4A-N5-C5A	3.70	121.08	116.72
2	B	9570	FMN	O3'-C3'-C4'	3.77	118.51	108.73
2	B	9570	FMN	O2'-C2'-C3'	3.91	119.01	108.96
3	B	9580	173	C1G-C2-C1	3.92	126.45	119.45
2	B	9570	FMN	C4-C4A-N5	4.12	123.71	118.70
2	B	9570	FMN	C6-C5A-N5	4.58	124.63	118.92
2	B	9570	FMN	O2'-C2'-C1'	4.69	121.52	109.93
2	A	5570	FMN	C4-N3-C2	13.55	126.47	115.16
2	B	9570	FMN	C4-N3-C2	14.73	127.45	115.16

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	5570	FMN	4	0
3	A	5580	173	1	0
2	B	9570	FMN	3	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 [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	392/511 (76%)	-0.78	2 (0%) 91 76	24, 43, 75, 88	0
1	B	389/511 (76%)	-0.59	3 (0%) 87 67	30, 54, 85, 89	0
All	All	781/1022 (76%)	-0.69	5 (0%) 90 73	24, 47, 83, 89	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	511	ALA	5.8
1	B	114	SER	4.1
1	A	114	SER	3.8
1	B	321	ALA	2.2
1	B	511	ALA	2.1

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
3	173	A	5580	11/11	0.93	0.24	2.83	57,62,64,65	0
3	173	B	9580	11/11	0.92	0.20	2.16	56,64,71,72	0
2	FMN	A	5570	31/31	0.99	0.10	-0.92	18,27,32,33	0
2	FMN	B	9570	31/31	0.99	0.10	-1.14	34,40,43,45	0

6.5 Other polymers [i](#)

There are no such residues in this entry.