



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 15, 2017 – 01:36 am GMT

PDB ID : 3AD8
Title : Heterotetrameric Sarcosine Oxidase from *Corynebacterium* sp. U-96 in complex with pyrrole 2-carboxylate
Authors : Suzuki, H.; Moriguchi, T.; Ida, K.
Deposited on : 2010-01-15
Resolution : 2.20 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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
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) : recalc28949

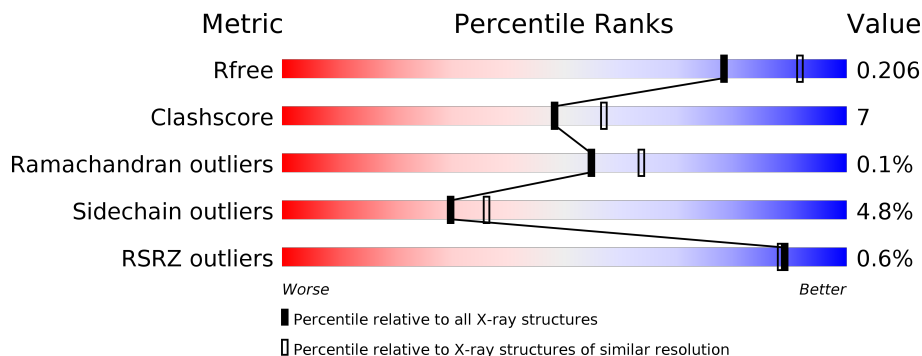
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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}	100719	4002 (2.20-2.20)
Clashscore	112137	4730 (2.20-2.20)
Ramachandran outliers	110173	4656 (2.20-2.20)
Sidechain outliers	110143	4657 (2.20-2.20)
RSRZ outliers	101464	4033 (2.20-2.20)

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	964	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 1%, green 98%);"></div> <div style="display: flex; justify-content: space-between; width: 90%; margin: 5px auto;"> 87% 11% </div> </div>
2	B	404	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 1%, green 98%);"></div> <div style="display: flex; justify-content: space-between; width: 90%; margin: 5px auto;"> 85% 12% </div> </div>
3	C	203	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, green 80%, yellow 13%, orange 5%, grey 2%);"></div> <div style="display: flex; justify-content: space-between; width: 90%; margin: 5px auto;"> 80% 13% </div> </div>
4	D	99	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, green 87%, yellow 5%, grey 8%);"></div> <div style="display: flex; justify-content: space-between; width: 90%; margin: 5px auto;"> 87% 5% 8% </div> </div>

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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
6	SO4	A	2506	-	-	-	X
6	SO4	A	2509	-	-	X	X
8	FMN	B	406	-	-	X	-
9	PYC	B	801	-	-	-	X

2 Entry composition

There are 11 unique types of molecules in this entry. The entry contains 13736 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 SARCOSINE OXIDASE ALPHA SUBUNIT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	963	Total	C	N	O	S	0	0	0
			7229	4507	1287	1413	22			

- Molecule 2 is a protein called SARCOSINE OXIDASE BETA SUBUNIT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	404	Total	C	N	O	S	0	0	0
			3108	1981	541	576	10			

- Molecule 3 is a protein called SARCOSINE OXIDASE GAMMA SUBUNIT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	195	Total	C	N	O	S	0	0	0
			1433	902	257	271	3			

- Molecule 4 is a protein called SARCOSINE OXIDASE DELTA SUBUNIT.

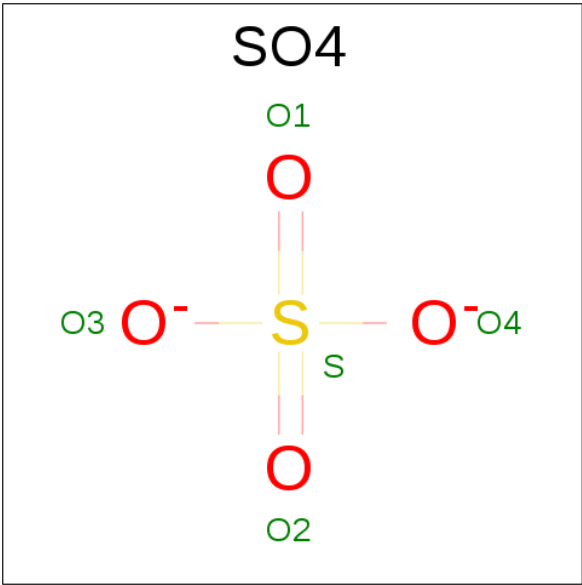
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	91	Total	C	N	O	S	0	0	0
			749	476	135	133	5			

- Molecule 5 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: C₂₁H₂₇N₇O₁₄P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

- Molecule 6 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



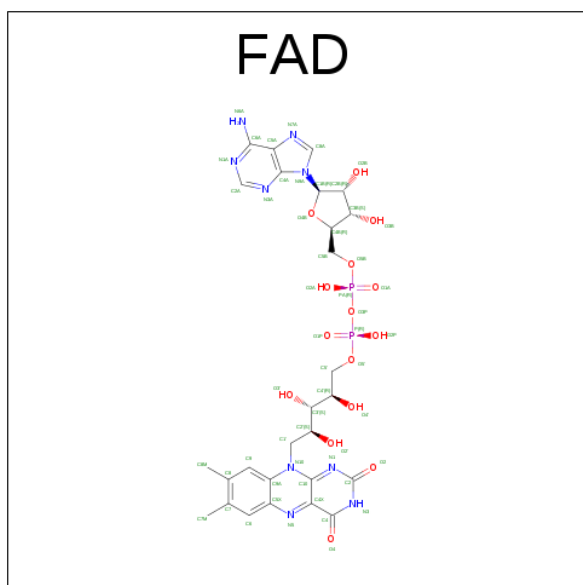
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	O	S	0	0
			5	4	1		
6	A	1	Total	O	S	0	0
			5	4	1		
6	A	1	Total	O	S	0	0
			5	4	1		

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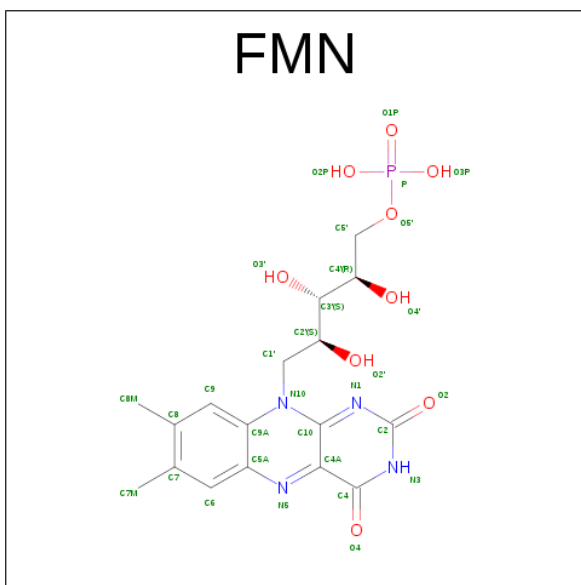
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	O	S	0	0
			5	4	1		
6	A	1	Total	O	S	0	0
			5	4	1		
6	A	1	Total	O	S	0	0
			5	4	1		
6	A	1	Total	O	S	0	0
			5	4	1		
6	B	1	Total	O	S	0	0
			5	4	1		
6	B	1	Total	O	S	0	0
			5	4	1		
6	B	1	Total	O	S	0	0
			5	4	1		
6	C	1	Total	O	S	0	0
			5	4	1		
6	D	1	Total	O	S	0	0
			5	4	1		

- Molecule 7 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).



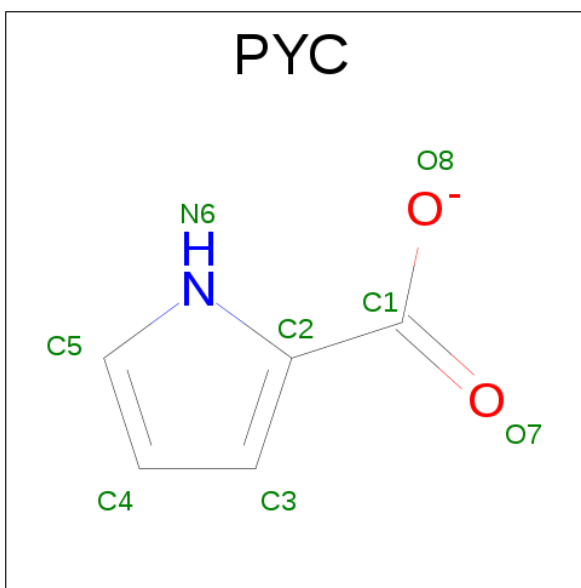
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	B	1	Total	C	N	O	P	
			53	27	9	15	2	

- Molecule 8 is FLAVIN MONONUCLEOTIDE (three-letter code: FMN) (formula: $C_{17}H_{21}N_4O_9P$).



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

- Molecule 9 is PYRROLE-2-CARBOXYLATE (three-letter code: PYC) (formula: $\text{C}_5\text{H}_4\text{NO}_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
9	B	1	Total	C	N	O	0	0
			8	5	1	2		

- Molecule 10 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	D	1	Total 1	Zn 1	0	0

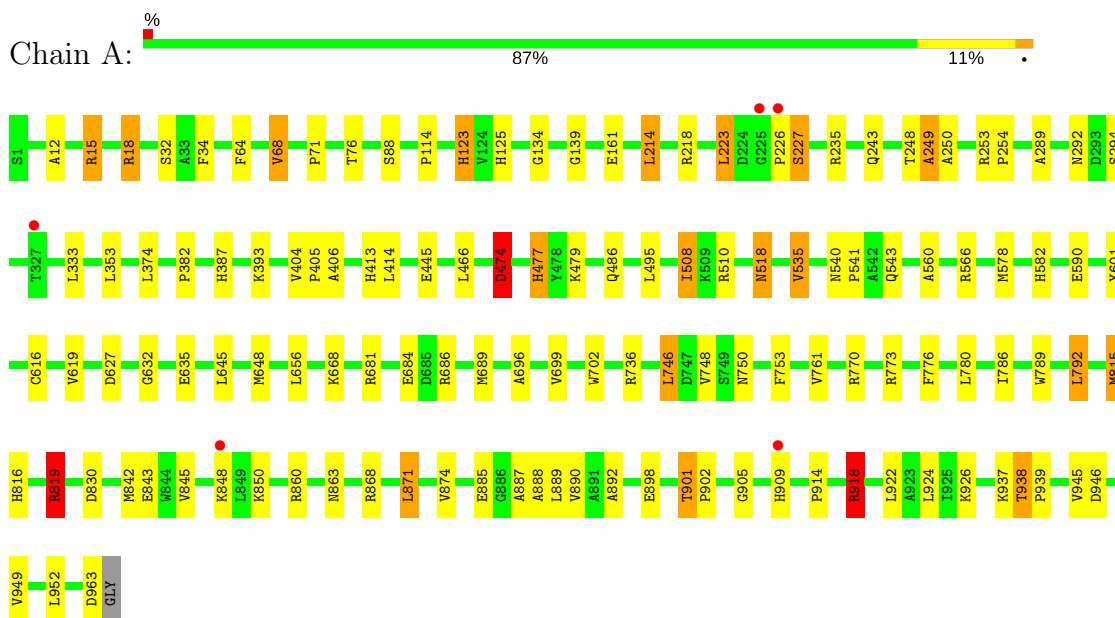
- Molecule 11 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	A	631	Total 631	O 631	0	0
11	B	189	Total 189	O 189	0	0
11	C	133	Total 133	O 133	0	0
11	D	67	Total 67	O 67	0	0

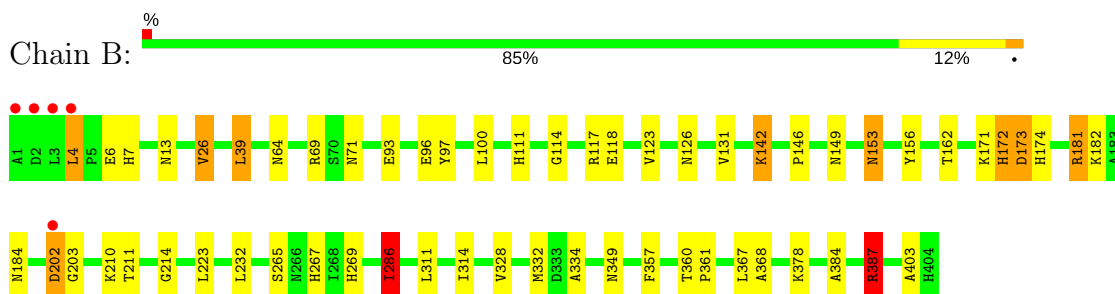
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes 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.

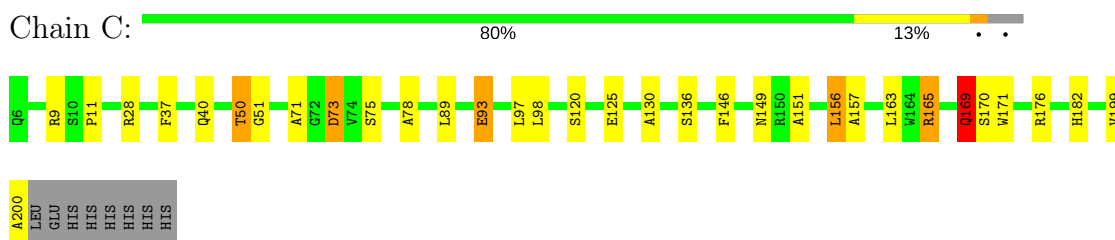
• Molecule 1: SARCOSINE OXIDASE ALPHA SUBUNIT




• Molecule 2: SARCOSINE OXIDASE BETA SUBUNIT

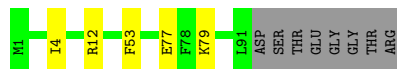


• Molecule 3: SARCOSINE OXIDASE GAMMA SUBUNIT



- Molecule 4: SARCOSINE OXIDASE DELTA SUBUNIT

Chain D:  87% 5% 8%



4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	198.80Å 198.80Å 196.82Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	61.31 – 2.20 61.31 – 2.20	Depositor EDS
% Data completeness (in resolution range)	100.0 (61.31-2.20) 100.0 (61.31-2.20)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	7.18 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.5.0072	Depositor
R, R_{free}	0.161 , 0.206 0.161 , 0.206	Depositor DCC
R_{free} test set	5799 reflections (5.28%)	DCC
Wilson B-factor (Å ²)	24.4	Xtriage
Anisotropy	0.054	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 43.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	13736	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.27% 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, 0.333 respectively for untwinned datasets, and 0.375, 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: ZN, NAD, FMN, SO4, PYC, FAD

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	1.04	5/7361 (0.1%)	0.93	19/10017 (0.2%)
2	B	1.01	2/3189 (0.1%)	0.91	6/4340 (0.1%)
3	C	1.11	3/1461 (0.2%)	1.00	6/1998 (0.3%)
4	D	1.03	1/772 (0.1%)	0.88	1/1040 (0.1%)
All	All	1.04	11/12783 (0.1%)	0.93	32/17395 (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	A	0	1
3	C	0	1
All	All	0	2

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	477	HIS	CB-CG	9.30	1.66	1.50
1	A	477	HIS	CA-CB	8.05	1.71	1.53
1	A	477	HIS	CA-C	6.84	1.70	1.52
4	D	53	PHE	CE1-CZ	6.38	1.49	1.37
1	A	249	ALA	CA-CB	-5.99	1.39	1.52
3	C	130	ALA	CA-CB	5.88	1.64	1.52
3	C	125	GLU	CB-CG	-5.67	1.41	1.52
1	A	445	GLU	CG-CD	5.64	1.60	1.51
3	C	169	GLN	CB-CG	5.41	1.67	1.52
2	B	387	ARG	CG-CD	5.31	1.65	1.51
2	B	334	ALA	CA-CB	5.17	1.63	1.52

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	387	ARG	NE-CZ-NH1	-14.34	113.13	120.30
1	A	15	ARG	NE-CZ-NH2	-12.63	113.98	120.30
1	A	819	ARG	NE-CZ-NH2	-12.61	113.99	120.30
1	A	819	ARG	NE-CZ-NH1	10.36	125.48	120.30
3	C	176	ARG	NE-CZ-NH2	-8.83	115.88	120.30
2	B	387	ARG	NE-CZ-NH2	8.37	124.49	120.30
1	A	477	HIS	CB-CA-C	8.22	126.83	110.40
1	A	918	ARG	NE-CZ-NH1	8.10	124.35	120.30
3	C	176	ARG	NE-CZ-NH1	8.08	124.34	120.30
3	C	165	ARG	NE-CZ-NH1	8.05	124.33	120.30
1	A	770	ARG	NE-CZ-NH2	-7.04	116.78	120.30
1	A	15	ARG	NE-CZ-NH1	6.93	123.77	120.30
2	B	117	ARG	NE-CZ-NH2	-6.77	116.91	120.30
1	A	214	LEU	CA-CB-CG	-6.67	99.95	115.30
3	C	28	ARG	NE-CZ-NH2	-6.57	117.02	120.30
1	A	918	ARG	NE-CZ-NH2	-6.45	117.08	120.30
1	A	819	ARG	CG-CD-NE	-6.43	98.30	111.80
3	C	28	ARG	NE-CZ-NH1	6.39	123.49	120.30
1	A	214	LEU	CB-CG-CD1	6.09	121.36	111.00
2	B	181	ARG	NE-CZ-NH2	-6.08	117.26	120.30
1	A	819	ARG	CD-NE-CZ	6.01	132.01	123.60
1	A	566	ARG	NE-CZ-NH2	-5.87	117.36	120.30
1	A	235	ARG	NE-CZ-NH2	-5.73	117.44	120.30
1	A	12	ALA	C-N-CA	-5.72	110.29	122.30
1	A	474	ASP	CB-CG-OD2	-5.68	113.19	118.30
2	B	286	ILE	CG1-CB-CG2	5.47	123.44	111.40
1	A	18	ARG	NE-CZ-NH2	-5.30	117.65	120.30
1	A	819	ARG	CB-CG-CD	5.28	125.33	111.60
4	D	12	ARG	NE-CZ-NH1	5.11	122.86	120.30
3	C	200	ALA	N-CA-C	5.11	124.80	111.00
2	B	173	ASP	CB-CG-OD1	5.09	122.88	118.30
1	A	68	VAL	CG1-CB-CG2	5.03	118.95	110.90

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	227	SER	Peptide
3	C	199	VAL	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	7229	0	7104	101	1
2	B	3108	0	3035	54	0
3	C	1433	0	1434	17	0
4	D	749	0	706	3	0
5	A	44	0	26	4	0
6	A	35	0	0	3	0
6	B	15	0	0	0	0
6	C	5	0	0	0	0
6	D	5	0	0	0	0
7	B	53	0	31	4	0
8	B	31	0	18	9	0
9	B	8	0	4	0	0
10	D	1	0	0	0	0
11	A	631	0	0	13	1
11	B	189	0	0	5	0
11	C	133	0	0	2	0
11	D	67	0	0	0	0
All	All	13736	0	12358	173	2

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 7.

All (173) 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)
2:B:172:HIS:NE2	8:B:406:FMN:C8M	1.71	1.50
1:A:830:ASP:OD1	1:A:909:HIS:CE1	1.88	1.26
1:A:696:ALA:HB3	6:A:2509:SO4:O3	1.41	1.17
2:B:384:ALA:O	2:B:387:ARG:HD3	1.44	1.12
2:B:387:ARG:HH11	2:B:387:ARG:HG3	0.98	1.11
1:A:830:ASP:CG	1:A:909:HIS:HE1	1.56	1.09
2:B:387:ARG:HH11	2:B:387:ARG:CG	1.66	1.06
1:A:885:GLU:HG2	1:A:909:HIS:HA	1.37	1.05
1:A:696:ALA:HB1	11:A:1461:HOH:O	1.57	1.04
2:B:387:ARG:NH1	2:B:387:ARG:HG3	1.76	0.98

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:736:ARG:NH2	1:A:963:ASP:O	1.99	0.95
1:A:696:ALA:CB	6:A:2509:SO4:O3	2.14	0.95
3:C:169:GLN:HE21	3:C:171:TRP:HE1	1.13	0.94
1:A:830:ASP:OD1	1:A:909:HIS:HE1	1.36	0.92
1:A:226:PRO:HB2	11:A:1256:HOH:O	1.70	0.90
3:C:169:GLN:NE2	3:C:171:TRP:HE1	1.72	0.88
1:A:15:ARG:CD	1:A:161:GLU:OE2	2.22	0.87
3:C:37:PHE:O	3:C:93:GLU:HG2	1.75	0.87
2:B:100:LEU:HD12	2:B:171:LYS:HD2	1.57	0.87
1:A:689:MET:HE2	1:A:699:VAL:HG11	1.56	0.87
1:A:15:ARG:HD2	1:A:161:GLU:OE2	1.75	0.85
1:A:868:ARG:O	1:A:926:LYS:O	1.95	0.84
1:A:689:MET:CE	1:A:699:VAL:HG11	2.08	0.83
1:A:819:ARG:HH22	1:A:909:HIS:CE1	1.98	0.81
1:A:830:ASP:CG	1:A:909:HIS:CE1	2.44	0.79
1:A:888:ALA:O	1:A:938:THR:HG23	1.83	0.79
2:B:172:HIS:HD1	2:B:173:ASP:H	1.32	0.78
1:A:249:ALA:HB1	11:A:978:HOH:O	1.83	0.78
1:A:830:ASP:OD2	1:A:909:HIS:HE1	1.66	0.78
1:A:510:ARG:HH21	8:B:406:FMN:H5'1	1.50	0.76
2:B:172:HIS:CD2	8:B:406:FMN:C8M	2.68	0.76
2:B:387:ARG:CG	2:B:387:ARG:NH1	2.37	0.75
1:A:892:ALA:HA	1:A:937:LYS:HD2	1.68	0.75
3:C:50:THR:HG22	11:C:482:HOH:O	1.87	0.74
2:B:269:HIS:CE1	2:B:403:ALA:H	2.06	0.73
1:A:292:ASN:HD22	1:A:294:SER:H	1.34	0.73
1:A:736:ARG:NH1	11:A:1486:HOH:O	2.21	0.72
1:A:816:HIS:O	1:A:819:ARG:HD3	1.89	0.72
2:B:202:ASP:OD1	2:B:203:GLY:N	2.22	0.72
1:A:909:HIS:NE2	1:A:922:LEU:HD12	2.05	0.72
1:A:901:THR:CG2	11:A:1118:HOH:O	2.39	0.71
1:A:15:ARG:HD3	1:A:161:GLU:OE2	1.89	0.70
1:A:510:ARG:HH21	8:B:406:FMN:C5'	2.04	0.70
1:A:918:ARG:HG3	1:A:918:ARG:HH11	1.57	0.70
2:B:153:ASN:H	2:B:153:ASN:HD22	1.38	0.70
1:A:696:ALA:HB3	6:A:2509:SO4:S	2.31	0.69
2:B:93:GLU:HG2	11:B:655:HOH:O	1.92	0.69
3:C:73:ASP:OD2	3:C:75:SER:HB2	1.92	0.69
1:A:292:ASN:ND2	1:A:294:SER:H	1.90	0.68
1:A:860:ARG:HH21	1:A:863:ASN:HD21	1.42	0.67
1:A:819:ARG:NH2	1:A:909:HIS:NE2	2.42	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:125:HIS:HD2	11:A:1191:HOH:O	1.77	0.66
1:A:736:ARG:HG2	1:A:748:VAL:HG23	1.78	0.65
1:A:937:LYS:NZ	1:A:946:ASP:OD2	2.29	0.65
3:C:37:PHE:O	3:C:93:GLU:CG	2.44	0.65
2:B:210:LYS:HE3	2:B:211:THR:O	1.97	0.65
2:B:149:ASN:HD21	2:B:153:ASN:HD21	1.45	0.64
1:A:830:ASP:OD2	1:A:909:HIS:CE1	2.48	0.64
1:A:243:GLN:OE1	1:A:413:HIS:HE1	1.80	0.64
2:B:269:HIS:HE1	2:B:403:ALA:H	1.46	0.64
1:A:901:THR:HG23	11:A:1118:HOH:O	1.98	0.63
1:A:223:LEU:HD23	1:A:226:PRO:HB3	1.80	0.63
2:B:171:LYS:HE2	2:B:174:HIS:CD2	2.34	0.62
1:A:632:GLY:HA3	1:A:696:ALA:HB2	1.81	0.61
3:C:157:ALA:O	3:C:182:HIS:HE1	1.84	0.60
1:A:750:ASN:HD22	1:A:773:ARG:HH12	1.49	0.60
2:B:142:LYS:HE2	2:B:146:PRO:O	2.02	0.59
1:A:945:VAL:HG13	11:A:1127:HOH:O	2.02	0.59
1:A:816:HIS:HA	1:A:819:ARG:HD2	1.85	0.59
2:B:172:HIS:HD1	2:B:173:ASP:N	1.98	0.58
1:A:842:MET:O	1:A:845:VAL:HG12	2.03	0.58
11:A:1227:HOH:O	3:C:182:HIS:HD2	1.86	0.58
1:A:582:HIS:CD2	1:A:601:TYR:OH	2.57	0.58
1:A:382:PRO:HG2	1:A:404:VAL:HG12	1.86	0.58
2:B:114:GLY:O	2:B:118:GLU:HG2	2.03	0.58
2:B:111:HIS:HD2	2:B:156:TYR:O	1.87	0.58
1:A:635:GLU:OE2	1:A:686:ARG:NH1	2.32	0.57
1:A:474:ASP:OD2	1:A:477:HIS:CD2	2.58	0.56
1:A:540:ASN:O	1:A:543:GLN:HB2	2.05	0.56
2:B:7:HIS:HE1	11:B:426:HOH:O	1.86	0.56
2:B:172:HIS:CE1	8:B:406:FMN:C8M	2.76	0.56
2:B:172:HIS:NE2	8:B:406:FMN:C8	2.63	0.55
1:A:510:ARG:NH2	8:B:406:FMN:C5'	2.70	0.55
1:A:888:ALA:O	1:A:938:THR:CG2	2.51	0.55
1:A:249:ALA:HB2	5:A:965:NAD:O3	2.07	0.55
1:A:750:ASN:ND2	1:A:773:ARG:HH12	2.06	0.54
2:B:64:ASN:HA	7:B:405:FAD:C6	2.37	0.54
1:A:746:LEU:HD13	1:A:748:VAL:HG12	1.89	0.54
2:B:39:LEU:HD13	2:B:368:ALA:HA	1.91	0.53
2:B:286:ILE:HD11	2:B:328:VAL:HG21	1.90	0.53
1:A:510:ARG:NH2	8:B:406:FMN:H5'1	2.22	0.53
2:B:360:THR:HB	2:B:361:PRO:HD3	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:249:ALA:HB2	5:A:965:NAD:PA	2.48	0.52
2:B:123:VAL:HG23	2:B:162:THR:HG22	1.92	0.52
1:A:945:VAL:CG1	11:A:1127:HOH:O	2.56	0.52
1:A:736:ARG:HG2	1:A:748:VAL:CG2	2.40	0.52
2:B:4:LEU:HD22	2:B:96:GLU:HA	1.91	0.52
3:C:11:PRO:HG2	3:C:120:SER:HB3	1.91	0.52
1:A:249:ALA:CB	5:A:965:NAD:O2A	2.58	0.52
2:B:118:GLU:OE2	11:B:802:HOH:O	2.19	0.51
2:B:26:VAL:HG13	2:B:223:LEU:HD23	1.92	0.51
1:A:114:PRO:HA	2:B:314:ILE:HG13	1.91	0.51
1:A:508:ILE:N	1:A:508:ILE:HD13	2.26	0.51
1:A:495:LEU:HD21	1:A:535:VAL:HG22	1.92	0.51
2:B:126:ASN:HB3	2:B:131:VAL:HG22	1.93	0.51
1:A:248:THR:O	1:A:249:ALA:HB3	2.12	0.50
1:A:486:GLN:HE22	1:A:518:ASN:ND2	2.10	0.49
1:A:871:LEU:HD13	1:A:952:LEU:HD11	1.93	0.49
4:D:77:GLU:CD	4:D:79:LYS:NZ	2.66	0.49
1:A:582:HIS:HE1	1:A:627:ASP:OD2	1.96	0.48
1:A:249:ALA:HB2	5:A:965:NAD:O2A	2.14	0.48
2:B:13:ASN:ND2	2:B:184:ASN:HD21	2.13	0.47
2:B:265:SER:O	2:B:269:HIS:HD2	1.98	0.47
1:A:689:MET:HE1	1:A:699:VAL:HG11	1.90	0.47
1:A:938:THR:HG22	1:A:939:PRO:HD2	1.97	0.47
1:A:736:ARG:HB2	1:A:780:LEU:HD21	1.98	0.46
1:A:64:PHE:HB3	1:A:71:PRO:HD2	1.97	0.46
3:C:169:GLN:NE2	3:C:171:TRP:NE1	2.53	0.46
1:A:901:THR:HG22	11:A:1118:HOH:O	2.10	0.46
2:B:111:HIS:CE1	2:B:265:SER:OG	2.69	0.46
3:C:40:GLN:HA	3:C:89:LEU:O	2.14	0.46
1:A:689:MET:HE2	1:A:699:VAL:CG1	2.37	0.46
2:B:265:SER:O	2:B:269:HIS:HA	2.16	0.46
1:A:510:ARG:NH2	8:B:406:FMN:H5'2	2.31	0.46
2:B:349:ASN:HB2	2:B:367:LEU:HD22	1.98	0.46
1:A:253:ARG:HB2	1:A:254:PRO:HD2	1.98	0.45
1:A:289:ALA:HB2	1:A:374:LEU:HD11	1.97	0.45
1:A:405:PRO:HB2	1:A:414:LEU:HD13	1.98	0.45
1:A:905:GLY:HA3	1:A:924:LEU:O	2.16	0.45
3:C:136:SER:OG	3:C:156:LEU:HD22	2.17	0.45
1:A:898:GLU:HG3	11:A:1298:HOH:O	2.16	0.45
2:B:69:ARG:HD2	2:B:71:ASN:OD1	2.14	0.45
1:A:815:MET:HE2	1:A:816:HIS:HD2	1.82	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:118:GLU:OE1	11:B:809:HOH:O	2.20	0.45
1:A:901:THR:HA	1:A:902:PRO:HD3	1.77	0.45
1:A:387:HIS:HD2	11:A:968:HOH:O	1.98	0.45
1:A:486:GLN:HE22	1:A:518:ASN:HD22	1.64	0.45
2:B:96:GLU:O	2:B:181:ARG:NH2	2.50	0.45
4:D:77:GLU:OE2	4:D:79:LYS:NZ	2.49	0.44
1:A:123:HIS:HE1	1:A:560:ALA:O	2.01	0.44
2:B:210:LYS:NZ	2:B:214:GLY:HA2	2.33	0.44
3:C:51:GLY:HA3	11:C:223:HOH:O	2.17	0.44
1:A:668:LYS:HD3	1:A:702:TRP:CH2	2.53	0.44
1:A:819:ARG:NH2	1:A:830:ASP:OD2	2.51	0.44
1:A:134:GLY:O	1:A:139:GLY:HA3	2.18	0.44
1:A:860:ARG:NH2	1:A:863:ASN:HD21	2.13	0.44
2:B:153:ASN:HD22	2:B:153:ASN:N	2.06	0.43
1:A:387:HIS:HE1	1:A:393:LYS:O	2.01	0.43
1:A:479:LYS:HE2	1:A:479:LYS:HB3	1.87	0.43
4:D:77:GLU:CD	4:D:79:LYS:HZ2	2.22	0.43
1:A:249:ALA:HB1	1:A:250:ALA:H	1.66	0.43
1:A:648:MET:HE2	1:A:648:MET:HA	2.01	0.43
1:A:689:MET:CE	1:A:699:VAL:CG1	2.88	0.43
2:B:149:ASN:HD21	2:B:153:ASN:ND2	2.13	0.43
1:A:616:CYS:HB2	1:A:914:PRO:HG2	2.01	0.43
2:B:39:LEU:HD13	2:B:368:ALA:CA	2.49	0.42
1:A:776:PHE:CE2	1:A:819:ARG:HG3	2.53	0.42
2:B:111:HIS:HE1	2:B:265:SER:OG	2.02	0.42
2:B:182:LYS:HD3	2:B:182:LYS:HA	1.80	0.42
3:C:71:ALA:O	3:C:78:ALA:HA	2.20	0.42
3:C:151:ALA:HA	3:C:163:LEU:O	2.20	0.42
2:B:210:LYS:HD2	2:B:211:THR:H	1.85	0.41
2:B:6:GLU:HB2	11:B:790:HOH:O	2.20	0.41
2:B:357:PHE:HB3	7:B:405:FAD:C2	2.50	0.41
2:B:64:ASN:HA	7:B:405:FAD:C5X	2.51	0.41
1:A:18:ARG:HA	1:A:34:PHE:CD1	2.55	0.41
1:A:792:LEU:HD12	3:C:9:ARG:HD3	2.03	0.41
2:B:64:ASN:HB2	7:B:405:FAD:C4X	2.50	0.41
1:A:887:ALA:HB1	1:A:938:THR:HG21	2.03	0.41
1:A:656:LEU:HD23	1:A:681:ARG:HB2	2.04	0.40
1:A:76:THR:HA	1:A:88:SER:HA	2.02	0.40
2:B:332:MET:HA	2:B:332:MET:CE	2.52	0.40
3:C:149:ASN:N	3:C:165:ARG:O	2.50	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the sym-

metry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:477:HIS:NE2	1:A:477:HIS:NE2[11_555]	1.27	0.93
11:A:1080:HOH:O	11:A:1080:HOH:O[9_555]	1.63	0.57

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	961/964 (100%)	941 (98%)	18 (2%)	2 (0%)	51	58
2	B	402/404 (100%)	387 (96%)	15 (4%)	0	100	100
3	C	193/203 (95%)	187 (97%)	6 (3%)	0	100	100
4	D	89/99 (90%)	87 (98%)	2 (2%)	0	100	100
All	All	1645/1670 (98%)	1602 (97%)	41 (2%)	2 (0%)	55	63

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	227	SER
1	A	406	ALA

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	747/747 (100%)	709 (95%)	38 (5%)	28	33

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	B	319/319 (100%)	305 (96%)	14 (4%)	33	40
3	C	143/151 (95%)	134 (94%)	9 (6%)	21	23
4	D	75/81 (93%)	74 (99%)	1 (1%)	73	85
All	All	1284/1298 (99%)	1222 (95%)	62 (5%)	30	36

All (62) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	32	SER
1	A	68	VAL
1	A	123	HIS
1	A	214	LEU
1	A	218	ARG
1	A	223	LEU
1	A	333	LEU
1	A	353	LEU
1	A	466	LEU
1	A	474	ASP
1	A	508	ILE
1	A	518	ASN
1	A	535	VAL
1	A	541	PRO
1	A	578	MET
1	A	590	GLU
1	A	619	VAL
1	A	645	LEU
1	A	684	GLU
1	A	746	LEU
1	A	753	PHE
1	A	761	VAL
1	A	786	ILE
1	A	789	TRP
1	A	792	LEU
1	A	815	MET
1	A	819	ARG
1	A	843	GLU
1	A	848	LYS
1	A	850	LYS
1	A	871	LEU
1	A	874	VAL
1	A	889	LEU

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Mol	Chain	Res	Type
1	A	890	VAL
1	A	901	THR
1	A	918	ARG
1	A	938	THR
1	A	949	VAL
2	B	4	LEU
2	B	26	VAL
2	B	39	LEU
2	B	97	TYR
2	B	142	LYS
2	B	153	ASN
2	B	172	HIS
2	B	202	ASP
2	B	232	LEU
2	B	267	HIS
2	B	286	ILE
2	B	311	LEU
2	B	378	LYS
2	B	387	ARG
3	C	50	THR
3	C	73	ASP
3	C	93	GLU
3	C	97	LEU
3	C	98	LEU
3	C	146	PHE
3	C	156	LEU
3	C	169	GLN
3	C	170	SER
4	D	4	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (30) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	10	GLN
1	A	82	GLN
1	A	123	HIS
1	A	125	HIS
1	A	181	GLN
1	A	238	HIS
1	A	292	ASN
1	A	387	HIS
1	A	413	HIS

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Mol	Chain	Res	Type
1	A	477	HIS
1	A	518	ASN
1	A	540	ASN
1	A	582	HIS
1	A	750	ASN
1	A	816	HIS
1	A	863	ASN
1	A	912	ASN
2	B	7	HIS
2	B	13	ASN
2	B	14	ASN
2	B	111	HIS
2	B	153	ASN
2	B	267	HIS
2	B	269	HIS
2	B	344	GLN
2	B	369	HIS
3	C	6	GLN
3	C	158	ASN
3	C	169	GLN
3	C	182	HIS

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](#)

Of 17 ligands modelled in this entry, 1 is monoatomic - leaving 16 for Mogul analysis.

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
6	SO4	A	2500	-	4,4,4	0.12	0	6,6,6	0.26	0
6	SO4	A	2501	-	4,4,4	0.36	0	6,6,6	0.56	0
6	SO4	A	2502	-	4,4,4	0.28	0	6,6,6	0.27	0
6	SO4	A	2504	-	4,4,4	0.11	0	6,6,6	0.33	0
6	SO4	A	2506	-	4,4,4	0.34	0	6,6,6	0.32	0
6	SO4	A	2507	-	4,4,4	0.22	0	6,6,6	0.44	0
6	SO4	A	2509	-	4,4,4	0.42	0	6,6,6	0.43	0
5	NAD	A	965	-	41,48,48	1.41	4 (9%)	43,73,73	1.69	4 (9%)
6	SO4	B	2505	-	4,4,4	0.21	0	6,6,6	0.24	0
6	SO4	B	2510	-	4,4,4	0.50	0	6,6,6	0.62	0
6	SO4	B	2511	-	4,4,4	0.23	0	6,6,6	0.44	0
7	FAD	B	405	-	51,58,58	1.83	7 (13%)	54,89,89	1.94	12 (22%)
8	FMN	B	406	-	31,33,33	2.21	11 (35%)	38,50,50	2.49	12 (31%)
9	PYC	B	801	-	5,8,8	2.95	1 (20%)	4,10,10	1.27	0
6	SO4	C	2508	-	4,4,4	0.66	0	6,6,6	0.27	0
6	SO4	D	2503	-	4,4,4	0.22	0	6,6,6	0.19	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
6	SO4	A	2500	-	-	0/0/0/0	0/0/0/0
6	SO4	A	2501	-	-	0/0/0/0	0/0/0/0
6	SO4	A	2502	-	-	0/0/0/0	0/0/0/0
6	SO4	A	2504	-	-	0/0/0/0	0/0/0/0
6	SO4	A	2506	-	-	0/0/0/0	0/0/0/0
6	SO4	A	2507	-	-	0/0/0/0	0/0/0/0
6	SO4	A	2509	-	-	0/0/0/0	0/0/0/0
5	NAD	A	965	-	-	0/22/62/62	0/5/5/5
6	SO4	B	2505	-	-	0/0/0/0	0/0/0/0
6	SO4	B	2510	-	-	0/0/0/0	0/0/0/0
6	SO4	B	2511	-	-	0/0/0/0	0/0/0/0
7	FAD	B	405	-	-	0/28/50/50	0/6/6/6
8	FMN	B	406	-	-	0/16/18/18	0/3/3/3
9	PYC	B	801	-	-	0/0/4/4	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	SO4	C	2508	-	-	0/0/0/0	0/0/0/0
6	SO4	D	2503	-	-	0/0/0/0	0/0/0/0

All (23) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	B	801	PYC	C3-C2	-6.07	1.32	1.40
8	B	406	FMN	O5'-C5'	-6.05	1.20	1.44
5	A	965	NAD	C2B-C1B	-3.06	1.48	1.53
8	B	406	FMN	C9A-C5A	-3.06	1.36	1.42
5	A	965	NAD	C5A-N7A	-2.45	1.31	1.39
8	B	406	FMN	C4A-C10	-2.21	1.37	1.41
8	B	406	FMN	C4-C4A	-2.17	1.37	1.41
8	B	406	FMN	O3'-C3'	2.05	1.47	1.43
7	B	405	FAD	O3B-C3B	2.17	1.47	1.43
7	B	405	FAD	C9-C8	2.23	1.43	1.37
8	B	406	FMN	C7M-C7	2.25	1.55	1.51
8	B	406	FMN	C4'-C3'	2.29	1.58	1.53
7	B	405	FAD	C5X-N5	2.41	1.39	1.35
5	A	965	NAD	C5B-C4B	2.43	1.59	1.51
8	B	406	FMN	C8M-C8	3.07	1.57	1.51
8	B	406	FMN	C4A-N5	3.50	1.38	1.33
7	B	405	FAD	C9A-N10	4.08	1.44	1.38
8	B	406	FMN	C10-N1	4.21	1.39	1.33
7	B	405	FAD	C4-N3	4.49	1.41	1.33
7	B	405	FAD	C10-N1	4.57	1.39	1.33
8	B	406	FMN	C4-N3	4.71	1.41	1.33
5	A	965	NAD	O7N-C7N	5.88	1.36	1.24
7	B	405	FAD	C4X-N5	7.86	1.44	1.33

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	965	NAD	N3A-C2A-N1A	-7.59	122.25	128.86
8	B	406	FMN	O4'-C4'-C5'	-6.23	96.13	110.00
7	B	405	FAD	N3A-C2A-N1A	-6.20	123.45	128.86
7	B	405	FAD	C4B-O4B-C1B	-5.56	103.85	109.77
8	B	406	FMN	C4A-C4-N3	-3.66	118.27	123.48
5	A	965	NAD	C5B-C4B-C3B	-3.44	102.17	115.29
5	A	965	NAD	C4A-C5A-N7A	-3.30	106.22	109.41
8	B	406	FMN	C4A-C10-N10	-3.24	118.27	120.52
8	B	406	FMN	O2P-P-O5'	-2.74	99.45	106.73

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	B	406	FMN	C9A-C5A-N5	-2.53	118.47	122.24
7	B	405	FAD	C4A-C5A-N7A	-2.40	107.09	109.41
7	B	405	FAD	C9A-C5X-N5	-2.34	118.75	122.24
7	B	405	FAD	C4X-C4-N3	-2.07	120.53	123.48
7	B	405	FAD	C4X-C10-N10	-2.04	119.10	120.52
8	B	406	FMN	P-O5'-C5'	2.28	124.58	118.30
5	A	965	NAD	C2N-C3N-C4N	2.43	121.03	118.26
7	B	405	FAD	C4-C4X-N5	2.46	121.38	118.68
7	B	405	FAD	C5X-C9A-N10	2.51	119.52	117.66
7	B	405	FAD	C6-C5X-C9A	2.59	122.37	119.00
8	B	406	FMN	O5'-C5'-C4'	3.10	117.63	109.36
8	B	406	FMN	C4A-N5-C5A	3.22	120.16	116.76
8	B	406	FMN	C4'-C3'-C2'	3.33	120.58	113.41
8	B	406	FMN	C5A-C9A-N10	3.59	120.32	117.66
7	B	405	FAD	C4-N3-C2	4.50	119.10	115.16
7	B	405	FAD	C1'-N10-C9A	4.84	122.79	118.35
7	B	405	FAD	C4X-N5-C5X	4.93	121.96	116.76
8	B	406	FMN	O5'-P-O1P	5.21	121.09	106.47
8	B	406	FMN	C4-N3-C2	7.46	121.69	115.16

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 20 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	2509	SO4	3	0
5	A	965	NAD	4	0
7	B	405	FAD	4	0
8	B	406	FMN	9	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	963/964 (99%)	-0.62	5 (0%) 90 90	11, 24, 41, 59	0
2	B	404/404 (100%)	-0.55	5 (1%) 79 77	13, 25, 39, 84	0
3	C	195/203 (96%)	-0.74	0 100 100	14, 22, 42, 49	0
4	D	91/99 (91%)	-0.80	0 100 100	15, 22, 42, 54	0
All	All	1653/1670 (98%)	-0.63	10 (0%) 89 88	11, 24, 41, 84	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	2	ASP	8.3
2	B	1	ALA	7.4
2	B	3	LEU	7.3
1	A	909	HIS	3.8
1	A	226	PRO	2.8
2	B	202	ASP	2.7
2	B	4	LEU	2.4
1	A	327	THR	2.4
1	A	848	LYS	2.1
1	A	225	GLY	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, 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(Å ²)	Q<0.9
6	SO4	A	2506	5/5	0.93	0.27	7.66	65,65,67,68	0
9	PYC	B	801	8/8	0.96	0.14	4.42	44,44,46,48	0
6	SO4	A	2509	5/5	0.98	0.13	3.15	44,46,49,52	0
6	SO4	B	2510	5/5	0.94	0.23	1.94	52,53,57,57	0
6	SO4	C	2508	5/5	0.99	0.13	1.24	30,33,35,35	0
7	FAD	B	405	53/53	0.97	0.10	0.76	11,22,27,28	0
6	SO4	A	2500	5/5	0.99	0.13	0.59	65,66,67,68	0
10	ZN	D	100	1/1	1.00	0.08	0.24	19,19,19,19	0
6	SO4	B	2511	5/5	0.90	0.11	0.22	73,73,75,76	0
8	FMN	B	406	31/31	0.98	0.09	0.02	11,15,19,23	0
5	NAD	A	965	44/44	0.99	0.09	-0.38	8,14,19,25	0
6	SO4	A	2501	5/5	0.97	0.12	-0.47	46,47,50,51	0
6	SO4	D	2503	5/5	0.94	0.13	-	82,82,83,83	0
6	SO4	B	2505	5/5	0.97	0.20	-	56,57,58,60	0
6	SO4	A	2504	5/5	1.00	0.08	-	29,30,35,37	0
6	SO4	A	2502	5/5	0.95	0.29	-	76,76,77,78	0
6	SO4	A	2507	5/5	0.96	0.20	-	59,60,62,64	0

6.5 Other polymers [i](#)

There are no such residues in this entry.