



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:16 PM GMT

PDB ID : 4OCG
Title : Structure of the Shewanella loihica PV-4 NADH-dependent persulfide reductase F161A Mutant
Authors : Lee, K.-H.; Sazinsky, M.H.; Crane, E.J.
Deposited on : 2014-01-09
Resolution : 2.75 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

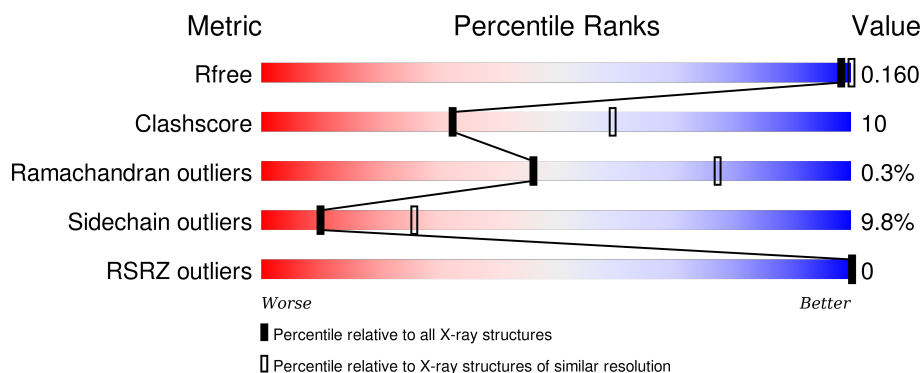
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION



The reported resolution of this entry is 2.75 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	3340 (2.80-2.72)
Clashscore	102246	3829 (2.80-2.72)
Ramachandran outliers	100387	3767 (2.80-2.72)
Sidechain outliers	100360	3770 (2.80-2.72)
RSRZ outliers	91569	3352 (2.80-2.72)

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	574	 75% 20% . .
1	B	574	 73% 22% . .

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 8950 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 FAD-dependent pyridine nucleotide-disulphide oxidoreductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	565	Total	C	N	O	S	0	0	0
			4292	2690	763	819	20			
1	B	565	Total	C	N	O	S	0	0	0
			4291	2690	762	819	20			

There are 18 discrepancies between the modelled and reference sequences:

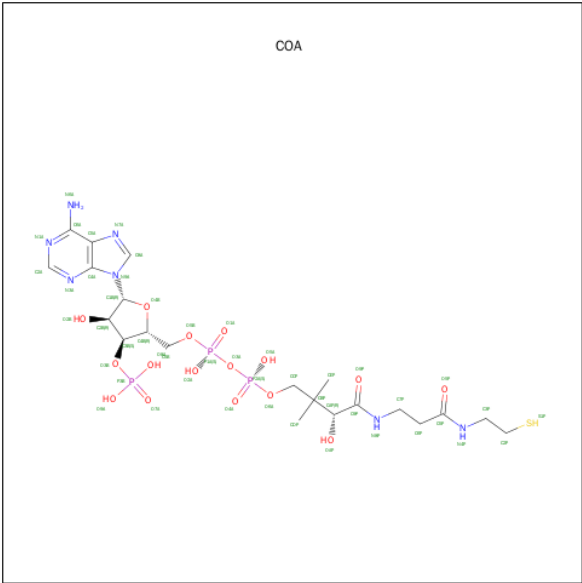
Chain	Residue	Modelled	Actual	Comment	Reference
A	161	ALA	PHE	ENGINEERED MUTATION	UNP A3QAV3
A	567	LEU	-	EXPRESSION TAG	UNP A3QAV3
A	568	GLU	-	EXPRESSION TAG	UNP A3QAV3
A	569	HIS	-	EXPRESSION TAG	UNP A3QAV3
A	570	HIS	-	EXPRESSION TAG	UNP A3QAV3
A	571	HIS	-	EXPRESSION TAG	UNP A3QAV3
A	572	HIS	-	EXPRESSION TAG	UNP A3QAV3
A	573	HIS	-	EXPRESSION TAG	UNP A3QAV3
A	574	HIS	-	EXPRESSION TAG	UNP A3QAV3
B	161	ALA	PHE	ENGINEERED MUTATION	UNP A3QAV3
B	567	LEU	-	EXPRESSION TAG	UNP A3QAV3
B	568	GLU	-	EXPRESSION TAG	UNP A3QAV3
B	569	HIS	-	EXPRESSION TAG	UNP A3QAV3
B	570	HIS	-	EXPRESSION TAG	UNP A3QAV3
B	571	HIS	-	EXPRESSION TAG	UNP A3QAV3
B	572	HIS	-	EXPRESSION TAG	UNP A3QAV3
B	573	HIS	-	EXPRESSION TAG	UNP A3QAV3
B	574	HIS	-	EXPRESSION TAG	UNP A3QAV3

- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 3 is COENZYME A (three-letter code: COA) (formula: C₂₁H₃₆N₇O₁₆P₃S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
3	A	1	Total	C	N	O	P	S	0	0
			48	21	7	16	3	1		
3	B	1	Total	C	N	O	P	S	0	0
			48	21	7	16	3	1		

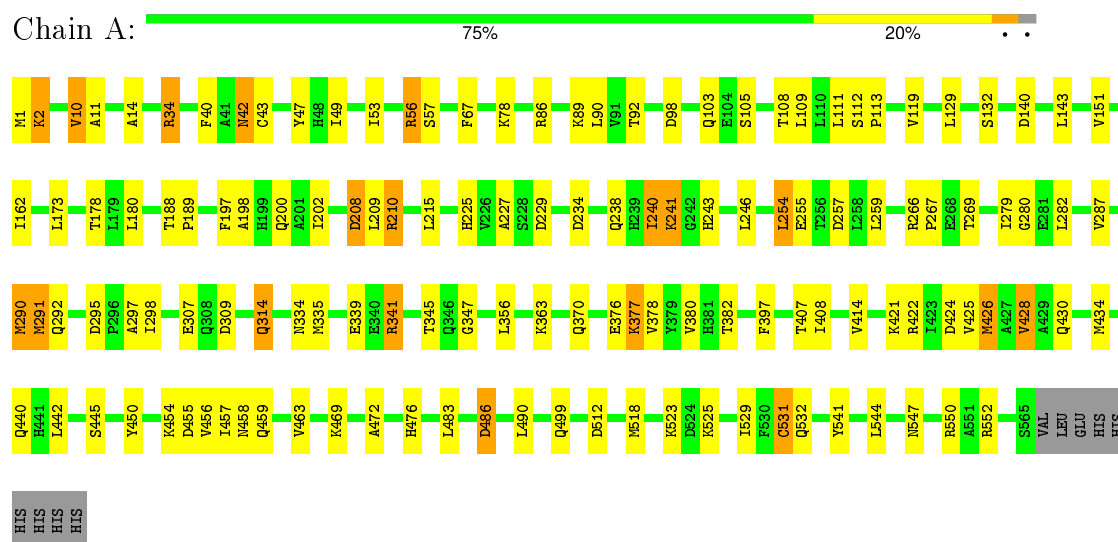
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	83	Total 83	O 83	0	0
4	B	82	Total 82	O 82	0	0

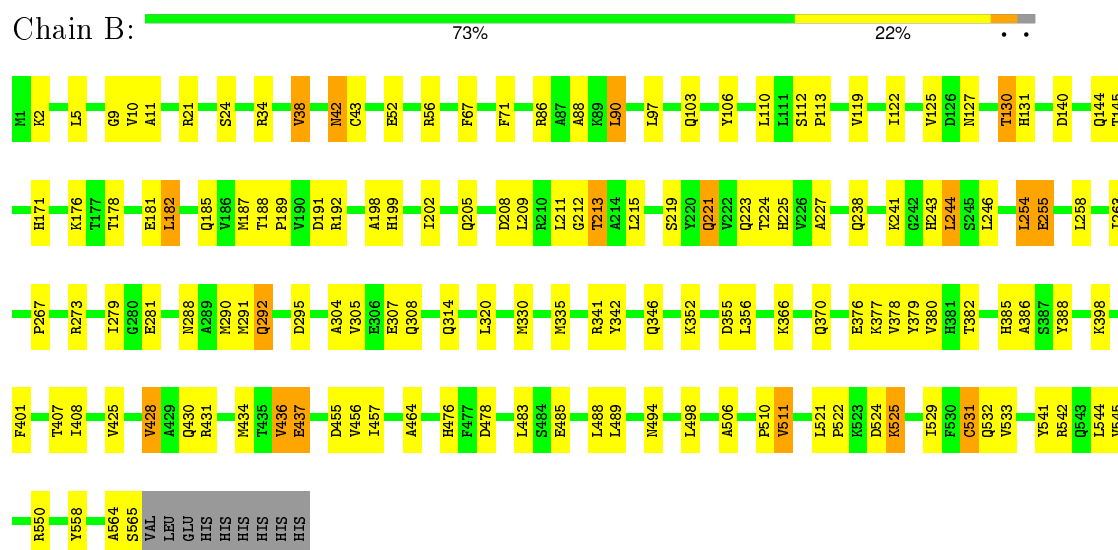
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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: FAD-dependent pyridine nucleotide-disulphide oxidoreductase



- Molecule 1: FAD-dependent pyridine nucleotide-disulphide oxidoreductase



4 Data and refinement statistics

Property	Value	Source
Space group	P 3	Depositor
Cell constants a, b, c, α , β , γ	133.71 Å 133.71 Å 84.32 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	50.00 – 2.75 35.10 – 2.75	Depositor EDS
% Data completeness (in resolution range)	96.1 (50.00-2.75) 96.3 (35.10-2.75)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	9.73 (at 2.76 Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.122 , 0.173 0.127 , 0.160	Depositor DCC
R_{free} test set	2138 reflections (5.33%)	DCC
Wilson B-factor (Å ²)	30.5	Xtriage
Anisotropy	1.495	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 3.4	EDS
Estimated twinning fraction	0.629 for H, K, L 0.371 for -H, H+K, -L 0.117 for -h,-k,l 0.117 for h,-h-k,-l 0.357 for -k,-h,-l	Xtriage
Reported twinning fraction	0.629 for H, K, L 0.371 for -H, H+K, -L	Depositor
L-test for twinning ²	$\langle L \rangle = 0.39$, $\langle L^2 \rangle = 0.21$	Xtriage
Outliers	0 of 42327 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	8950	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.63% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² Theoretical 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: COA, 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	1.01	0/4361	0.91	4/5904 (0.1%)
1	B	1.06	3/4360 (0.1%)	0.93	5/5902 (0.1%)
All	All	1.03	3/8721 (0.0%)	0.92	9/11806 (0.1%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	305	VAL	CB-CG1	-5.55	1.41	1.52
1	B	558	TYR	CD2-CE2	-5.24	1.31	1.39
1	B	428	VAL	CB-CG2	-5.02	1.42	1.52

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	478	ASP	CB-CG-OD1	5.89	123.60	118.30
1	B	244	LEU	CB-CG-CD1	-5.54	101.58	111.00
1	B	550	ARG	NE-CZ-NH1	-5.44	117.58	120.30
1	A	426	MET	CG-SD-CE	-5.27	91.77	100.20
1	B	511	VAL	CB-CA-C	-5.21	101.51	111.40
1	A	290	MET	N-CA-C	5.13	124.86	111.00
1	A	98	ASP	CB-CG-OD2	5.09	122.88	118.30
1	B	355	ASP	CB-CG-OD1	5.06	122.86	118.30
1	A	89	LYS	C-N-CA	5.02	134.25	121.70

There are no chirality outliers.

There are no planarity outliers.

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	4292	0	4304	94	1
1	B	4291	0	4300	88	0
2	A	53	0	31	5	0
2	B	53	0	31	3	0
3	A	48	0	32	4	0
3	B	48	0	32	7	0
4	A	83	0	0	3	0
4	B	82	0	0	2	0
All	All	8950	0	8730	183	1

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 10.

All (183) 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:341:ARG:HH11	1:A:341:ARG:HG2	1.19	1.03
1:B:346:GLN:HE21	1:B:430:GLN:HE21	0.96	0.93
1:A:382:THR:OG1	1:A:456:VAL:HG22	1.68	0.92
1:B:43:CYS:HG	3:B:901:COA:HS1	1.02	0.92
1:B:346:GLN:HE21	1:B:430:GLN:NE2	1.70	0.90
1:A:314:GLN:HA	1:A:314:GLN:HE21	1.42	0.82
1:A:290:MET:O	1:A:334:ASN:ND2	2.13	0.81
1:B:243:HIS:CE1	1:B:255:GLU:HG3	2.16	0.80
1:A:425:VAL:HG13	1:B:428:VAL:HG21	1.63	0.79
1:A:531:CYS:SG	1:A:532:GLN:N	2.56	0.79
1:A:376:GLU:HG3	1:A:377:LYS:N	1.99	0.78
1:A:377:LYS:HD3	1:A:476:HIS:CE1	2.18	0.78
1:A:43:CYS:SG	3:A:901:COA:S1P	2.58	0.77
1:B:223:GLN:HE21	1:B:224:THR:N	1.83	0.76
1:B:330:MET:HE1	1:B:342:TYR:HB2	1.69	0.75
1:B:86:ARG:HD2	1:B:295:ASP:OD1	1.87	0.75
1:B:223:GLN:HE21	1:B:224:THR:H	1.33	0.75
1:A:376:GLU:HG3	1:A:377:LYS:H	1.52	0.74
1:B:524:ASP:OD2	1:B:525:LYS:HE3	1.87	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:10:VAL:HG12	1:A:11:ALA:N	2.04	0.73
1:B:346:GLN:NE2	1:B:430:GLN:HE21	1.80	0.72
1:B:2:LYS:HE2	1:B:335:MET:O	1.90	0.72
1:B:288:ASN:HD21	1:B:292:GLN:HG3	1.55	0.71
1:B:431:ARG:O	1:B:431:ARG:HG2	1.91	0.69
1:A:408:ILE:HD12	1:A:434:MET:O	1.93	0.69
1:A:290:MET:CE	1:A:339:GLU:HA	2.24	0.68
1:A:341:ARG:HH11	1:A:341:ARG:CG	2.02	0.67
1:B:182:LEU:O	1:B:212:GLY:HA2	1.94	0.67
1:A:397:PHE:CZ	1:A:426:MET:HE1	2.29	0.67
1:B:288:ASN:ND2	1:B:292:GLN:HG3	2.10	0.67
1:A:341:ARG:HG2	1:A:341:ARG:NH1	1.97	0.66
1:A:486:ASP:OD1	1:A:550:ARG:NH2	2.27	0.66
1:B:243:HIS:HE1	1:B:255:GLU:HG3	1.60	0.66
1:A:397:PHE:HZ	1:A:426:MET:HE1	1.60	0.65
1:A:314:GLN:CA	1:A:314:GLN:HE21	2.09	0.65
1:A:290:MET:HE1	1:A:339:GLU:HA	1.77	0.65
1:A:243:HIS:HE1	1:A:255:GLU:OE2	1.79	0.65
1:A:455:ASP:O	1:A:458:ASN:N	2.29	0.63
1:B:43:CYS:SG	3:B:901:COA:S1P	2.72	0.63
1:A:162:ILE:HD12	2:A:900:FAD:C7	2.28	0.62
1:B:494:ASN:HD22	1:B:510:PRO:HB3	1.64	0.62
1:A:422:ARG:NH2	1:A:455:ASP:OD1	2.27	0.62
1:A:442:LEU:HD13	1:B:428:VAL:HG22	1.82	0.62
1:B:531:CYS:SG	1:B:532:GLN:N	2.73	0.62
1:A:56:ARG:NH1	1:A:140:ASP:OD1	2.33	0.61
1:A:90:LEU:HD23	1:A:105:SER:HA	1.80	0.61
1:A:112:SER:N	1:A:113:PRO:HD3	2.16	0.60
1:A:10:VAL:HG12	1:A:11:ALA:H	1.66	0.60
1:B:541:TYR:CZ	1:B:545:VAL:HG21	2.36	0.60
1:A:314:GLN:HA	1:A:314:GLN:NE2	2.15	0.59
1:B:5:LEU:HG	1:B:106:TYR:HB3	1.83	0.59
1:A:86:ARG:HD2	1:A:295:ASP:OD1	2.02	0.59
1:B:21:ARG:HG2	1:B:71:PHE:O	2.01	0.59
1:B:377:LYS:HD2	1:B:379:TYR:CE2	2.38	0.59
1:B:382:THR:HB	1:B:456:VAL:HG22	1.83	0.59
1:A:240:ILE:O	1:A:240:ILE:HG12	2.01	0.59
1:B:43:CYS:HG	3:B:901:COA:C2P	2.15	0.58
1:B:370:GLN:O	1:B:370:GLN:HG2	2.03	0.58
1:B:212:GLY:C	1:B:213:THR:CG2	2.71	0.58
1:B:223:GLN:HE21	1:B:225:HIS:H	1.51	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:431:ARG:CG	1:B:431:ARG:O	2.52	0.57
1:B:223:GLN:NE2	1:B:225:HIS:H	2.03	0.57
1:A:109:LEU:HD23	1:A:298:ILE:HD13	1.87	0.57
1:B:127:ASN:OD1	1:B:130:THR:HG23	2.05	0.56
1:A:459:GLN:O	1:A:463:VAL:HG23	2.04	0.56
1:A:10:VAL:O	1:A:14:ALA:HB3	2.05	0.56
1:B:212:GLY:O	1:B:213:THR:HG22	2.06	0.56
1:B:564:ALA:O	1:B:565:SER:HB3	2.06	0.55
1:A:426:MET:HE1	1:A:457:ILE:HD12	1.88	0.55
1:A:290:MET:N	1:A:291:MET:HA	2.22	0.55
1:B:125:VAL:O	1:B:125:VAL:HG22	2.06	0.55
1:B:188:THR:N	1:B:189:PRO:CD	2.69	0.55
1:B:223:GLN:NE2	1:B:224:THR:H	2.03	0.54
1:A:92:THR:OG1	1:A:103:GLN:NE2	2.37	0.54
1:A:112:SER:N	1:A:113:PRO:CD	2.70	0.53
1:B:9:GLY:O	1:B:38:VAL:HG22	2.09	0.53
1:B:288:ASN:HD21	1:B:292:GLN:HE21	1.56	0.53
1:A:287:VAL:HA	1:A:292:GLN:O	2.09	0.53
1:A:47:TYR:HB3	1:A:53:ILE:HG13	1.90	0.53
1:A:43:CYS:HG	3:A:901:COA:HS1	0.55	0.52
1:A:240:ILE:HG22	4:A:1052:HOH:O	2.09	0.52
1:A:197:PHE:HB2	1:A:414:VAL:HG11	1.92	0.52
1:B:377:LYS:HD2	1:B:379:TYR:HE2	1.75	0.51
1:A:108:THR:CG2	1:A:297:ALA:O	2.58	0.51
1:A:108:THR:HG23	1:A:297:ALA:O	2.10	0.51
1:B:191:ASP:OD1	1:B:366:LYS:HG3	2.11	0.51
1:A:47:TYR:HB2	1:A:53:ILE:HD12	1.93	0.51
1:B:212:GLY:C	1:B:213:THR:HG22	2.32	0.51
1:B:494:ASN:ND2	1:B:510:PRO:HB3	2.26	0.51
1:A:454:LYS:NZ	3:B:901:COA:O1A	2.44	0.50
1:A:215:LEU:CD1	1:A:246:LEU:HB3	2.42	0.50
1:A:424:ASP:O	1:A:428:VAL:HG12	2.13	0.49
1:A:180:LEU:CD2	1:A:210:ARG:HG3	2.42	0.49
1:B:88:ALA:HB1	1:B:90:LEU:HD22	1.95	0.49
1:B:131:HIS:HE1	1:B:145:THR:OG1	1.96	0.48
1:A:469:LYS:NZ	4:A:1045:HOH:O	2.46	0.48
1:A:380:VAL:HB	1:A:382:THR:HG23	1.95	0.48
1:A:151:VAL:HG22	1:A:257:ASP:HB2	1.96	0.48
1:B:408:ILE:HD12	1:B:434:MET:O	2.13	0.48
1:B:187:MET:C	1:B:189:PRO:HD2	2.34	0.48
1:B:56:ARG:NH2	1:B:140:ASP:OD2	2.45	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:110:LEU:HB2	1:B:335:MET:CE	2.44	0.48
1:A:472:ALA:HA	1:A:541:TYR:CE2	2.48	0.48
1:A:518:MET:SD	1:A:544:LEU:HD23	2.54	0.48
1:A:290:MET:HE3	1:A:339:GLU:HA	1.95	0.47
1:A:10:VAL:CG1	1:A:11:ALA:H	2.24	0.47
1:B:227:ALA:H	1:B:238:GLN:NE2	2.11	0.47
1:A:47:TYR:HB2	1:A:53:ILE:CD1	2.44	0.47
1:A:178:THR:HG21	1:A:254:LEU:HD11	1.97	0.47
1:B:178:THR:HG21	1:B:254:LEU:HD11	1.97	0.46
1:B:127:ASN:HD21	1:B:130:THR:CG2	2.28	0.46
1:B:42:ASN:HB2	3:B:901:COA:H22	1.98	0.46
1:B:489:LEU:HD23	1:B:506:ALA:HB2	1.97	0.46
1:B:112:SER:N	1:B:113:PRO:HD3	2.31	0.46
1:B:542:ARG:HD2	1:B:542:ARG:HA	1.75	0.46
1:A:180:LEU:HD23	1:A:210:ARG:HG3	1.97	0.46
1:A:440:GLN:NE2	4:A:1045:HOH:O	2.42	0.46
1:B:401:PHE:CZ	1:B:464:ALA:HB1	2.51	0.46
1:B:379:TYR:CE1	1:B:398:LYS:HD2	2.50	0.46
1:A:42:ASN:OD1	3:A:901:COA:H22	2.15	0.46
1:A:162:ILE:CD1	2:A:900:FAD:C7	2.94	0.46
1:A:34:ARG:O	1:A:78:LYS:HA	2.16	0.45
1:A:49:ILE:O	1:A:173:LEU:HD11	2.16	0.45
1:B:130:THR:HG22	1:B:244:LEU:HD11	1.98	0.45
1:B:436:VAL:O	1:B:437:GLU:C	2.53	0.45
1:A:2:LYS:HE2	1:A:335:MET:O	2.16	0.45
1:A:241:LYS:HD2	1:A:241:LYS:HA	1.50	0.45
1:B:131:HIS:HD2	4:B:1046:HOH:O	1.99	0.45
1:A:198:ALA:O	1:A:202:ILE:HG13	2.16	0.45
1:A:178:THR:HA	1:A:208:ASP:O	2.17	0.45
1:B:125:VAL:CG2	1:B:125:VAL:O	2.65	0.45
1:B:191:ASP:O	1:B:192:ARG:C	2.55	0.45
1:B:290:MET:O	1:B:291:MET:HB2	2.17	0.45
1:A:314:GLN:CA	1:A:314:GLN:NE2	2.78	0.44
1:A:309:ASP:OD1	1:A:363:LYS:HE3	2.18	0.44
1:B:330:MET:CE	1:B:342:TYR:HB2	2.45	0.44
1:B:185:GLN:HA	1:B:199:HIS:CE1	2.53	0.44
1:A:397:PHE:CZ	1:A:426:MET:CE	2.99	0.44
1:B:320:LEU:HA	1:B:320:LEU:HD23	1.80	0.44
1:A:397:PHE:HZ	1:A:426:MET:CE	2.30	0.43
1:B:529:ILE:HD11	1:B:544:LEU:HD12	1.99	0.43
1:B:221:GLN:HB3	1:B:221:GLN:HE21	1.69	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:314:GLN:OE1	1:B:314:GLN:HA	2.18	0.43
1:A:455:ASP:OD2	1:A:456:VAL:N	2.51	0.43
1:B:11:ALA:HB3	2:B:900:FAD:O5'	2.18	0.43
1:A:345:THR:HG22	1:A:347:GLY:H	1.83	0.43
1:B:122:ILE:O	1:B:125:VAL:HG12	2.18	0.43
1:B:215:LEU:CD1	1:B:246:LEU:HB3	2.49	0.43
3:B:901:COA:H131	3:B:901:COA:O9P	2.18	0.43
1:A:162:ILE:HD12	2:A:900:FAD:C7M	2.49	0.43
1:A:234:ASP:C	1:A:234:ASP:OD1	2.57	0.43
1:A:426:MET:CE	1:A:457:ILE:HD12	2.48	0.43
1:B:455:ASP:HB2	4:B:1058:HOH:O	2.19	0.43
1:A:421:LYS:O	1:A:425:VAL:HG23	2.19	0.43
1:B:377:LYS:HZ2	1:B:476:HIS:CG	2.37	0.43
1:A:188:THR:N	1:A:189:PRO:CD	2.81	0.43
1:A:450:TYR:OH	1:B:352:LYS:HG3	2.20	0.42
1:B:211:LEU:O	1:B:213:THR:HG23	2.19	0.42
1:B:127:ASN:ND2	1:B:130:THR:HG23	2.34	0.42
1:A:529:ILE:HG13	1:A:552:ARG:O	2.19	0.42
1:B:171:HIS:CE1	1:B:205:GLN:HE21	2.38	0.42
1:A:279:ILE:HG22	1:A:280:GLY:O	2.20	0.42
2:A:900:FAD:H9	2:A:900:FAD:H1'1	1.83	0.41
1:B:376:GLU:HG3	1:B:377:LYS:H	1.85	0.41
1:A:518:MET:HE1	1:A:547:ASN:OD1	2.19	0.41
2:B:900:FAD:H1'1	2:B:900:FAD:H9	1.82	0.41
1:B:215:LEU:HD11	1:B:246:LEU:HB3	2.01	0.41
1:B:425:VAL:HG11	1:B:457:ILE:HD12	2.01	0.41
1:A:227:ALA:H	1:A:238:GLN:NE2	2.18	0.41
1:A:434:MET:HB3	1:A:434:MET:HE2	1.95	0.41
1:B:386:ALA:HB1	1:B:388:TYR:CE2	2.55	0.41
1:A:455:ASP:O	1:A:456:VAL:C	2.59	0.41
1:A:40:PHE:O	3:A:901:COA:H62	2.21	0.41
1:A:499:GLN:NE2	1:A:499:GLN:H	2.18	0.41
1:A:430:GLN:O	1:A:430:GLN:HG2	2.20	0.41
1:B:112:SER:O	2:B:900:FAD:H52A	2.20	0.41
1:A:188:THR:N	1:A:189:PRO:HD2	2.36	0.41
1:A:162:ILE:H	1:A:162:ILE:HG12	1.72	0.40
1:B:267:PRO:HB3	1:B:304:ALA:HA	2.03	0.40
1:B:521:LEU:HA	1:B:522:PRO:HD2	1.86	0.40
1:B:198:ALA:O	1:B:202:ILE:HG13	2.22	0.40
3:B:901:COA:HO2A	3:B:901:COA:P3B	2.44	0.40
1:A:112:SER:O	2:A:900:FAD:H52A	2.21	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:225:HIS:ND1	1:A:512:ASP:OD2[2_555]	2.16	0.04

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	563/574 (98%)	541 (96%)	21 (4%)	1 (0%)	52 83
1	B	563/574 (98%)	537 (95%)	24 (4%)	2 (0%)	39 72
All	All	1126/1148 (98%)	1078 (96%)	45 (4%)	3 (0%)	46 77

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	10	VAL
1	B	10	VAL
1	B	263	ILE

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	449/458 (98%)	407 (91%)	42 (9%)	11 28
1	B	448/458 (98%)	402 (90%)	46 (10%)	9 23

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	897/916 (98%)	809 (90%)	88 (10%)	10	26

All (88) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	2	LYS
1	A	34	ARG
1	A	42	ASN
1	A	56	ARG
1	A	57	SER
1	A	67	PHE
1	A	111	LEU
1	A	119	VAL
1	A	129	LEU
1	A	132	SER
1	A	143	LEU
1	A	200	GLN
1	A	208	ASP
1	A	209	LEU
1	A	210	ARG
1	A	229	ASP
1	A	240	ILE
1	A	241	LYS
1	A	254	LEU
1	A	259	LEU
1	A	266	ARG
1	A	267	PRO
1	A	269	THR
1	A	282	LEU
1	A	291	MET
1	A	307	GLU
1	A	314	GLN
1	A	341	ARG
1	A	356	LEU
1	A	370	GLN
1	A	377	LYS
1	A	378	VAL
1	A	407	THR
1	A	428	VAL
1	A	445	SER
1	A	483	LEU

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Mol	Chain	Res	Type
1	A	486	ASP
1	A	490	LEU
1	A	523	LYS
1	A	525	LYS
1	A	531	CYS
1	B	24	SER
1	B	34	ARG
1	B	38	VAL
1	B	42	ASN
1	B	52	GLU
1	B	67	PHE
1	B	90	LEU
1	B	97	LEU
1	B	103	GLN
1	B	119	VAL
1	B	130	THR
1	B	144	GLN
1	B	176	LYS
1	B	181	GLU
1	B	182	LEU
1	B	208	ASP
1	B	209	LEU
1	B	213	THR
1	B	219	SER
1	B	221	GLN
1	B	241	LYS
1	B	254	LEU
1	B	255	GLU
1	B	258	LEU
1	B	273	ARG
1	B	279	ILE
1	B	281	GLU
1	B	292	GLN
1	B	307	GLU
1	B	308	GLN
1	B	341	ARG
1	B	356	LEU
1	B	378	VAL
1	B	380	VAL
1	B	385	HIS
1	B	407	THR
1	B	436	VAL

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Mol	Chain	Res	Type
1	B	437	GLU
1	B	483	LEU
1	B	485	GLU
1	B	488	LEU
1	B	498	LEU
1	B	511	VAL
1	B	525	LYS
1	B	531	CYS
1	B	533	VAL

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	223	GLN
1	A	238	GLN
1	A	243	HIS
1	A	292	GLN
1	A	314	GLN
1	A	385	HIS
1	A	440	GLN
1	A	441	HIS
1	A	499	GLN
1	A	543	GLN
1	A	546	ASN
1	B	62	GLN
1	B	131	HIS
1	B	200	GLN
1	B	205	GLN
1	B	221	GLN
1	B	223	GLN
1	B	238	GLN
1	B	243	HIS
1	B	292	GLN
1	B	308	GLN
1	B	370	GLN
1	B	430	GLN
1	B	440	GLN
1	B	482	ASN
1	B	494	ASN
1	B	543	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

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	FAD	A	900	-	48,58,58	1.28	4 (8%)	54,89,89	2.66	17 (31%)
3	COA	A	901	-	40,50,50	1.72	3 (7%)	50,75,75	2.26	12 (24%)
2	FAD	B	900	-	48,58,58	1.29	5 (10%)	54,89,89	2.00	10 (18%)
3	COA	B	901	-	40,50,50	1.63	3 (7%)	50,75,75	2.00	10 (20%)

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	FAD	A	900	-	-	0/30/50/50	0/6/6/6
3	COA	A	901	-	-	0/44/64/64	0/3/3/3
2	FAD	B	900	-	-	0/30/50/50	0/6/6/6
3	COA	B	901	-	-	0/44/64/64	0/3/3/3

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	900	FAD	O4B-C1B	2.40	1.44	1.41
2	A	900	FAD	C4-N3	2.71	1.38	1.33
3	B	901	COA	C2A-N1A	2.71	1.39	1.33
2	B	900	FAD	C4X-N5	2.76	1.37	1.33
3	A	901	COA	C2A-N1A	2.89	1.39	1.33
3	A	901	COA	C2A-N3A	3.01	1.37	1.32
2	A	900	FAD	C2A-N3A	3.23	1.37	1.32
2	B	900	FAD	C4-N3	3.29	1.39	1.33
3	B	901	COA	C2A-N3A	3.46	1.38	1.32
2	B	900	FAD	C2A-N1A	3.48	1.40	1.33
2	A	900	FAD	C5X-N5	3.51	1.40	1.35
2	B	900	FAD	C2A-N3A	4.31	1.39	1.32
2	A	900	FAD	C4X-N5	4.34	1.40	1.33
3	B	901	COA	O9P-C9P	8.15	1.39	1.23
3	A	901	COA	O9P-C9P	9.05	1.41	1.23

All (49) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	900	FAD	N3A-C2A-N1A	-13.43	118.61	128.89
3	A	901	COA	N3A-C2A-N1A	-11.82	119.85	128.89
3	B	901	COA	N3A-C2A-N1A	-9.21	121.84	128.89
2	B	900	FAD	N3A-C2A-N1A	-7.72	122.98	128.89
2	A	900	FAD	C4B-O4B-C1B	-4.82	104.42	109.72
2	A	900	FAD	C2B-C1B-N9A	-4.35	107.65	114.29
2	B	900	FAD	C4X-C4-N3	-3.76	118.44	123.59
3	B	901	COA	P2A-O3A-P1A	-3.25	123.59	132.73
2	A	900	FAD	C1B-N9A-C4A	-3.06	122.33	126.94
2	A	900	FAD	O2'-C2'-C1'	-2.98	102.62	109.94
3	A	901	COA	C2B-C1B-N9A	-2.78	110.05	114.29
3	A	901	COA	O3A-P2A-O6A	-2.72	95.72	102.94
2	A	900	FAD	C4-C4X-C10	-2.55	118.31	119.94
3	A	901	COA	C6P-C5P-N4P	-2.51	112.10	116.46
3	B	901	COA	O2B-C2B-C3B	-2.42	104.16	111.16
3	B	901	COA	C4A-C5A-N7A	-2.42	107.25	109.48
3	B	901	COA	C2P-C3P-N4P	-2.40	107.64	112.37
2	A	900	FAD	C9A-C5X-N5	-2.38	118.84	122.36
2	B	900	FAD	C9A-C5X-N5	-2.37	118.84	122.36
3	A	901	COA	O5B-C5B-C4B	-2.31	100.59	109.12
3	A	901	COA	P2A-O3A-P1A	-2.31	126.25	132.73
2	A	900	FAD	C4A-C5A-N7A	-2.22	107.44	109.48
3	A	901	COA	C4A-C5A-N7A	-2.19	107.46	109.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	900	FAD	C4X-C10-N10	2.10	121.76	120.52
2	A	900	FAD	C2A-N1A-C6A	2.11	122.54	118.77
2	B	900	FAD	C6-C5X-C9A	2.18	121.85	118.98
2	A	900	FAD	O4B-C1B-N9A	2.26	112.83	108.10
2	A	900	FAD	O4B-C4B-C3B	2.29	109.76	105.15
3	A	901	COA	O6A-CCP-CBP	2.30	114.24	110.55
3	B	901	COA	O2A-P1A-O3A	2.32	115.61	105.09
2	B	900	FAD	C1B-N9A-C4A	2.36	130.50	126.94
3	A	901	COA	CEP-CBP-CCP	2.52	111.77	108.50
3	B	901	COA	O5A-P2A-O3A	2.68	117.27	105.09
2	A	900	FAD	O2A-PA-O3P	2.73	117.46	105.09
3	B	901	COA	CAP-C9P-N8P	2.80	122.67	116.47
3	A	901	COA	O2A-P1A-O3A	2.84	117.98	105.09
3	A	901	COA	O5A-P2A-O3A	2.94	118.41	105.09
2	B	900	FAD	C4-C4X-C10	3.05	121.90	119.94
2	A	900	FAD	C4X-N5-C5X	3.39	120.66	116.76
2	A	900	FAD	C4-C4X-N5	3.46	122.91	118.72
2	B	900	FAD	O3P-PA-O5B	3.61	112.52	102.94
2	A	900	FAD	C5X-C9A-N10	3.62	120.37	117.62
3	A	901	COA	O9A-P3B-O8A	3.66	121.30	107.38
3	B	901	COA	OAP-CAP-C9P	3.67	118.81	110.38
2	B	900	FAD	C4X-N5-C5X	3.68	121.00	116.76
2	B	900	FAD	C5X-C9A-N10	4.50	121.04	117.62
3	B	901	COA	C3P-N4P-C5P	4.67	131.98	122.79
2	A	900	FAD	C4-N3-C2	5.51	120.01	115.25
2	B	900	FAD	C4-N3-C2	5.89	120.34	115.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 19 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	900	FAD	5	0
3	A	901	COA	4	0
2	B	900	FAD	3	0
3	B	901	COA	7	0

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 [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	565/574 (98%)	-0.57	0 100 100	26, 36, 54, 72	0
1	B	565/574 (98%)	-0.54	0 100 100	24, 35, 53, 76	0
All	All	1130/1148 (98%)	-0.55	0 100 100	24, 35, 54, 76	0

There are no RSRZ outliers to report.

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
2	FAD	A	900	53/53	0.99	0.13	0.20	23,30,33,35	0
2	FAD	B	900	53/53	0.99	0.13	0.11	18,27,31,32	0
3	COA	B	901	48/48	0.99	0.13	-0.24	24,32,41,46	0
3	COA	A	901	48/48	0.99	0.13	-0.46	20,30,35,45	0

6.5 Other polymers

There are no such residues in this entry.