



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 3, 2023 – 02:01 AM EDT

PDB ID : 3VQR
Title : Structure of a dye-linked L-proline dehydrogenase mutant from the aerobic hyperthermophilic archaeon, *Aeropyrum pernix*
Authors : Sakuraba, H.; Ohshima, T.; Satomura, T.; Yoneda, K.
Deposited on : 2012-03-29
Resolution : 2.01 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.36
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36

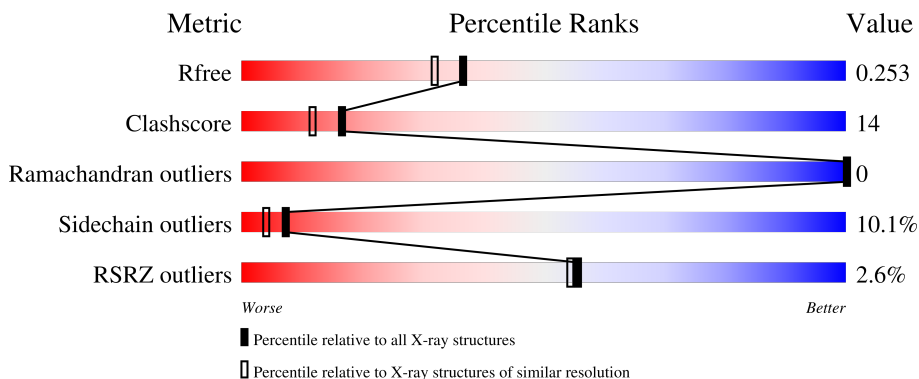
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.01 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}	130704	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	447	
1	B	447	

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	ACT	A	1002	-	-	X	-

2 Entry composition i

There are 5 unique types of molecules in this entry. The entry contains 6893 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 Putative oxidoreductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	425	3275	2081	573	612	9	0	0	0
1	B	425	3275	2081	573	612	9	0	0	0

There are 42 discrepancies between the modelled and reference sequences:

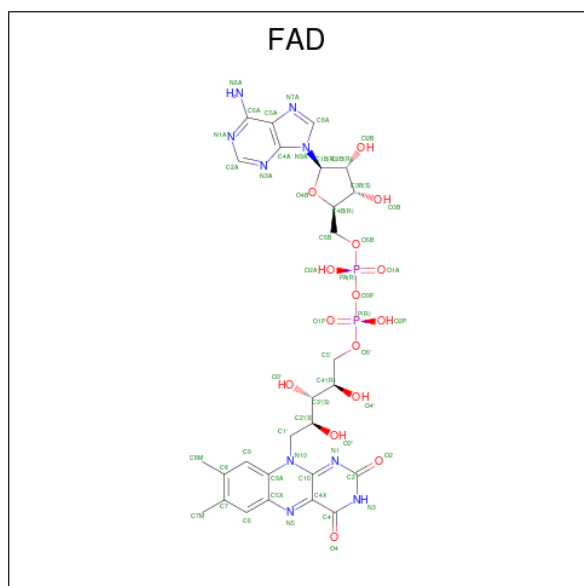
Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	expression tag	UNP Q9YCJ0
A	-18	GLY	-	expression tag	UNP Q9YCJ0
A	-17	SER	-	expression tag	UNP Q9YCJ0
A	-16	SER	-	expression tag	UNP Q9YCJ0
A	-15	HIS	-	expression tag	UNP Q9YCJ0
A	-14	HIS	-	expression tag	UNP Q9YCJ0
A	-13	HIS	-	expression tag	UNP Q9YCJ0
A	-12	HIS	-	expression tag	UNP Q9YCJ0
A	-11	HIS	-	expression tag	UNP Q9YCJ0
A	-10	HIS	-	expression tag	UNP Q9YCJ0
A	-9	SER	-	expression tag	UNP Q9YCJ0
A	-8	SER	-	expression tag	UNP Q9YCJ0
A	-7	GLY	-	expression tag	UNP Q9YCJ0
A	-6	LEU	-	expression tag	UNP Q9YCJ0
A	-5	VAL	-	expression tag	UNP Q9YCJ0
A	-4	PRO	-	expression tag	UNP Q9YCJ0
A	-3	ARG	-	expression tag	UNP Q9YCJ0
A	-2	GLY	-	expression tag	UNP Q9YCJ0
A	-1	SER	-	expression tag	UNP Q9YCJ0
A	0	HIS	-	expression tag	UNP Q9YCJ0
A	?	-	LEU	deletion	UNP Q9YCJ0
B	-19	MET	-	expression tag	UNP Q9YCJ0
B	-18	GLY	-	expression tag	UNP Q9YCJ0
B	-17	SER	-	expression tag	UNP Q9YCJ0
B	-16	SER	-	expression tag	UNP Q9YCJ0

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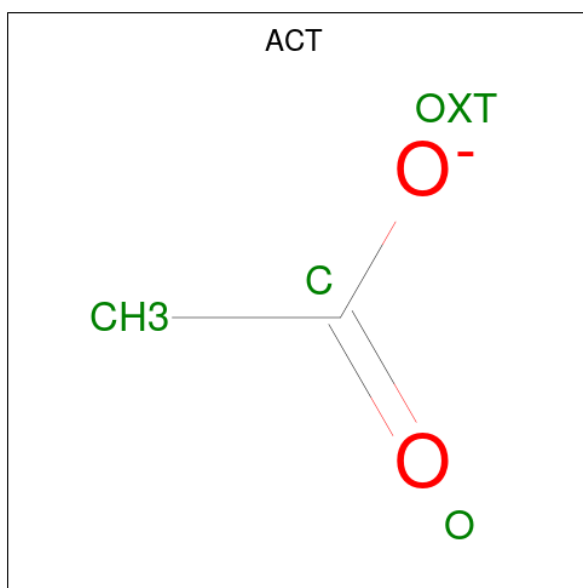
Chain	Residue	Modelled	Actual	Comment	Reference
B	-15	HIS	-	expression tag	UNP Q9YJCJ0
B	-14	HIS	-	expression tag	UNP Q9YJCJ0
B	-13	HIS	-	expression tag	UNP Q9YJCJ0
B	-12	HIS	-	expression tag	UNP Q9YJCJ0
B	-11	HIS	-	expression tag	UNP Q9YJCJ0
B	-10	HIS	-	expression tag	UNP Q9YJCJ0
B	-9	SER	-	expression tag	UNP Q9YJCJ0
B	-8	SER	-	expression tag	UNP Q9YJCJ0
B	-7	GLY	-	expression tag	UNP Q9YJCJ0
B	-6	LEU	-	expression tag	UNP Q9YJCJ0
B	-5	VAL	-	expression tag	UNP Q9YJCJ0
B	-4	PRO	-	expression tag	UNP Q9YJCJ0
B	-3	ARG	-	expression tag	UNP Q9YJCJ0
B	-2	GLY	-	expression tag	UNP Q9YJCJ0
B	-1	SER	-	expression tag	UNP Q9YJCJ0
B	0	HIS	-	expression tag	UNP Q9YJCJ0
B	?	-	LEU	deletion	UNP Q9YJCJ0

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



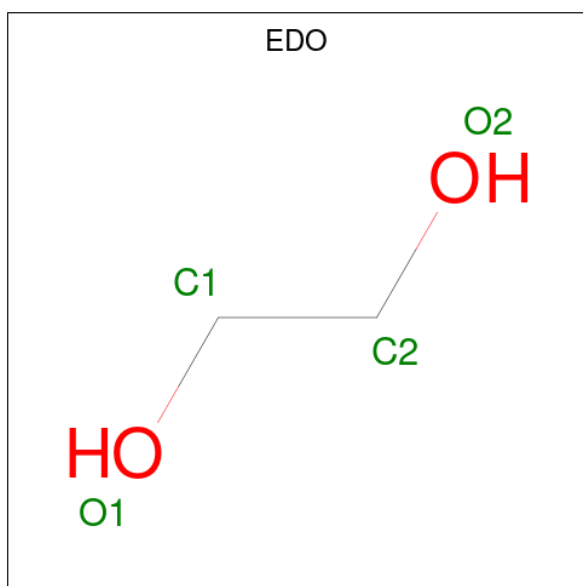
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf		
			Total	C	N	O			P	
2	A	1	Total	53	27	9	15	2	0	0
2	B	1	Total	53	27	9	15	2	0	0

- Molecule 3 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf	
3	A	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		

- Molecule 4 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).



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

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O 4 2 2	0	0
4	B	1	Total C O 4 2 2	0	0
4	B	1	Total C O 4 2 2	0	0

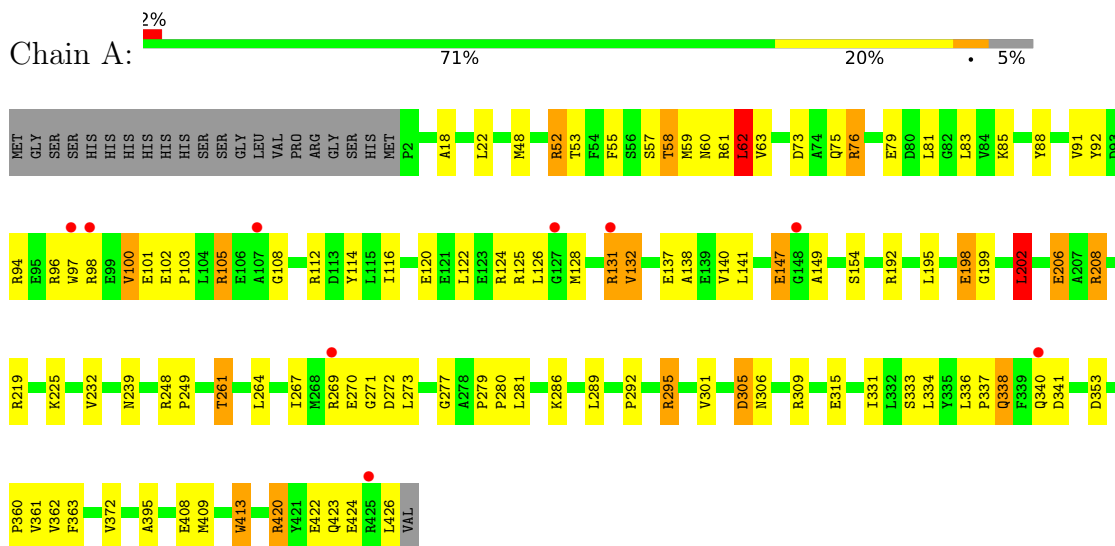
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	104	Total O 104 104	0	0
5	B	109	Total O 109 109	0	0

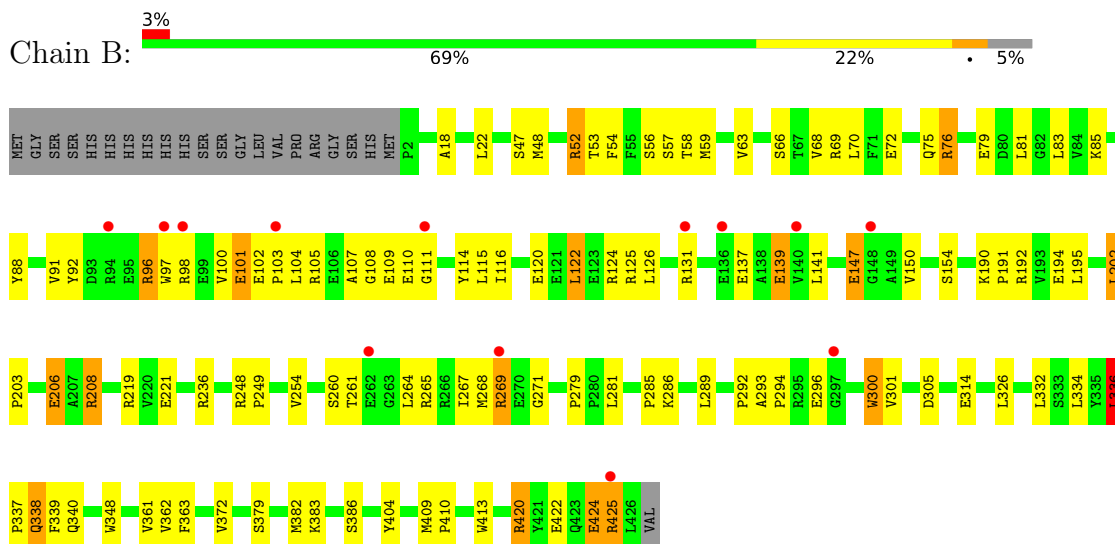
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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: Putative oxidoreductase



- Molecule 1: Putative oxidoreductase



4 Data and refinement statistics i

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	47.99Å 70.57Å 73.09Å 71.65° 70.85° 70.15°	Depositor
Resolution (Å)	24.20 – 2.01 24.20 – 2.01	Depositor EDS
% Data completeness (in resolution range)	(Not available) (24.20-2.01) 95.1 (24.20-2.01)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	8.16 (at 2.01Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.210 , 0.255 0.212 , 0.253	Depositor DCC
R_{free} test set	2654 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	26.8	Xtriage
Anisotropy	0.655	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 34.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	0.489 for h,h-k,h-l 0.015 for -h,-l,-k 0.014 for -h,-h+l,-h+k	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6893	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.89% 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: FAD, EDO, ACT

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	1.07	2/3349 (0.1%)	1.02	5/4538 (0.1%)
1	B	1.04	2/3349 (0.1%)	1.01	6/4538 (0.1%)
All	All	1.06	4/6698 (0.1%)	1.01	11/9076 (0.1%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	97	TRP	CD2-CE2	5.81	1.48	1.41
1	B	348	TRP	CD2-CE2	5.81	1.48	1.41
1	B	300	TRP	CD2-CE2	5.49	1.48	1.41
1	A	413	TRP	CD2-CE2	5.37	1.47	1.41

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	202	LEU	CB-CG-CD1	5.93	121.08	111.00
1	A	52	ARG	NE-CZ-NH1	5.88	123.24	120.30
1	A	126	LEU	CA-CB-CG	5.86	128.78	115.30
1	B	122	LEU	CA-CB-CG	5.61	128.20	115.30
1	B	202	LEU	CB-CG-CD2	5.56	120.45	111.00
1	B	236	ARG	NE-CZ-NH1	-5.43	117.58	120.30
1	A	62	LEU	CA-CB-CG	5.38	127.68	115.30
1	B	52	ARG	NE-CZ-NH1	5.34	122.97	120.30
1	B	336	LEU	CB-CG-CD1	-5.33	101.94	111.00
1	A	309	ARG	NE-CZ-NH1	5.25	122.93	120.30
1	B	126	LEU	CA-CB-CG	5.09	127.00	115.30

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	3275	0	3234	86	0
1	B	3275	0	3234	102	0
2	A	53	0	31	0	0
2	B	53	0	31	3	0
3	A	4	0	3	6	0
3	B	4	0	3	1	0
4	A	8	0	12	0	0
4	B	8	0	12	2	0
5	A	104	0	0	8	0
5	B	109	0	0	4	0
All	All	6893	0	6560	188	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 (188) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:336:LEU:HD13	1:B:339:PHE:CD2	1.83	1.13
1:A:261:THR:HG22	1:A:264:LEU:H	1.09	1.09
1:B:332:LEU:HD12	1:B:336:LEU:HD11	1.25	1.07
1:B:191:PRO:HG2	1:B:194:GLU:HG3	1.39	1.05
1:A:76:ARG:HG2	1:A:76:ARG:HH11	1.30	0.96
1:A:131:ARG:NE	1:A:131:ARG:O	1.99	0.95
1:B:336:LEU:HB2	1:B:338:GLN:HE22	1.29	0.94
1:B:332:LEU:CD1	1:B:336:LEU:HD11	2.00	0.92
1:B:75:GLN:HE21	1:B:81:LEU:H	1.21	0.89
1:A:76:ARG:HH11	1:A:76:ARG:CG	1.88	0.85
1:A:75:GLN:HE21	1:A:81:LEU:H	1.25	0.85
1:A:336:LEU:HB3	1:A:338:GLN:HE22	1.39	0.84
1:A:267:ILE:HD11	1:A:338:GLN:NE2	1.92	0.84
1:B:57:SER:HB3	1:B:424:GLU:HG2	1.60	0.83
1:A:261:THR:HG22	1:A:264:LEU:N	1.93	0.82
1:B:336:LEU:HD13	1:B:339:PHE:HD2	1.42	0.81
1:B:75:GLN:HE22	1:B:83:LEU:H	1.29	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:102:GLU:HB2	1:B:103:PRO:HD3	1.65	0.78
1:B:265:ARG:O	1:B:269:ARG:HD2	1.83	0.78
1:B:191:PRO:HG2	1:B:194:GLU:CG	2.13	0.77
1:A:116:ILE:HD12	1:A:116:ILE:O	1.83	0.77
1:A:75:GLN:HE22	1:A:83:LEU:H	1.36	0.74
1:A:261:THR:CG2	1:A:264:LEU:H	1.96	0.72
1:A:61:ARG:HH11	1:A:61:ARG:HG3	1.56	0.70
1:B:332:LEU:HD12	1:B:336:LEU:CD1	2.15	0.70
1:A:132:VAL:HG13	1:A:138:ALA:HB2	1.75	0.69
1:B:105:ARG:HA	5:B:1137:HOH:O	1.93	0.68
1:B:268:MET:HE3	1:B:279:PRO:HG2	1.73	0.68
1:B:268:MET:CE	1:B:279:PRO:CD	2.71	0.68
1:A:267:ILE:HD11	1:A:338:GLN:HE21	1.56	0.68
1:A:102:GLU:HB2	1:A:103:PRO:HD3	1.76	0.68
1:A:105:ARG:HA	5:A:1188:HOH:O	1.94	0.67
1:B:59:MET:HE3	1:B:413:TRP:HB3	1.76	0.67
1:A:131:ARG:O	1:A:131:ARG:CD	2.43	0.66
1:A:92:TYR:CE1	1:A:147:GLU:HB2	2.31	0.66
1:A:122:LEU:HD23	1:A:128:MET:SD	2.37	0.65
1:B:336:LEU:HD13	1:B:339:PHE:CE2	2.32	0.65
1:A:57:SER:HB3	1:A:424:GLU:OE1	1.96	0.65
1:B:279:PRO:O	1:B:292:PRO:HG3	1.96	0.64
1:B:75:GLN:HE21	1:B:81:LEU:N	1.96	0.64
1:A:125:ARG:HD3	1:A:270:GLU:O	1.97	0.64
1:A:52:ARG:HD3	1:A:88:TYR:CE2	2.33	0.64
1:A:295:ARG:HG3	5:A:1195:HOH:O	1.98	0.64
1:B:76:ARG:CB	1:B:76:ARG:HH11	2.13	0.62
1:B:338:GLN:H	1:B:338:GLN:CD	2.01	0.62
1:B:261:THR:CG2	1:B:264:LEU:H	2.13	0.62
1:B:190:LYS:HE2	1:B:194:GLU:OE1	1.99	0.62
1:A:76:ARG:HG2	1:A:76:ARG:NH1	2.02	0.61
1:B:337:PRO:O	1:B:340:GLN:HG3	2.00	0.61
1:B:92:TYR:CE1	1:B:147:GLU:HB2	2.35	0.61
1:A:105:ARG:HH11	1:A:105:ARG:HG3	1.65	0.61
1:B:363:PHE:CE1	4:B:1004:EDO:H11	2.36	0.61
1:A:58:THR:OG1	1:A:59:MET:N	2.34	0.61
1:A:137:GLU:HG2	1:A:334:LEU:HD13	1.82	0.60
1:B:111:GLY:HA2	1:B:115:LEU:HD23	1.83	0.60
1:B:268:MET:HE3	1:B:279:PRO:CG	2.29	0.60
1:B:382:MET:HG2	2:B:1001:FAD:O2	2.01	0.60
1:A:131:ARG:O	1:A:131:ARG:CZ	2.48	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:83:LEU:O	1:A:83:LEU:HD23	2.02	0.60
1:A:75:GLN:HA	1:A:79:GLU:O	2.01	0.60
1:A:333:SER:HB2	1:A:340:GLN:HA	1.84	0.60
1:A:125:ARG:HD2	1:A:273:LEU:HD23	1.83	0.59
1:B:75:GLN:HA	1:B:79:GLU:O	2.02	0.58
1:B:139:GLU:HG2	5:B:1154:HOH:O	2.03	0.58
1:B:59:MET:CE	1:B:413:TRP:HB3	2.33	0.58
1:B:192:ARG:HG3	1:B:206:GLU:HB3	1.86	0.58
1:B:59:MET:CE	1:B:413:TRP:CE3	2.87	0.58
1:B:305:ASP:N	1:B:305:ASP:OD1	2.34	0.57
1:B:70:LEU:HD12	1:B:404:TYR:CD1	2.40	0.57
1:A:272:ASP:OD2	1:A:277:GLY:N	2.39	0.56
1:A:279:PRO:O	1:A:292:PRO:HG3	2.06	0.56
1:A:338:GLN:H	1:A:338:GLN:CD	2.08	0.56
1:B:268:MET:HE3	1:B:279:PRO:CD	2.34	0.56
1:A:141:LEU:HD12	1:A:334:LEU:HD22	1.88	0.56
1:B:116:ILE:HG12	5:B:1204:HOH:O	2.04	0.56
1:A:62:LEU:HD12	1:A:63:VAL:N	2.20	0.56
1:A:105:ARG:HH11	1:A:105:ARG:CG	2.19	0.55
1:A:122:LEU:CD2	1:A:128:MET:SD	2.94	0.55
1:A:267:ILE:O	1:A:271:GLY:HA3	2.05	0.55
1:B:116:ILE:HG13	1:B:116:ILE:O	2.06	0.55
1:A:353:ASP:HB2	3:A:1002:ACT:H2	1.88	0.55
1:B:58:THR:O	1:B:59:MET:C	2.45	0.55
1:A:61:ARG:HG3	1:A:61:ARG:NH1	2.19	0.55
1:B:268:MET:HE2	1:B:279:PRO:HD3	1.89	0.54
1:B:59:MET:CE	1:B:413:TRP:HE3	2.20	0.54
1:B:362:VAL:HA	1:B:372:VAL:O	2.08	0.54
1:A:295:ARG:CG	5:A:1195:HOH:O	2.56	0.54
1:B:332:LEU:O	1:B:336:LEU:HD12	2.07	0.54
1:B:261:THR:HG23	1:B:264:LEU:H	1.74	0.53
1:A:53:THR:CG2	1:A:154:SER:OG	2.56	0.53
1:B:293:ALA:HB1	1:B:296:GLU:HG3	1.90	0.53
1:B:59:MET:HE1	1:B:413:TRP:HE3	1.72	0.53
1:B:137:GLU:HG2	1:B:334:LEU:HD13	1.89	0.53
1:A:232:VAL:HG13	3:A:1002:ACT:C	2.38	0.53
3:A:1002:ACT:CH3	5:A:1185:HOH:O	2.56	0.53
1:B:97:TRP:NE1	1:B:101:GLU:HG3	2.23	0.53
1:B:116:ILE:O	1:B:116:ILE:CG1	2.58	0.52
1:B:102:GLU:CB	1:B:103:PRO:HD3	2.37	0.52
1:A:75:GLN:HE21	1:A:81:LEU:N	2.01	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:268:MET:CE	1:B:279:PRO:HD3	2.40	0.52
3:A:1002:ACT:H1	5:A:1185:HOH:O	2.08	0.52
1:B:267:ILE:O	1:B:271:GLY:HA3	2.10	0.52
1:A:305:ASP:OD1	1:A:305:ASP:N	2.40	0.52
1:B:379:SER:O	1:B:383:LYS:HG3	2.10	0.51
1:B:409:MET:HG3	1:B:413:TRP:CE3	2.45	0.51
1:B:332:LEU:CD1	1:B:336:LEU:CD1	2.81	0.51
1:A:73:ASP:HA	1:A:76:ARG:NH1	2.25	0.51
1:A:102:GLU:CB	1:A:103:PRO:HD3	2.40	0.51
1:A:83:LEU:HD23	1:A:83:LEU:C	2.31	0.51
1:A:92:TYR:HE1	1:A:147:GLU:HB2	1.75	0.51
1:A:239:ASN:HD22	1:B:248:ARG:HH21	1.59	0.51
1:A:239:ASN:ND2	1:B:248:ARG:HH21	2.09	0.50
1:A:362:VAL:HA	1:A:372:VAL:O	2.12	0.50
1:B:66:SER:HA	1:B:69:ARG:HG2	1.93	0.50
1:B:261:THR:HG22	1:B:264:LEU:HB2	1.92	0.50
1:A:337:PRO:O	1:A:340:GLN:HG3	2.12	0.49
1:B:96:ARG:HD3	1:B:96:ARG:N	2.26	0.49
1:A:208:ARG:HD3	5:A:1173:HOH:O	2.11	0.49
1:B:97:TRP:CE2	1:B:101:GLU:HG3	2.48	0.49
1:B:268:MET:HE1	1:B:294:PRO:HG3	1.95	0.48
1:B:363:PHE:HE1	4:B:1004:EDO:H11	1.77	0.48
1:B:336:LEU:CD1	1:B:339:PHE:CD2	2.76	0.48
1:B:102:GLU:HB2	1:B:103:PRO:CD	2.38	0.48
1:A:108:GLY:O	1:A:114:TYR:HE2	1.98	0.47
1:A:132:VAL:HG13	1:A:138:ALA:CB	2.44	0.47
1:A:58:THR:HG21	1:A:420:ARG:NH2	2.30	0.47
1:A:105:ARG:HB2	1:A:105:ARG:NH1	2.30	0.47
1:A:353:ASP:CB	3:A:1002:ACT:H2	2.45	0.47
1:B:420:ARG:O	1:B:420:ARG:HG3	2.14	0.47
1:A:100:VAL:O	1:A:103:PRO:HD2	2.15	0.47
1:A:141:LEU:O	1:A:286:LYS:NZ	2.43	0.46
1:B:59:MET:HE3	1:B:420:ARG:HB3	1.97	0.46
1:A:18:ALA:O	1:A:22:LEU:HB2	2.16	0.46
1:A:198:GLU:HB3	1:B:326:LEU:HD13	1.98	0.45
1:A:76:ARG:HH11	1:A:76:ARG:CB	2.28	0.45
1:B:141:LEU:HD12	1:B:334:LEU:HD22	1.99	0.45
1:B:202:LEU:HB3	1:B:203:PRO:HD2	1.98	0.45
1:A:192:ARG:HG3	1:A:206:GLU:HB3	1.97	0.45
1:B:52:ARG:HD3	1:B:88:TYR:CE2	2.52	0.45
1:B:111:GLY:HA2	1:B:115:LEU:CD2	2.46	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:289:LEU:O	1:A:301:VAL:HG13	2.18	0.44
1:A:248:ARG:HG3	1:A:249:PRO:HD2	1.99	0.44
1:B:58:THR:OG1	1:B:59:MET:N	2.50	0.44
1:A:409:MET:HG3	1:A:413:TRP:CE3	2.53	0.44
1:B:63:VAL:HG13	1:B:386:SER:HB2	1.99	0.44
1:B:248:ARG:HG3	1:B:249:PRO:HD2	1.98	0.44
1:A:360:PRO:HD3	5:A:1120:HOH:O	2.17	0.44
1:B:102:GLU:CB	1:B:103:PRO:CD	2.95	0.44
1:B:59:MET:HE2	1:B:413:TRP:CE3	2.53	0.44
1:B:219:ARG:NH1	1:B:221:GLU:HG2	2.32	0.43
1:B:336:LEU:CB	1:B:338:GLN:HE22	2.15	0.43
1:A:199:GLY:O	1:A:202:LEU:HD13	2.19	0.43
1:B:338:GLN:CD	1:B:338:GLN:N	2.71	0.43
1:A:131:ARG:O	1:A:131:ARG:CG	2.67	0.43
1:A:53:THR:HG21	1:A:154:SER:OG	2.19	0.43
1:B:18:ALA:O	1:B:22:LEU:HB2	2.18	0.42
1:A:116:ILE:O	1:A:116:ILE:CD1	2.62	0.42
1:A:105:ARG:CG	1:A:105:ARG:NH1	2.83	0.42
1:B:76:ARG:HH11	1:B:76:ARG:HB3	1.84	0.42
1:A:424:GLU:HG3	1:A:426:LEU:H	1.84	0.42
1:B:59:MET:HG2	1:B:420:ARG:HG2	2.00	0.42
1:B:68:VAL:O	1:B:72:GLU:HG3	2.20	0.42
1:B:261:THR:HG22	1:B:264:LEU:CB	2.49	0.42
1:A:280:PRO:HD3	5:A:1129:HOH:O	2.20	0.42
1:B:54:PHE:CZ	1:B:107:ALA:HB3	2.54	0.42
1:B:254:VAL:CG1	1:B:300:TRP:HB3	2.50	0.41
1:A:361:VAL:HG12	1:A:363:PHE:CD2	2.56	0.41
1:B:56:SER:OG	1:B:425:ARG:HG2	2.20	0.41
1:B:59:MET:HE1	1:B:413:TRP:CE3	2.52	0.41
1:B:409:MET:HA	1:B:410:PRO:HD3	1.99	0.41
1:A:122:LEU:HD21	1:A:149:ALA:HB2	2.03	0.41
1:B:208:ARG:HD3	5:B:1177:HOH:O	2.20	0.41
1:A:337:PRO:HD2	1:A:338:GLN:OE1	2.20	0.41
1:B:104:LEU:HD11	1:B:150:VAL:HG21	2.02	0.41
1:B:91:VAL:HG22	1:B:285:PRO:HD3	2.03	0.41
2:B:1001:FAD:C5'	3:B:1002:ACT:H2	2.50	0.41
1:A:306:ASN:HD22	1:A:306:ASN:HA	1.67	0.41
1:B:47:SER:HA	2:B:1001:FAD:C6	2.50	0.41
1:A:55:PHE:CD2	1:A:60:ASN:HB3	2.56	0.41
1:B:53:THR:CG2	1:B:154:SER:OG	2.69	0.41
1:B:83:LEU:C	1:B:83:LEU:HD23	2.41	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:361:VAL:HG12	1:B:363:PHE:CD2	2.56	0.41
1:A:225:LYS:HD2	1:A:395:ALA:O	2.21	0.41
1:A:208:ARG:HE	1:A:208:ARG:HB2	1.59	0.40
1:B:208:ARG:HE	1:B:208:ARG:HB2	1.54	0.40
1:B:289:LEU:O	1:B:301:VAL:HA	2.21	0.40
1:A:232:VAL:HG13	3:A:1002:ACT:O	2.20	0.40
1:B:108:GLY:O	1:B:114:TYR:HE2	2.04	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	423/447 (95%)	411 (97%)	12 (3%)	0	100	100
1	B	423/447 (95%)	417 (99%)	6 (1%)	0	100	100
All	All	846/894 (95%)	828 (98%)	18 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	335/354 (95%)	297 (89%)	38 (11%)	6	3

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	335/354 (95%)	305 (91%)	30 (9%)	9	6
All	All	670/708 (95%)	602 (90%)	68 (10%)	7	4

All (68) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	48	MET
1	A	58	THR
1	A	62	LEU
1	A	76	ARG
1	A	85	LYS
1	A	91	VAL
1	A	94	ARG
1	A	96	ARG
1	A	98	ARG
1	A	100	VAL
1	A	101	GLU
1	A	105	ARG
1	A	112	ARG
1	A	120	GLU
1	A	124	ARG
1	A	131	ARG
1	A	132	VAL
1	A	140	VAL
1	A	147	GLU
1	A	195	LEU
1	A	198	GLU
1	A	202	LEU
1	A	206	GLU
1	A	208	ARG
1	A	219	ARG
1	A	261	THR
1	A	269	ARG
1	A	281	LEU
1	A	295	ARG
1	A	305	ASP
1	A	315	GLU
1	A	331	ILE
1	A	338	GLN
1	A	341	ASP
1	A	408	GLU
1	A	420	ARG

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Mol	Chain	Res	Type
1	A	422	GLU
1	A	423	GLN
1	B	48	MET
1	B	76	ARG
1	B	85	LYS
1	B	96	ARG
1	B	98	ARG
1	B	100	VAL
1	B	101	GLU
1	B	109	GLU
1	B	110	GLU
1	B	120	GLU
1	B	122	LEU
1	B	124	ARG
1	B	125	ARG
1	B	131	ARG
1	B	139	GLU
1	B	147	GLU
1	B	195	LEU
1	B	206	GLU
1	B	208	ARG
1	B	260	SER
1	B	269	ARG
1	B	281	LEU
1	B	286	LYS
1	B	314	GLU
1	B	336	LEU
1	B	338	GLN
1	B	420	ARG
1	B	422	GLU
1	B	424	GLU
1	B	425	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	75	GLN
1	A	239	ASN
1	A	306	ASN
1	B	75	GLN
1	B	239	ASN
1	B	306	ASN

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Mol	Chain	Res	Type
1	B	338	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

8 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$
3	ACT	A	1002	-	3,3,3	0.60	0	3,3,3	3.18	2 (66%)
2	FAD	A	1001	-	53,58,58	1.76	13 (24%)	68,89,89	1.54	9 (13%)
4	EDO	B	1004	-	3,3,3	0.56	0	2,2,2	2.36	1 (50%)
4	EDO	A	1003	-	3,3,3	0.94	0	2,2,2	0.40	0
4	EDO	A	1004	-	3,3,3	0.45	0	2,2,2	1.24	0
4	EDO	B	1003	-	3,3,3	1.03	0	2,2,2	0.45	0
3	ACT	B	1002	-	3,3,3	2.55	1 (33%)	3,3,3	1.73	1 (33%)
2	FAD	B	1001	-	53,58,58	1.73	9 (16%)	68,89,89	1.60	10 (14%)

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	1001	-	-	2/30/50/50	0/6/6/6
4	EDO	B	1004	-	-	0/1/1/1	-
4	EDO	A	1003	-	-	0/1/1/1	-
4	EDO	A	1004	-	-	0/1/1/1	-
4	EDO	B	1003	-	-	0/1/1/1	-
2	FAD	B	1001	-	-	2/30/50/50	0/6/6/6

All (23) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1001	FAD	C4X-N5	5.77	1.42	1.30
2	B	1001	FAD	C4X-N5	5.64	1.41	1.30
2	B	1001	FAD	C10-N1	5.15	1.43	1.33
2	A	1001	FAD	C10-N1	4.98	1.43	1.33
2	B	1001	FAD	C2A-N3A	4.73	1.39	1.32
3	B	1002	ACT	O-C	4.19	1.41	1.22
2	A	1001	FAD	O4B-C1B	3.81	1.46	1.41
2	A	1001	FAD	C2A-N1A	3.55	1.40	1.33
2	B	1001	FAD	C2A-N1A	3.50	1.40	1.33
2	A	1001	FAD	C2A-N3A	3.35	1.37	1.32
2	B	1001	FAD	P-O2P	-2.53	1.43	1.55
2	A	1001	FAD	P-O2P	-2.47	1.43	1.55
2	A	1001	FAD	O4'-C4'	-2.41	1.38	1.43
2	B	1001	FAD	C8A-N7A	2.25	1.38	1.34
2	B	1001	FAD	O4B-C1B	2.24	1.44	1.41
2	A	1001	FAD	C10-N10	2.18	1.42	1.37
2	A	1001	FAD	C1'-N10	2.11	1.53	1.48
2	A	1001	FAD	C5A-N7A	-2.11	1.32	1.39
2	B	1001	FAD	C1'-N10	2.10	1.53	1.48
2	A	1001	FAD	C5'-C4'	-2.05	1.48	1.51
2	A	1001	FAD	C2-N1	2.02	1.41	1.36
2	B	1001	FAD	C5A-N7A	-2.02	1.32	1.39
2	A	1001	FAD	C8A-N7A	2.01	1.38	1.34

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1001	FAD	N3A-C2A-N1A	-4.81	121.15	128.68
2	B	1001	FAD	N3A-C2A-N1A	-4.63	121.44	128.68
3	A	1002	ACT	OXT-C-CH3	4.10	132.14	115.18
2	B	1001	FAD	C9A-C5X-N5	-3.97	118.11	122.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1001	FAD	C9A-C5X-N5	-3.88	118.22	122.43
3	A	1002	ACT	O-C-CH3	-3.62	108.24	122.33
2	B	1001	FAD	C5'-C4'-C3'	-3.42	105.59	112.20
2	B	1001	FAD	C4X-C4-N3	3.31	121.59	113.19
2	A	1001	FAD	C4A-C5A-N7A	-3.24	106.02	109.40
4	B	1004	EDO	O2-C2-C1	-3.23	88.69	111.91
2	A	1001	FAD	C4X-C4-N3	3.17	121.24	113.19
2	B	1001	FAD	C4A-C5A-N7A	-3.07	106.20	109.40
2	B	1001	FAD	O4'-C4'-C5'	-2.90	103.40	109.92
2	A	1001	FAD	C5'-C4'-C3'	-2.87	106.65	112.20
2	A	1001	FAD	C4-N3-C2	-2.64	120.76	125.64
2	A	1001	FAD	O4-C4-C4X	-2.37	120.30	126.60
2	B	1001	FAD	C4X-C10-N1	-2.37	119.24	124.73
2	A	1001	FAD	C5X-C9A-N10	2.35	120.38	117.95
2	B	1001	FAD	C4-N3-C2	-2.28	121.43	125.64
2	B	1001	FAD	O4-C4-C4X	-2.26	120.60	126.60
3	B	1002	ACT	OXT-C-CH3	-2.26	105.85	115.18
2	B	1001	FAD	C5X-C9A-N10	2.17	120.19	117.95
2	A	1001	FAD	O3B-C3B-C4B	-2.08	105.04	111.05

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	1001	FAD	O4B-C4B-C5B-O5B
2	A	1001	FAD	O4B-C4B-C5B-O5B
2	A	1001	FAD	O3'-C3'-C4'-C5'
2	B	1001	FAD	O3'-C3'-C4'-C5'

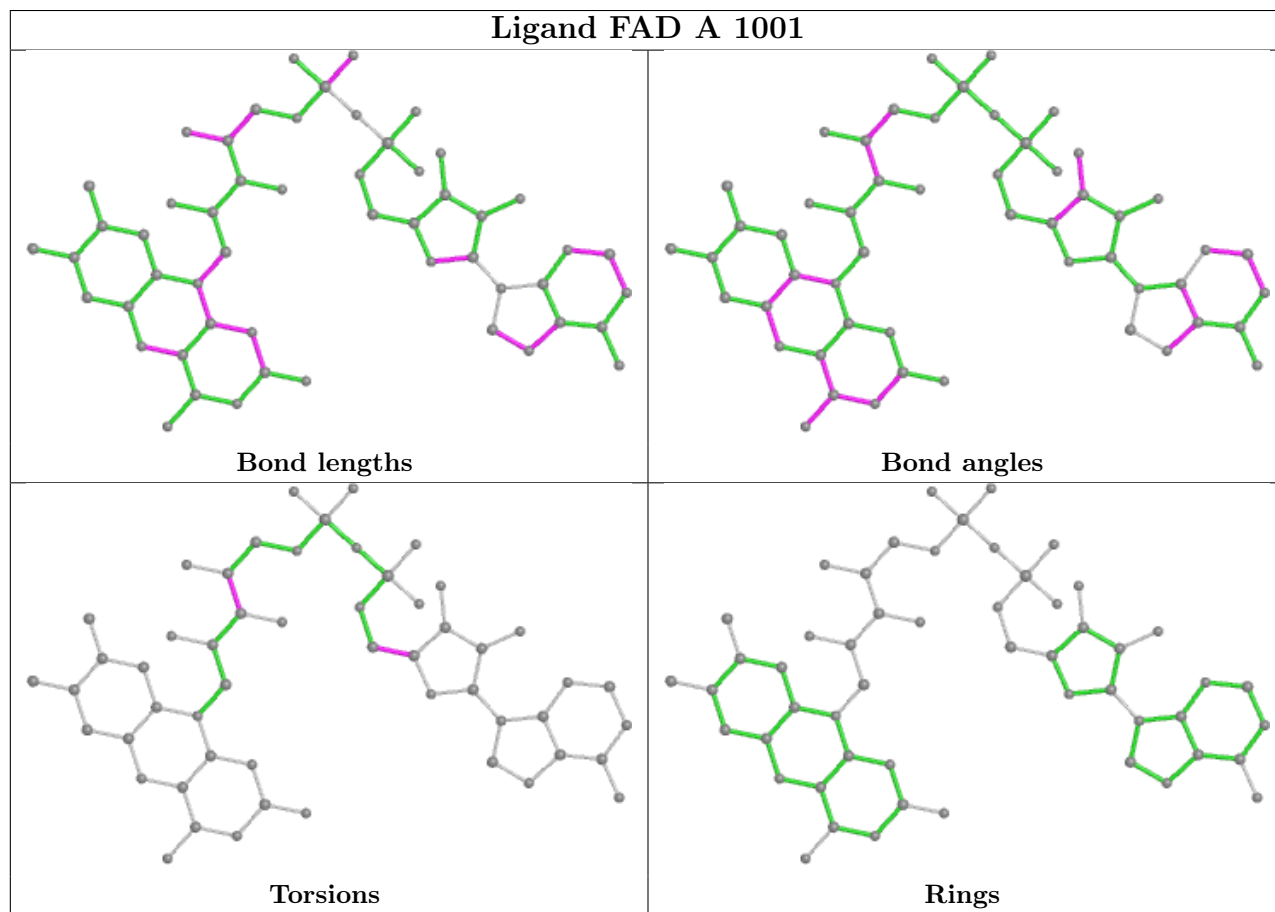
There are no ring outliers.

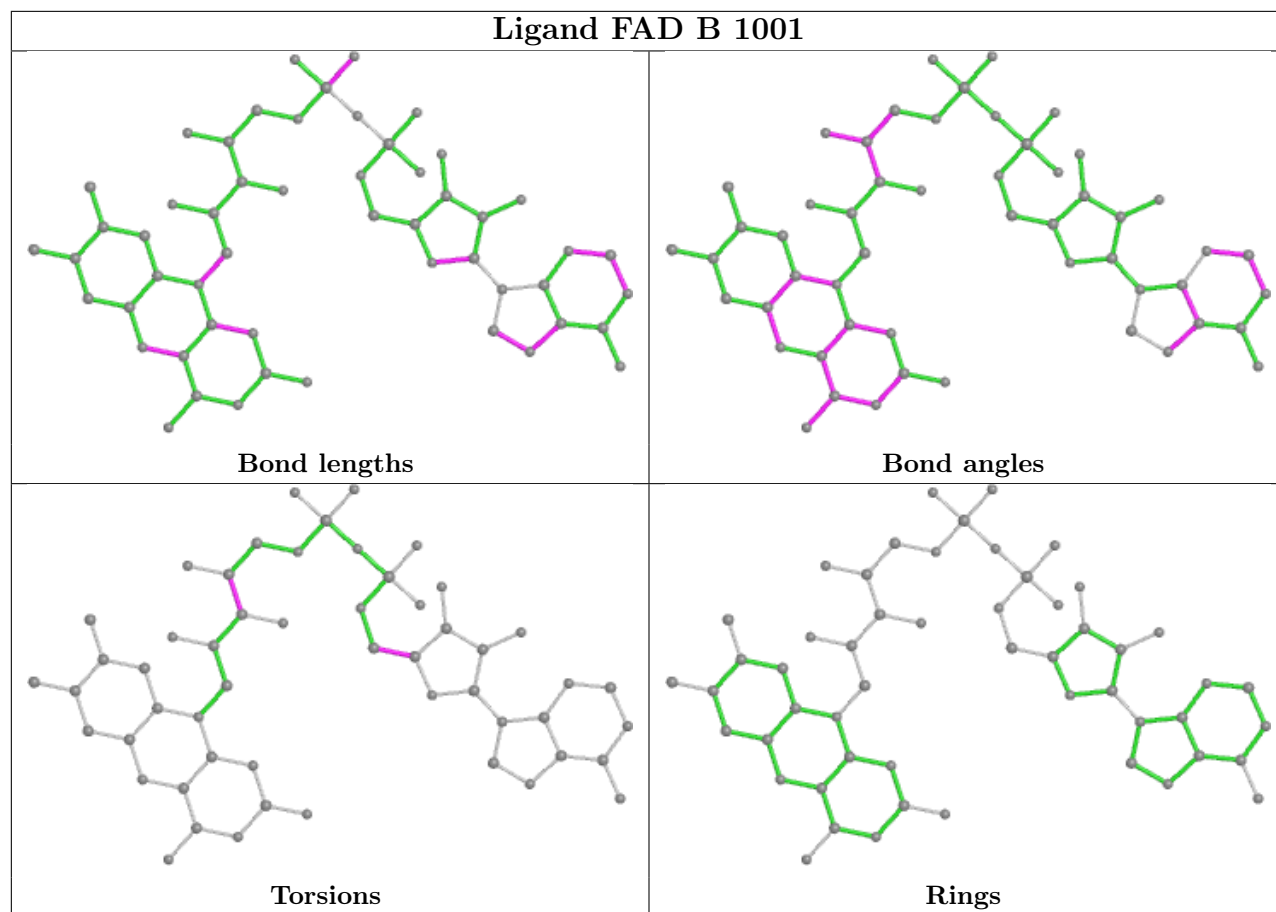
4 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1002	ACT	6	0
4	B	1004	EDO	2	0
3	B	1002	ACT	1	0
2	B	1001	FAD	3	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will

also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





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

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	425/447 (95%)	0.07	9 (2%) 63 62	16, 35, 67, 93	0
1	B	425/447 (95%)	0.08	13 (3%) 49 48	16, 35, 68, 99	0
All	All	850/894 (95%)	0.07	22 (2%) 56 54	16, 35, 68, 99	0

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	269	ARG	3.5
1	A	131	ARG	3.4
1	A	98	ARG	3.4
1	A	97	TRP	3.0
1	A	425	ARG	2.9
1	A	340	GLN	2.7
1	A	107	ALA	2.7
1	B	131	ARG	2.5
1	B	136	GLU	2.5
1	B	297	GLY	2.5
1	B	111	GLY	2.4
1	A	127	GLY	2.4
1	B	103	PRO	2.3
1	B	148	GLY	2.3
1	B	425	ARG	2.2
1	B	97	TRP	2.2
1	A	148	GLY	2.1
1	B	140	VAL	2.1
1	A	269	ARG	2.1
1	B	262	GLU	2.1
1	B	98	ARG	2.0
1	B	94	ARG	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 monosaccharides in this entry.

6.4 Ligands [i](#)

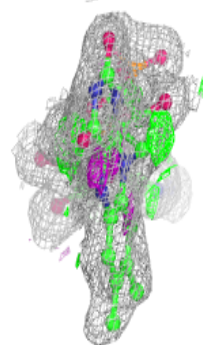
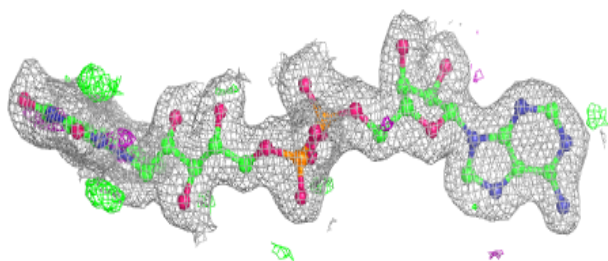
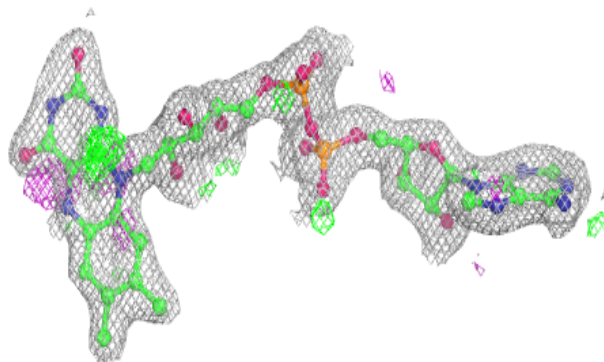
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	ACT	B	1002	4/4	0.71	0.18	26,30,34,49	0
3	ACT	A	1002	4/4	0.76	0.28	20,25,36,52	0
4	EDO	A	1003	4/4	0.93	0.12	18,21,22,22	0
4	EDO	A	1004	4/4	0.93	0.13	21,27,27,32	0
4	EDO	B	1003	4/4	0.93	0.09	18,20,21,22	0
2	FAD	B	1001	53/53	0.95	0.11	16,20,31,41	0
2	FAD	A	1001	53/53	0.97	0.11	15,20,32,42	0
4	EDO	B	1004	4/4	0.98	0.10	20,29,37,42	0

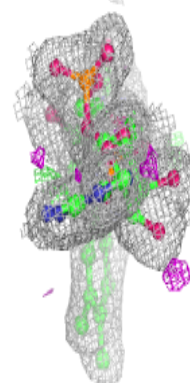
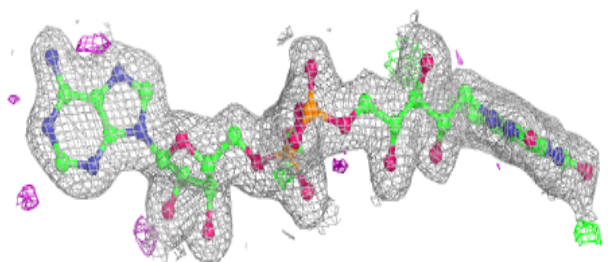
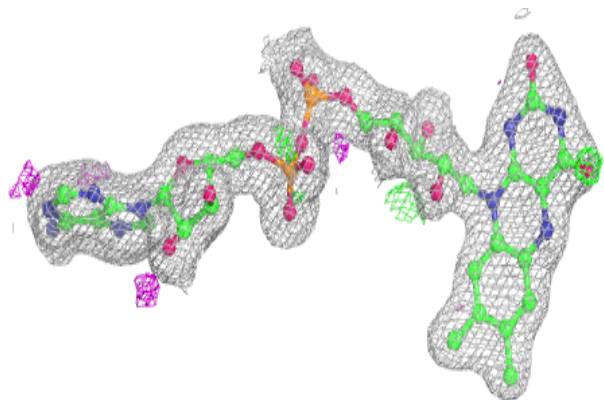
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around FAD B 1001:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around FAD A 1001:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.