



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 07:19 PM UTC

PDB ID : 2RGO / pdb_00002rgo
Title : Structure of Alpha-Glycerophosphate Oxidase from Streptococcus sp.: A Template for the Mitochondrial Alpha-Glycerophosphate Dehydrogenase
Authors : Colussi, T.; Boles, W.; Mallett, T.C.; Karplus, P.A.; Claiborne, A.
Deposited on : 2007-10-04
Resolution : 2.40 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 : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

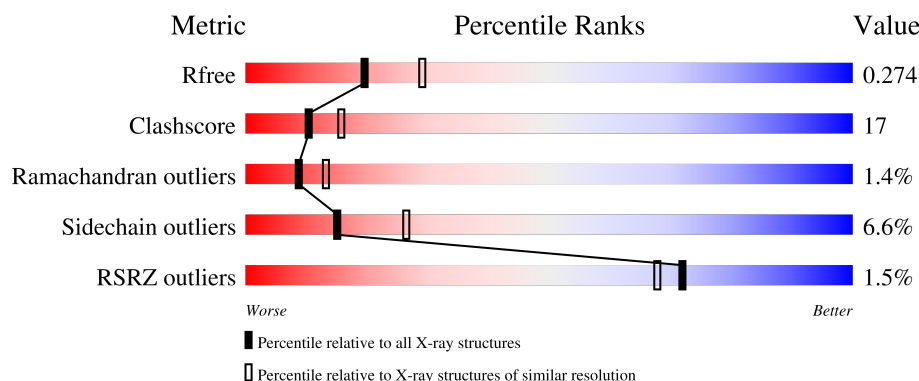
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

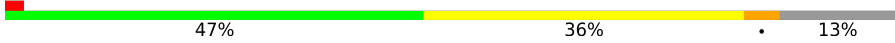
The reported resolution of this entry is 2.40 Å.

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}	180053	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	607	
1	B	607	

2 Entry composition [i](#)

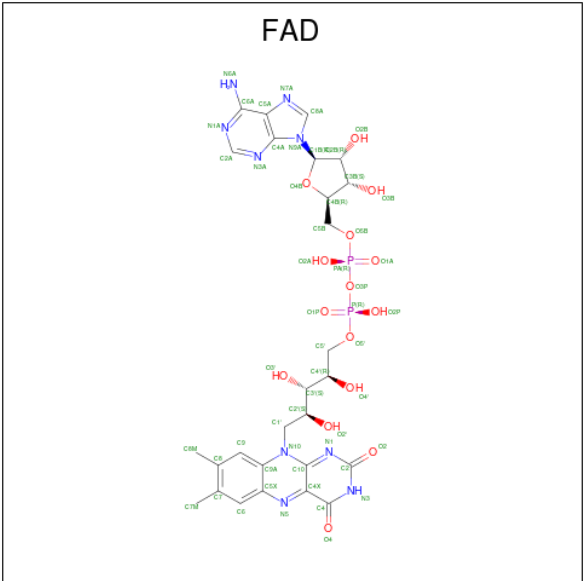
There are 3 unique types of molecules in this entry. The entry contains 8862 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Alpha-Glycerophosphate Oxidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	557	Total	C	N	O	S	0	0	0
			4361	2759	733	858	11			
1	B	530	Total	C	N	O	S	0	0	0
			4163	2640	701	811	11			

- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (CCD ID: 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		

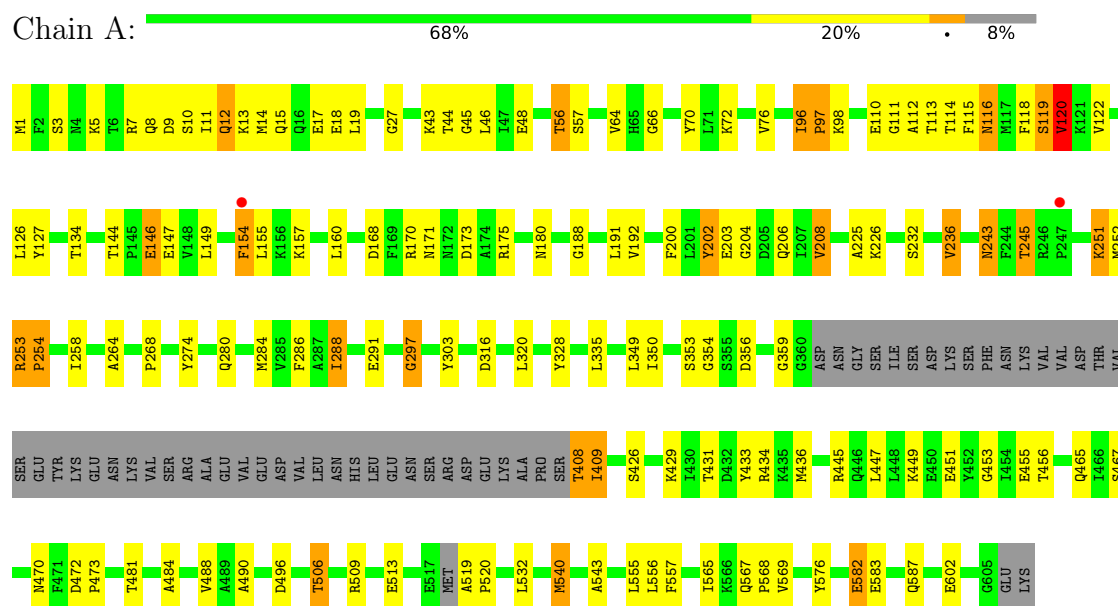
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	183	Total 183	O 183	0	0
3	B	49	Total 49	O 49	0	0

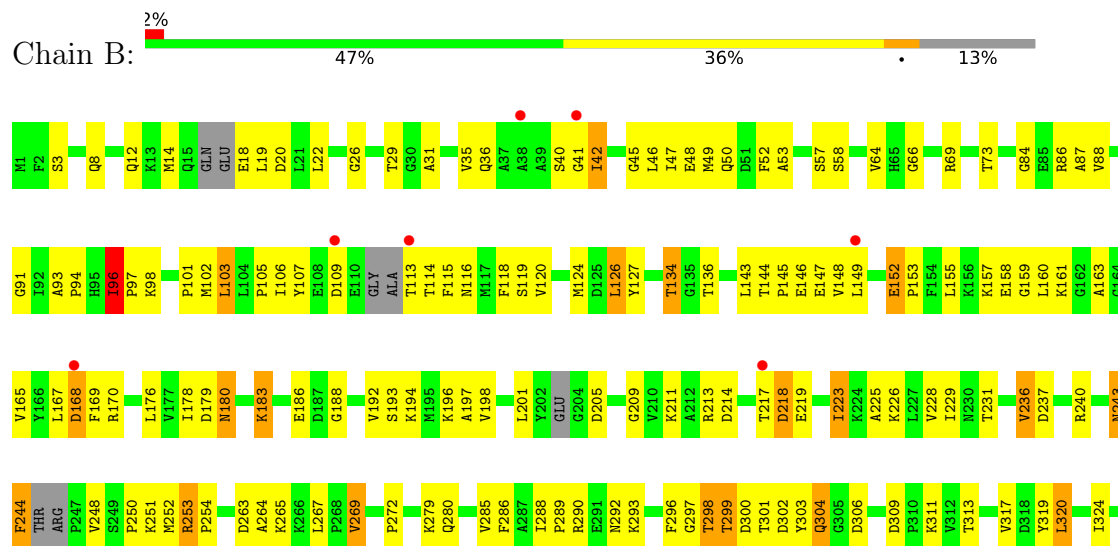
3 Residue-property plots

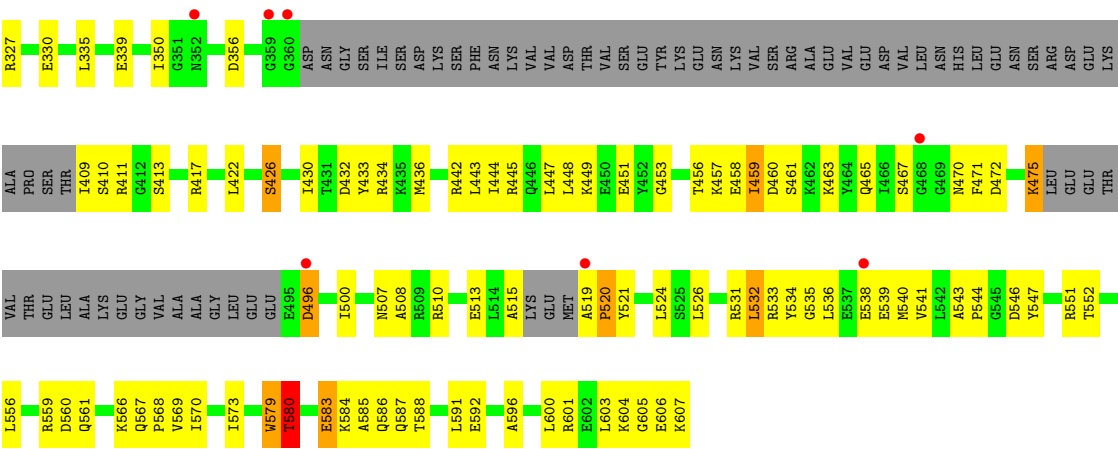
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Alpha-Glycerophosphate Oxidase



• Molecule 1: Alpha-Glycerophosphate Oxidase





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	128.36Å 106.78Å 58.89Å 90.00° 99.18° 90.00°	Depositor
Resolution (Å)	12.00 – 2.40 12.00 – 2.40	Depositor EDS
% Data completeness (in resolution range)	94.9 (12.00-2.40) 94.1 (12.00-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.81 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.2	Depositor
R, R_{free}	0.238 , 0.247 0.230 , 0.274	Depositor DCC
R_{free} test set	2883 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	47.6	Xtriage
Anisotropy	0.549	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 75.4	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	0.94	EDS
Total number of atoms	8862	wwPDB-VP
Average B, all atoms (Å ²)	62.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.40% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 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	0.88	1/4434 (0.0%)	1.16	38/5991 (0.6%)
1	B	0.61	0/4231	1.03	20/5706 (0.4%)
All	All	0.76	1/8665 (0.0%)	1.10	58/11697 (0.5%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2
1	B	0	2
All	All	0	4

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	97	PRO	CA-C	5.41	1.58	1.52

All (58) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	253	ARG	CA-C-N	8.96	128.99	119.76
1	A	253	ARG	C-N-CA	8.96	128.99	119.76
1	A	188	GLY	N-CA-C	7.76	125.01	114.92
1	B	253	ARG	CA-C-N	7.71	127.72	119.78
1	B	253	ARG	C-N-CA	7.71	127.72	119.78
1	B	538	GLU	N-CA-C	7.25	124.19	114.12
1	A	180	ASN	N-CA-C	-7.04	103.02	111.69
1	A	204	GLY	N-CA-C	-6.98	102.87	111.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	44	THR	N-CA-C	6.95	120.84	109.24
1	A	297	GLY	O-C-N	-6.82	115.43	123.55
1	A	496	ASP	N-CA-C	6.80	118.69	111.28
1	B	152	GLU	CA-C-N	6.71	126.97	119.32
1	B	152	GLU	C-N-CA	6.71	126.97	119.32
1	A	409	ILE	N-CA-C	-6.65	106.37	111.62
1	A	252	MET	N-CA-C	6.62	120.31	110.52
1	A	206	GLN	N-CA-C	6.56	119.42	108.20
1	A	48	GLU	CA-C-N	6.43	128.90	120.28
1	A	48	GLU	C-N-CA	6.43	128.90	120.28
1	A	119	SER	N-CA-C	6.24	119.06	111.82
1	A	27	GLY	N-CA-C	-6.12	103.85	112.60
1	B	196	LYS	N-CA-C	5.95	119.33	109.46
1	B	192	VAL	N-CA-C	5.93	116.15	106.72
1	A	543	ALA	CA-C-N	5.92	127.24	119.84
1	A	543	ALA	C-N-CA	5.92	127.24	119.84
1	A	232	SER	N-CA-C	5.82	119.45	112.93
1	B	180	ASN	N-CA-C	-5.79	104.41	112.45
1	A	236	VAL	N-CA-C	5.78	116.52	110.62
1	A	254	PRO	N-CA-C	5.77	120.15	111.14
1	A	556	LEU	N-CA-C	5.75	117.34	111.14
1	A	354	GLY	N-CA-C	-5.75	107.21	114.16
1	B	169	PHE	N-CA-C	5.66	118.44	109.50
1	A	303	TYR	N-CA-C	5.63	117.96	109.23
1	A	3	SER	N-CA-C	5.63	115.54	108.45
1	B	426	SER	N-CA-C	5.57	118.33	109.81
1	A	353	SER	N-CA-C	5.52	117.58	109.24
1	A	120	VAL	CB-CA-C	-5.47	103.69	112.16
1	B	205	ASP	N-CA-C	-5.46	106.47	114.39
1	B	556	LEU	N-CA-C	5.45	117.30	111.36
1	A	587	GLN	N-CA-C	-5.44	105.51	111.82
1	B	126	LEU	N-CA-C	-5.41	105.46	111.36
1	A	76	VAL	N-CA-C	5.41	117.85	111.09
1	A	192	VAL	N-CA-C	5.40	115.64	107.75
1	B	311	LYS	N-CA-C	5.40	118.36	109.72
1	A	451	GLU	N-CA-C	5.39	119.58	112.89
1	A	555	LEU	N-CA-C	5.38	116.83	110.97
1	A	202	TYR	N-CA-C	5.36	117.33	109.24
1	A	264	ALA	N-CA-C	5.34	117.85	111.71
1	B	583	GLU	N-CA-C	-5.32	105.65	111.82
1	B	248	VAL	N-CA-C	5.32	115.02	108.53
1	B	269	VAL	CA-C-N	5.32	125.03	119.82

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	269	VAL	C-N-CA	5.32	125.03	119.82
1	A	426	SER	N-CA-C	5.31	118.25	109.85
1	B	472	ASP	CA-C-N	5.27	126.43	119.84
1	B	472	ASP	C-N-CA	5.27	126.43	119.84
1	A	291	GLU	CB-CA-C	-5.21	109.05	116.34
1	A	557	PHE	N-CA-C	5.19	119.76	113.38
1	A	467	SER	N-CA-C	5.18	117.85	110.24
1	A	540	MET	N-CA-C	5.18	118.58	111.54

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	297	GLY	Peptide
1	A	328	TYR	Mainchain
1	B	297	GLY	Peptide
1	B	298	THR	Peptide

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4361	0	4361	88	0
1	B	4163	0	4167	204	0
2	A	53	0	31	4	0
2	B	53	0	31	3	0
3	A	183	0	0	4	0
3	B	49	0	0	2	0
All	All	8862	0	8590	293	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 17.

All (293) 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:144:THR:HG22	1:B:147:GLU:HG2	1.15	1.13
1:B:14:MET:HE2	1:B:223:ILE:HD12	1.50	0.94
1:B:109:ASP:OD2	1:B:159:GLY:HA3	1.70	0.90
1:A:449:LYS:HE2	1:A:455:GLU:HG2	1.53	0.89
1:B:46:LEU:HD23	1:B:47:ILE:N	1.90	0.87
1:B:471:PHE:HE1	1:B:475:LYS:HG2	1.44	0.83
1:B:579:TRP:O	1:B:580:THR:HG23	1.78	0.83
1:B:22:LEU:HD12	1:B:45:GLY:O	1.78	0.82
1:B:144:THR:HG22	1:B:147:GLU:CG	2.05	0.81
1:B:3:SER:HB3	1:B:536:LEU:O	1.79	0.80
1:B:14:MET:HE2	1:B:223:ILE:CD1	2.09	0.80
1:B:460:ASP:OD2	1:B:463:LYS:HB2	1.82	0.79
1:B:566:LYS:O	1:B:569:VAL:HG12	1.82	0.79
1:B:20:ASP:HA	1:B:226:LYS:HD3	1.66	0.77
1:B:228:VAL:O	1:B:422:LEU:HD12	1.84	0.77
1:B:116:ASN:HD22	1:B:118:PHE:H	1.32	0.77
1:A:116:ASN:HD22	1:A:118:PHE:H	1.35	0.74
1:A:203:GLU:O	1:A:203:GLU:HG3	1.86	0.73
1:B:243:ASN:C	1:B:243:ASN:HD22	1.95	0.73
1:A:98:LYS:HG3	1:A:170:ARG:NH1	2.03	0.72
1:B:144:THR:CG2	1:B:147:GLU:HG2	2.09	0.72
1:B:410:SER:HB3	1:B:413:SER:OG	1.90	0.71
1:B:544:PRO:HG2	1:B:591:LEU:HD21	1.71	0.70
1:B:584:LYS:C	1:B:584:LYS:HD3	2.16	0.70
1:B:96:ILE:HB	1:B:97:PRO:CD	2.21	0.69
1:B:47:ILE:HD11	1:B:197:ALA:HB2	1.73	0.69
1:B:73:THR:HG22	1:B:73:THR:O	1.93	0.69
1:A:251:LYS:O	1:A:350:ILE:HG13	1.94	0.68
1:B:31:ALA:O	1:B:35:VAL:HG23	1.93	0.68
1:B:115:PHE:HB3	1:B:120:VAL:HG13	1.76	0.68
1:A:243:ASN:C	1:A:243:ASN:HD22	2.02	0.68
1:A:116:ASN:HD22	1:A:118:PHE:N	1.92	0.67
1:B:515:ALA:O	1:B:526:LEU:HD11	1.94	0.67
1:B:269:VAL:HG23	1:B:269:VAL:O	1.94	0.66
1:B:443:LEU:O	1:B:443:LEU:HG	1.94	0.65
1:B:106:ILE:HD11	1:B:120:VAL:HG11	1.79	0.65
1:B:584:LYS:HD3	1:B:585:ALA:N	2.11	0.65
1:A:116:ASN:ND2	1:A:118:PHE:H	1.95	0.65
1:B:103:LEU:HD12	1:B:165:VAL:HG22	1.78	0.64
1:B:532:LEU:HD11	1:B:569:VAL:HG23	1.79	0.64
1:B:47:ILE:HG13	1:B:47:ILE:O	1.98	0.64
1:B:237:ASP:OD1	1:B:240:ARG:NH1	2.31	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:510:ARG:HH21	1:B:513:GLU:HB2	1.63	0.63
1:B:14:MET:HG2	1:B:19:LEU:HD11	1.81	0.63
1:B:298:THR:OG1	1:B:299:THR:N	2.31	0.63
1:B:335:LEU:O	1:B:335:LEU:HD23	1.99	0.63
1:B:532:LEU:HD21	1:B:573:ILE:HD11	1.81	0.62
1:A:116:ASN:ND2	1:A:119:SER:H	1.98	0.62
1:A:203:GLU:HG2	1:A:208:VAL:HG11	1.82	0.62
1:B:14:MET:CE	1:B:223:ILE:HD12	2.28	0.62
1:A:509:ARG:O	1:A:513:GLU:HG3	2.00	0.62
1:B:569:VAL:O	1:B:573:ILE:HG12	2.00	0.61
1:B:579:TRP:CG	1:B:583:GLU:HG2	2.34	0.61
1:A:445:ARG:HD3	1:A:456:THR:OG1	2.00	0.61
1:B:109:ASP:CG	1:B:159:GLY:HA3	2.26	0.61
1:B:471:PHE:CE1	1:B:475:LYS:HG2	2.32	0.61
1:B:40:SER:OG	1:B:459:ILE:HG12	2.00	0.61
1:A:258:ILE:HD13	1:A:316:ASP:HB3	1.82	0.61
1:B:217:THR:O	1:B:219:GLU:N	2.31	0.61
1:A:116:ASN:HD21	1:A:118:PHE:HB3	1.66	0.61
1:A:98:LYS:HG3	1:A:170:ARG:HH12	1.65	0.60
1:B:145:PRO:HG3	1:B:161:LYS:C	2.26	0.60
1:B:243:ASN:C	1:B:243:ASN:ND2	2.57	0.60
1:B:19:LEU:O	1:B:225:ALA:HA	2.00	0.60
1:A:154:PHE:CZ	1:A:268:PRO:O	2.55	0.60
1:A:116:ASN:ND2	1:A:118:PHE:N	2.50	0.59
1:A:114:THR:HG23	1:A:115:PHE:CD1	2.37	0.59
1:B:52:PHE:CZ	1:B:193:SER:HB3	2.38	0.59
1:A:449:LYS:CE	1:A:455:GLU:HG2	2.29	0.59
1:B:18:GLU:OE2	1:B:226:LYS:HG3	2.02	0.59
1:B:263:ASP:OD2	1:B:265:LYS:HG2	2.02	0.59
1:B:98:LYS:HB2	1:B:170:ARG:HD3	1.83	0.59
1:B:567:GLN:HB3	1:B:568:PRO:HD3	1.85	0.59
1:A:519:ALA:N	1:A:520:PRO:HD3	2.18	0.58
1:B:475:LYS:HD2	1:B:475:LYS:C	2.28	0.58
1:A:286:PHE:HB3	1:A:288:ILE:HD11	1.84	0.58
1:B:417:ARG:CZ	1:B:447:LEU:HD21	2.33	0.58
1:A:46:LEU:HB3	1:A:191:LEU:HD23	1.83	0.58
1:B:179:ASP:O	1:B:467:SER:HB3	2.04	0.58
1:B:432:ASP:O	1:B:436:MET:HG3	2.04	0.58
1:A:540:MET:O	1:A:540:MET:HG3	2.03	0.57
1:A:56:THR:HG23	2:A:1001:FAD:O2A	2.04	0.57
1:B:288:ILE:HG23	1:B:289:PRO:HD2	1.86	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:217:THR:C	1:B:219:GLU:H	2.12	0.57
1:A:408:THR:HG23	1:A:409:ILE:H	1.68	0.57
1:B:531:ARG:NH2	1:B:552:THR:HG22	2.19	0.57
1:A:9:ASP:O	1:A:13:LYS:HD3	2.05	0.57
1:B:540:MET:HG3	1:B:540:MET:O	2.05	0.57
1:B:116:ASN:ND2	1:B:119:SER:H	2.03	0.57
1:A:66:GLY:HA3	1:A:127:TYR:OH	2.05	0.56
1:B:586:GLN:HA	1:B:586:GLN:NE2	2.21	0.56
1:B:496:ASP:O	1:B:500:ILE:CD1	2.54	0.56
1:A:154:PHE:CD1	1:A:154:PHE:N	2.73	0.56
1:A:115:PHE:CB	1:A:120:VAL:HG13	2.36	0.56
1:A:433:TYR:HA	1:A:436:MET:HE3	1.88	0.56
1:A:116:ASN:O	1:A:120:VAL:HG22	2.06	0.55
1:A:115:PHE:HB2	1:A:120:VAL:HG13	1.87	0.55
1:B:36:GLN:CD	1:B:461:SER:HB3	2.30	0.55
1:B:519:ALA:N	1:B:520:PRO:HD3	2.22	0.55
1:B:417:ARG:NH1	1:B:447:LEU:HD11	2.22	0.55
1:B:596:ALA:HB2	1:B:601:ARG:HH21	1.73	0.55
1:B:201:LEU:HD12	1:B:209:GLY:HA3	1.89	0.54
1:A:114:THR:HG23	1:A:115:PHE:N	2.22	0.54
1:B:144:THR:HG23	1:B:146:GLU:H	1.72	0.54
1:B:579:TRP:O	1:B:580:THR:CG2	2.55	0.54
1:B:586:GLN:HA	1:B:586:GLN:HE21	1.72	0.54
1:B:96:ILE:HB	1:B:97:PRO:HD3	1.90	0.54
1:B:583:GLU:O	1:B:587:GLN:HG2	2.08	0.54
1:B:551:ARG:HG2	1:B:551:ARG:HH11	1.73	0.53
1:A:175:ARG:HB3	3:A:1076:HOH:O	2.09	0.53
1:A:149:LEU:HD21	1:A:160:LEU:HD23	1.90	0.52
1:B:105:PRO:HA	1:B:163:ALA:HB2	1.91	0.52
1:A:14:MET:HE3	1:A:45:GLY:HA3	1.90	0.52
1:B:46:LEU:HD23	1:B:46:LEU:C	2.34	0.52
1:B:113:THR:OG1	1:B:114:THR:N	2.42	0.52
1:B:252:MET:O	1:B:350:ILE:HD11	2.09	0.52
1:A:116:ASN:HD22	1:A:116:ASN:C	2.17	0.52
1:B:97:PRO:HA	1:B:170:ARG:O	2.10	0.51
1:B:448:LEU:HD12	1:B:456:THR:HG21	1.92	0.51
1:A:8:GLN:O	1:A:12:GLN:HG2	2.09	0.51
1:B:457:LYS:HG2	1:B:458:GLU:H	1.75	0.51
1:B:145:PRO:O	1:B:149:LEU:HD23	2.10	0.51
1:B:496:ASP:O	1:B:500:ILE:HG13	2.10	0.51
1:B:236:VAL:HG21	1:B:426:SER:HB3	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:244:PHE:HD1	1:B:244:PHE:H	1.58	0.51
1:A:96:ILE:HB	1:A:97:PRO:CD	2.41	0.51
1:B:8:GLN:O	1:B:12:GLN:HG2	2.10	0.51
1:B:335:LEU:HD22	1:B:604:LYS:HG2	1.91	0.50
1:A:582:GLU:HG2	1:A:583:GLU:N	2.25	0.50
1:B:449:LYS:HA	1:B:453:GLY:H	1.76	0.50
1:B:600:LEU:HB3	1:B:603:LEU:HD12	1.94	0.50
1:A:484:ALA:O	1:A:488:VAL:HG23	2.12	0.50
1:A:203:GLU:O	1:A:203:GLU:CG	2.57	0.50
1:A:465:GLN:OE1	1:A:470:ASN:HA	2.11	0.50
1:A:96:ILE:HB	1:A:97:PRO:HD2	1.94	0.50
1:B:102:MET:HG3	1:B:272:PRO:HB2	1.93	0.50
1:B:143:LEU:HB3	1:B:148:VAL:HG23	1.94	0.50
1:B:604:LYS:HB3	1:B:607:LYS:OXT	2.12	0.49
1:A:18:GLU:OE1	1:A:226:LYS:HG3	2.11	0.49
1:B:604:LYS:HD2	1:B:607:LYS:HA	1.94	0.49
1:A:144:THR:HG23	1:A:147:GLU:OE1	2.13	0.49
1:A:146:GLU:CD	1:A:146:GLU:H	2.19	0.49
1:B:167:LEU:HD12	1:B:168:ASP:H	1.76	0.49
1:A:243:ASN:HD21	1:A:245:THR:HB	1.78	0.49
1:A:154:PHE:HZ	1:A:268:PRO:O	1.95	0.49
1:A:243:ASN:HD22	1:A:245:THR:H	1.59	0.49
1:B:253:ARG:NH1	1:B:253:ARG:HG2	2.27	0.49
1:B:120:VAL:O	1:B:124:MET:HG3	2.12	0.49
1:B:48:GLU:HG3	1:B:50:GLN:O	2.13	0.48
1:B:41:GLY:O	1:B:42:ILE:C	2.56	0.48
1:B:98:LYS:HB2	1:B:170:ARG:CG	2.44	0.48
1:A:11:ILE:O	1:A:15:GLN:HG3	2.13	0.48
1:B:288:ILE:CG2	1:B:289:PRO:HD2	2.44	0.48
1:B:320:LEU:O	1:B:324:ILE:HD12	2.14	0.48
1:B:457:LYS:HG2	1:B:458:GLU:N	2.29	0.48
1:B:98:LYS:HB2	1:B:170:ARG:CD	2.44	0.48
1:B:510:ARG:HH21	1:B:513:GLU:CB	2.26	0.48
1:A:70:TYR:CZ	1:A:359:GLY:HA2	2.49	0.48
1:B:532:LEU:CD1	1:B:569:VAL:HG23	2.43	0.47
1:B:605:GLY:O	1:B:606:GLU:HB3	2.13	0.47
1:A:7:ARG:HD3	3:A:1080:HOH:O	2.15	0.47
1:A:436:MET:HE1	3:A:1154:HOH:O	2.13	0.47
1:B:279:LYS:NZ	1:B:319:TYR:HA	2.29	0.47
1:B:449:LYS:C	1:B:451:GLU:H	2.22	0.47
1:B:500:ILE:HG22	1:B:508:ALA:HB1	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:19:LEU:O	1:A:225:ALA:HA	2.15	0.47
1:B:250:PRO:O	1:B:251:LYS:HD2	2.15	0.47
1:A:154:PHE:HD1	1:A:154:PHE:H	1.55	0.47
1:B:244:PHE:CD1	1:B:244:PHE:N	2.83	0.47
1:A:116:ASN:ND2	1:A:116:ASN:C	2.74	0.46
1:B:531:ARG:O	1:B:547:TYR:HE1	1.98	0.46
1:A:57:SER:O	1:A:173:ASP:HB2	2.15	0.46
1:B:22:LEU:HD13	1:B:223:ILE:HG21	1.97	0.46
1:B:465:GLN:OE1	1:B:470:ASN:O	2.34	0.46
1:A:208:VAL:O	1:A:208:VAL:HG23	2.13	0.46
1:B:229:ILE:HG23	1:B:229:ILE:O	2.16	0.46
1:B:198:VAL:HG22	1:B:211:LYS:O	2.16	0.46
1:B:579:TRP:HB3	1:B:583:GLU:HG2	1.96	0.46
1:A:243:ASN:ND2	1:A:245:THR:H	2.13	0.46
1:A:429:LYS:HB2	2:A:1001:FAD:O2	2.16	0.46
1:B:26:GLY:HA3	1:B:48:GLU:HB2	1.98	0.46
1:B:116:ASN:HD22	1:B:118:PHE:N	2.07	0.46
1:B:178:ILE:HG23	1:B:179:ASP:N	2.31	0.46
1:B:69:ARG:HB2	3:B:1036:HOH:O	2.16	0.46
1:B:73:THR:O	1:B:73:THR:CG2	2.62	0.46
1:B:604:LYS:HZ2	1:B:607:LYS:CE	2.29	0.46
1:B:543:ALA:HB1	1:B:544:PRO:HD2	1.97	0.46
1:B:579:TRP:HB3	1:B:583:GLU:CG	2.46	0.46
1:B:107:TYR:CE2	1:B:327:ARG:NH1	2.84	0.46
1:A:64:VAL:HG23	1:A:171:ASN:HB2	1.97	0.45
1:B:303:TYR:C	1:B:304:GLN:NE2	2.74	0.45
1:B:285:VAL:CG1	1:B:286:PHE:N	2.80	0.45
1:B:604:LYS:HD2	1:B:607:LYS:HD2	1.98	0.45
1:A:567:GLN:HB3	1:A:568:PRO:HD3	1.98	0.45
1:B:29:THR:CG2	1:B:231:THR:HB	2.47	0.45
1:B:183:LYS:HZ3	1:B:183:LYS:HB2	1.82	0.45
1:B:293:LYS:HE3	1:B:339:GLU:OE1	2.17	0.45
1:B:66:GLY:HA3	1:B:127:TYR:OH	2.16	0.45
1:B:507:ASN:HB3	1:B:534:TYR:OH	2.17	0.45
1:A:253:ARG:N	1:A:254:PRO:HD3	2.30	0.45
1:B:19:LEU:HB2	1:B:225:ALA:HB2	1.99	0.45
1:B:433:TYR:CG	1:B:434:ARG:N	2.85	0.45
1:B:35:VAL:HG21	1:B:180:ASN:OD1	2.17	0.44
1:B:313:THR:O	1:B:317:VAL:HG23	2.16	0.44
1:A:565:ILE:O	1:A:569:VAL:HG23	2.17	0.44
1:B:535:GLY:O	1:B:541:VAL:HG23	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:472:ASP:HA	1:A:473:PRO:HD2	1.89	0.44
1:B:64:VAL:HG11	1:B:86:ARG:HG3	1.98	0.44
1:B:116:ASN:HD21	1:B:118:PHE:HB3	1.83	0.44
1:B:324:ILE:HD12	1:B:324:ILE:H	1.83	0.44
1:B:269:VAL:O	1:B:269:VAL:CG2	2.65	0.44
1:B:569:VAL:HG13	1:B:570:ILE:N	2.33	0.44
1:B:579:TRP:C	1:B:580:THR:OG1	2.60	0.44
1:A:5:LYS:HA	1:A:5:LYS:HD3	1.82	0.44
1:A:10:SER:O	1:A:14:MET:HG3	2.18	0.43
1:B:320:LEU:O	1:B:324:ILE:CD1	2.66	0.43
1:A:576:TYR:CD1	1:A:576:TYR:C	2.96	0.43
1:B:253:ARG:HG2	1:B:253:ARG:HH11	1.84	0.43
1:B:49:MET:HB2	2:B:1001:FAD:C2A	2.48	0.43
1:B:442:ARG:C	1:B:444:ILE:H	2.27	0.43
1:A:157:LYS:HE3	3:A:1104:HOH:O	2.18	0.43
1:A:433:TYR:CG	1:A:434:ARG:N	2.87	0.43
1:B:532:LEU:HD21	1:B:573:ILE:CD1	2.47	0.43
1:B:296:PHE:CD2	1:B:324:ILE:HD11	2.54	0.43
1:B:350:ILE:HD12	1:B:350:ILE:N	2.34	0.43
1:B:449:LYS:C	1:B:451:GLU:N	2.74	0.43
1:B:250:PRO:C	1:B:251:LYS:HD2	2.44	0.43
1:A:116:ASN:HD21	1:A:119:SER:H	1.67	0.43
1:A:470:ASN:OD1	1:A:506:THR:HG21	2.19	0.43
1:B:84:GLY:O	1:B:87:ALA:HB3	2.19	0.43
1:B:116:ASN:ND2	1:B:118:PHE:HB3	2.34	0.43
1:A:114:THR:CG2	1:A:115:PHE:N	2.82	0.42
1:B:324:ILE:HD12	1:B:324:ILE:N	2.34	0.42
1:B:300:ASP:C	1:B:301:THR:HG23	2.45	0.42
1:B:496:ASP:O	1:B:500:ILE:HD12	2.18	0.42
1:A:12:GLN:HE21	1:A:12:GLN:HB3	1.63	0.42
1:A:203:GLU:HG2	1:A:208:VAL:CG1	2.46	0.42
1:B:98:LYS:O	1:B:170:ARG:HG2	2.18	0.42
1:B:167:LEU:HD12	1:B:168:ASP:N	2.34	0.42
1:B:253:ARG:NH2	1:B:302:ASP:OD1	2.51	0.42
2:B:1001:FAD:H1'1	2:B:1001:FAD:H9	1.77	0.42
1:B:53:ALA:O	1:B:58:SER:HB3	2.19	0.42
1:B:579:TRP:CZ3	1:B:587:GLN:HG3	2.55	0.42
1:B:588:THR:O	1:B:592:GLU:HG2	2.19	0.42
1:B:306:ASP:OD2	1:B:309:ASP:N	2.42	0.42
1:B:535:GLY:HA2	1:B:539:GLU:HG3	2.02	0.42
1:B:430:ILE:HG22	2:B:1001:FAD:N1	2.35	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:521:TYR:HB2	1:B:524:LEU:HD12	2.01	0.42
1:A:70:TYR:OH	1:A:356:ASP:O	2.36	0.42
1:B:546:ASP:O	1:B:551:ARG:HG2	2.19	0.42
1:B:579:TRP:CG	1:B:583:GLU:CG	3.03	0.42
1:B:29:THR:HG22	1:B:231:THR:HB	2.02	0.42
1:B:586:GLN:HE21	1:B:586:GLN:CA	2.28	0.42
1:B:445:ARG:NE	1:B:457:LYS:O	2.53	0.42
1:A:349:LEU:O	1:A:350:ILE:C	2.61	0.42
1:B:214:ASP:HB3	1:B:217:THR:O	2.20	0.42
1:A:245:THR:O	1:A:245:THR:CG2	2.67	0.41
1:B:42:ILE:HD12	1:B:448:LEU:HD11	2.02	0.41
1:A:72:LYS:HA	1:A:72:LYS:HD2	1.80	0.41
1:A:408:THR:HG23	1:A:409:ILE:N	2.35	0.41
1:A:429:LYS:HE3	2:A:1001:FAD:O2	2.20	0.41
1:A:274:TYR:CE1	1:A:284:MET:HB3	2.55	0.41
1:A:111:GLY:O	1:A:112:ALA:HB3	2.20	0.41
1:B:304:GLN:NE2	1:B:304:GLN:N	2.69	0.41
1:B:579:TRP:CB	1:B:583:GLU:HG2	2.51	0.41
1:B:579:TRP:CD1	1:B:583:GLU:HG2	2.55	0.41
1:B:157:LYS:O	1:B:159:GLY:N	2.54	0.41
1:B:201:LEU:HB2	1:B:209:GLY:CA	2.50	0.41
1:B:253:ARG:N	1:B:254:PRO:HD3	2.35	0.41
1:B:531:ARG:HH21	1:B:552:THR:HG22	1.84	0.41
1:B:559:ARG:C	1:B:561:GLN:H	2.29	0.41
1:A:431:THR:HG23	2:A:1001:FAD:O2	2.21	0.41
1:B:519:ALA:O	1:B:520:PRO:O	2.38	0.41
1:B:604:LYS:NZ	1:B:607:LYS:NZ	2.68	0.41
1:B:152:GLU:HA	1:B:153:PRO:HD2	1.84	0.41
1:B:194:LYS:HD3	1:B:194:LYS:HA	1.95	0.41
1:B:217:THR:C	1:B:219:GLU:N	2.77	0.41
1:B:102:MET:HE3	1:B:102:MET:HB2	1.99	0.41
1:B:532:LEU:HA	1:B:547:TYR:CE1	2.56	0.41
1:B:551:ARG:HG2	1:B:551:ARG:NH1	2.34	0.41
1:B:20:ASP:CB	1:B:42:ILE:HG23	2.50	0.41
1:B:134:THR:HG23	1:B:134:THR:O	2.21	0.41
1:B:279:LYS:HZ2	1:B:319:TYR:HA	1.86	0.41
1:A:200:PHE:HB2	1:A:202:TYR:OH	2.22	0.40
1:B:604:LYS:CD	1:B:607:LYS:HA	2.51	0.40
1:A:46:LEU:HD23	1:A:46:LEU:C	2.46	0.40
1:B:53:ALA:HA	3:B:1010:HOH:O	2.20	0.40
1:B:213:ARG:NH2	1:B:218:ASP:HB3	2.35	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:290:ARG:HD3	1:B:290:ARG:HA	1.96	0.40
1:B:93:ALA:N	1:B:94:PRO:CD	2.85	0.40
1:B:264:ALA:HB2	1:B:292:ASN:O	2.21	0.40
1:A:488:VAL:C	1:A:490:ALA:H	2.29	0.40
1:B:88:VAL:O	1:B:91:GLY:N	2.54	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	551/607 (91%)	522 (95%)	27 (5%)	2 (0%)	30	43
1	B	514/607 (85%)	449 (87%)	52 (10%)	13 (2%)	4	5
All	All	1065/1214 (88%)	971 (91%)	79 (7%)	15 (1%)	9	13

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	96	ILE
1	B	158	GLU
1	B	218	ASP
1	B	520	PRO
1	B	579	TRP
1	B	42	ILE
1	B	134	THR
1	B	560	ASP
1	B	580	THR
1	B	57	SER
1	B	96	ILE
1	B	136	THR
1	B	186	GLU

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Mol	Chain	Res	Type
1	B	188	GLY
1	A	453	GLY

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	465/512 (91%)	433 (93%)	32 (7%)	14	24
1	B	445/512 (87%)	417 (94%)	28 (6%)	16	29
All	All	910/1024 (89%)	850 (93%)	60 (7%)	15	26

All (60) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	12	GLN
1	A	17	GLU
1	A	43	LYS
1	A	56	THR
1	A	110	GLU
1	A	113	THR
1	A	116	ASN
1	A	120	VAL
1	A	122	VAL
1	A	126	LEU
1	A	134	THR
1	A	146	GLU
1	A	154	PHE
1	A	155	LEU
1	A	168	ASP
1	A	208	VAL
1	A	236	VAL
1	A	243	ASN
1	A	245	THR
1	A	251	LYS

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Mol	Chain	Res	Type
1	A	280	GLN
1	A	288	ILE
1	A	320	LEU
1	A	335	LEU
1	A	408	THR
1	A	447	LEU
1	A	481	THR
1	A	506	THR
1	A	532	LEU
1	A	582	GLU
1	A	602	GLU
1	B	96	ILE
1	B	101	PRO
1	B	103	LEU
1	B	126	LEU
1	B	155	LEU
1	B	160	LEU
1	B	168	ASP
1	B	176	LEU
1	B	183	LYS
1	B	223	ILE
1	B	236	VAL
1	B	243	ASN
1	B	244	PHE
1	B	267	LEU
1	B	280	GLN
1	B	299	THR
1	B	304	GLN
1	B	320	LEU
1	B	330	GLU
1	B	356	ASP
1	B	409	ILE
1	B	411	ARG
1	B	459	ILE
1	B	475	LYS
1	B	496	ASP
1	B	532	LEU
1	B	533	ARG
1	B	580	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	4	ASN
1	A	8	GLN
1	A	12	GLN
1	A	116	ASN
1	A	171	ASN
1	A	230	ASN
1	A	241	ASN
1	A	243	ASN
1	A	280	GLN
1	A	554	HIS
1	B	8	GLN
1	B	12	GLN
1	B	116	ASN
1	B	243	ASN
1	B	304	GLN
1	B	314	GLN
1	B	332	ASN
1	B	554	HIS
1	B	586	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 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	1001	-	58,58,58	1.91	14 (24%)	85,89,89	1.99	23 (27%)
2	FAD	B	1001	-	58,58,58	1.25	6 (10%)	85,89,89	1.60	18 (21%)

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	-	-	1/34/50/50	0/6/6/6
2	FAD	B	1001	-	-	0/34/50/50	0/6/6/6

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1001	FAD	PA-O3P	-5.09	1.54	1.59
2	B	1001	FAD	C4X-N5	4.41	1.40	1.30
2	A	1001	FAD	C9-C9A	4.13	1.46	1.39
2	A	1001	FAD	C5A-N7A	-4.12	1.31	1.39
2	A	1001	FAD	P-O3P	3.75	1.63	1.59
2	A	1001	FAD	C10-N10	3.67	1.45	1.37
2	B	1001	FAD	C10-N1	3.61	1.40	1.33
2	A	1001	FAD	C9A-N10	3.61	1.47	1.41
2	A	1001	FAD	C10-N1	3.54	1.40	1.33
2	A	1001	FAD	C6-C7	3.34	1.44	1.39
2	B	1001	FAD	C5A-N7A	-3.06	1.33	1.39
2	A	1001	FAD	C6-C5X	3.03	1.44	1.40
2	A	1001	FAD	C4X-N5	2.80	1.36	1.30
2	A	1001	FAD	C7M-C7	2.72	1.56	1.51
2	A	1001	FAD	P-O2P	-2.58	1.43	1.55
2	A	1001	FAD	O2B-C2B	2.40	1.48	1.43
2	B	1001	FAD	C8A-N9A	-2.27	1.33	1.37
2	B	1001	FAD	PA-O3P	2.23	1.61	1.59
2	A	1001	FAD	C8A-N9A	-2.21	1.33	1.37
2	B	1001	FAD	C4A-N9A	-2.04	1.33	1.37

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1001	FAD	N3A-C2A-N1A	-6.47	118.78	128.58
2	A	1001	FAD	O2P-P-O3P	-5.97	91.14	107.27
2	B	1001	FAD	C5A-C4A-N3A	-5.48	119.17	126.72
2	B	1001	FAD	N3A-C2A-N1A	-4.69	121.48	128.58
2	A	1001	FAD	N3A-C4A-N9A	4.30	134.48	127.17
2	A	1001	FAD	C2A-N3A-C4A	3.74	120.96	111.83
2	B	1001	FAD	C2A-N3A-C4A	3.72	120.91	111.83
2	A	1001	FAD	C5A-C4A-N3A	-3.67	121.67	126.72
2	B	1001	FAD	N3A-C4A-N9A	3.61	133.31	127.17
2	A	1001	FAD	C5'-C4'-C3'	-3.53	105.56	112.22
2	A	1001	FAD	C9A-C5X-N5	-3.44	118.80	122.45
2	A	1001	FAD	O5'-C5'-C4'	3.42	118.50	109.36
2	A	1001	FAD	O3P-PA-O1A	3.36	120.81	110.70
2	B	1001	FAD	C4-N3-C2	-3.32	119.74	125.64
2	A	1001	FAD	C2'-C1'-N10	3.15	125.10	110.20
2	A	1001	FAD	O2P-P-O1P	3.04	126.61	112.44
2	A	1001	FAD	C5X-N5-C4X	3.04	123.00	118.09
2	B	1001	FAD	C4A-C5A-N7A	-2.91	107.25	110.58
2	A	1001	FAD	O3'-C3'-C4'	-2.84	102.47	108.93
2	B	1001	FAD	N9A-C8A-N7A	-2.76	110.03	113.94
2	A	1001	FAD	O5B-C5B-C4B	-2.70	99.82	108.99
2	B	1001	FAD	C5A-N7A-C8A	2.69	107.67	103.45
2	B	1001	FAD	C4X-C4-N3	2.67	120.05	113.25
2	B	1001	FAD	C5X-C9A-N10	2.66	120.37	117.97
2	B	1001	FAD	O4-C4-C4X	-2.63	119.58	126.53
2	B	1001	FAD	C4B-O4B-C1B	-2.59	103.75	109.47
2	A	1001	FAD	C4X-C4-N3	2.57	119.80	113.25
2	A	1001	FAD	O4B-C4B-C5B	2.57	117.57	109.33
2	A	1001	FAD	C4A-N9A-C8A	2.53	108.40	105.74
2	A	1001	FAD	O4B-C1B-N9A	2.50	112.90	108.09
2	A	1001	FAD	C4-N3-C2	-2.50	121.21	125.64
2	A	1001	FAD	C4-C4X-N5	2.46	121.61	118.21
2	A	1001	FAD	C4X-C10-N1	-2.40	118.70	124.59
2	A	1001	FAD	C10-N1-C2	2.34	121.91	116.85
2	B	1001	FAD	C4X-C10-N10	2.33	119.81	116.48
2	A	1001	FAD	C10-C4X-N5	-2.31	120.09	124.81
2	B	1001	FAD	C9A-C5X-N5	-2.26	120.05	122.45
2	B	1001	FAD	C4-C4X-C10	2.24	120.78	116.93
2	B	1001	FAD	C4X-C10-N1	-2.11	119.41	124.59
2	B	1001	FAD	C10-C4X-N5	-2.09	120.55	124.81
2	B	1001	FAD	C2'-C1'-N10	2.04	119.82	110.20

There are no chirality outliers.

All (1) torsion outliers are listed below:

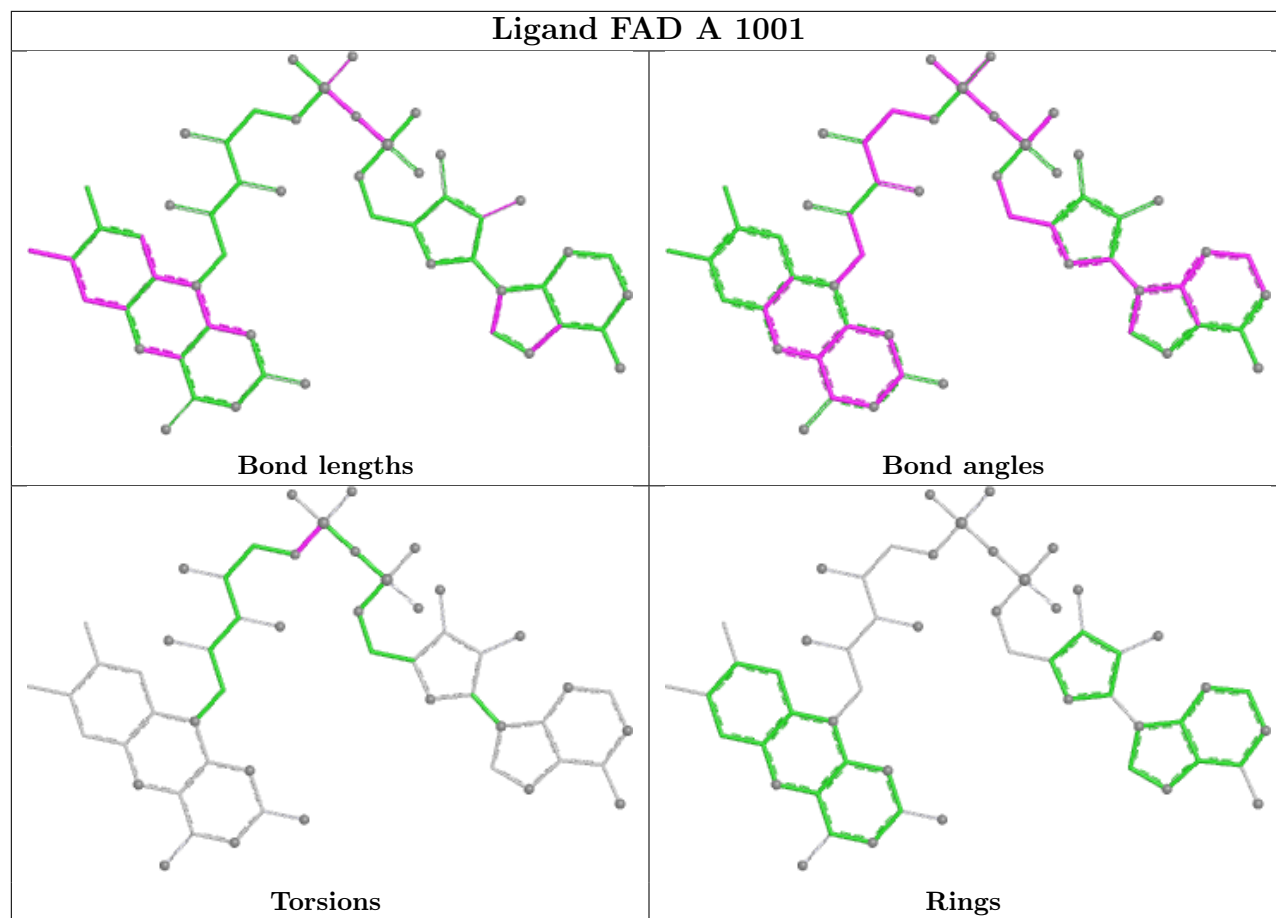
Mol	Chain	Res	Type	Atoms
2	A	1001	FAD	C5'-O5'-P-O1P

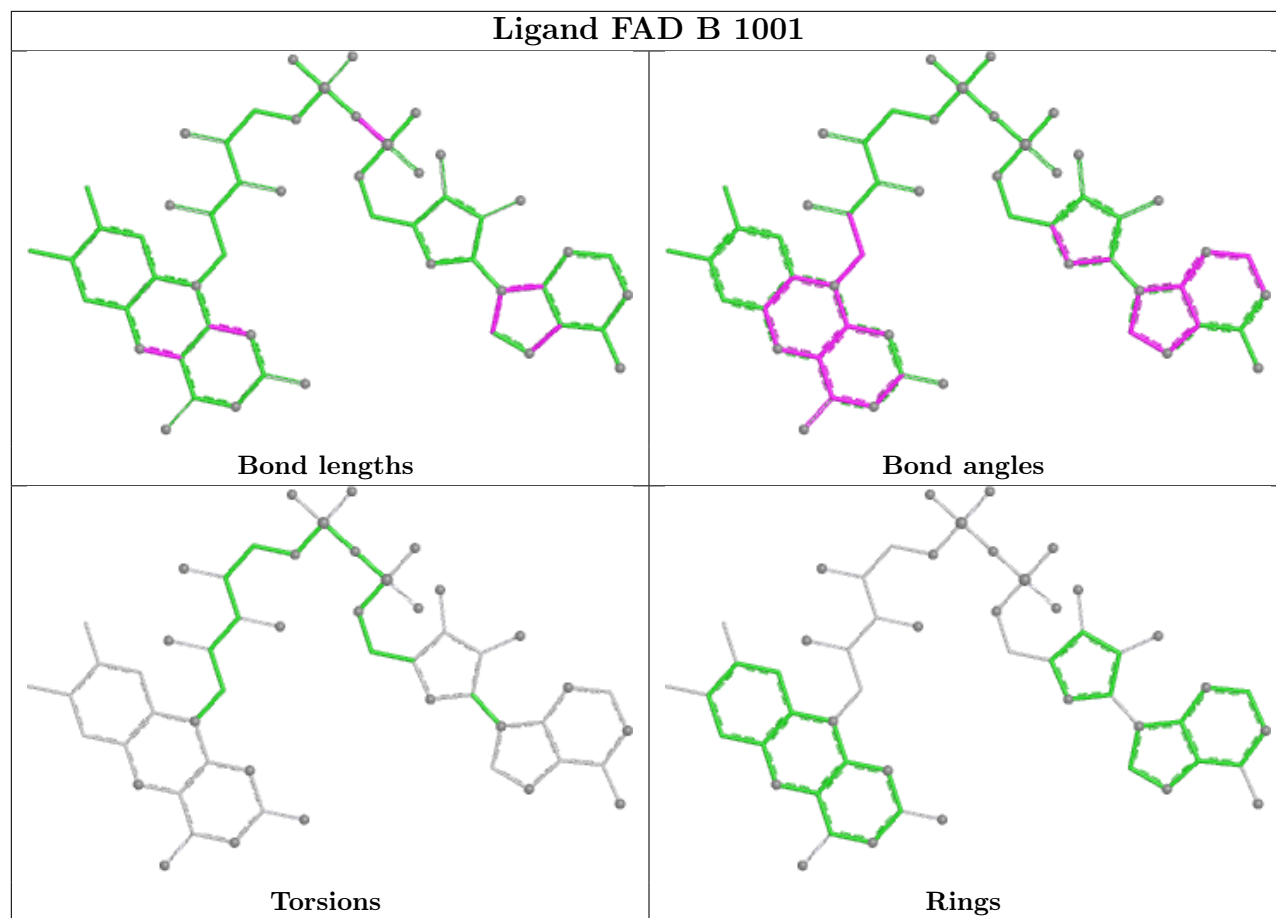
There are no ring outliers.

2 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1001	FAD	4	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 [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	557/607 (91%)	-0.39	2 (0%) 88 86	17, 45, 72, 103	0
1	B	530/607 (87%)	0.35	14 (2%) 57 53	39, 78, 120, 157	1 (0%)
All	All	1087/1214 (89%)	-0.03	16 (1%) 72 68	17, 59, 109, 157	1 (0%)

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	113	THR	5.3
1	B	38	ALA	4.4
1	B	519	ALA	3.4
1	B	496	ASP	3.2
1	B	41	GLY	2.6
1	B	359	GLY	2.6
1	A	154	PHE	2.4
1	B	109	ASP	2.3
1	B	360	GLY	2.2
1	B	149	LEU	2.2
1	B	538	GLU	2.2
1	B	168	ASP	2.2
1	B	468	GLY	2.1
1	B	217	THR	2.1
1	B	352	ASN	2.0
1	A	247	PRO	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](#)

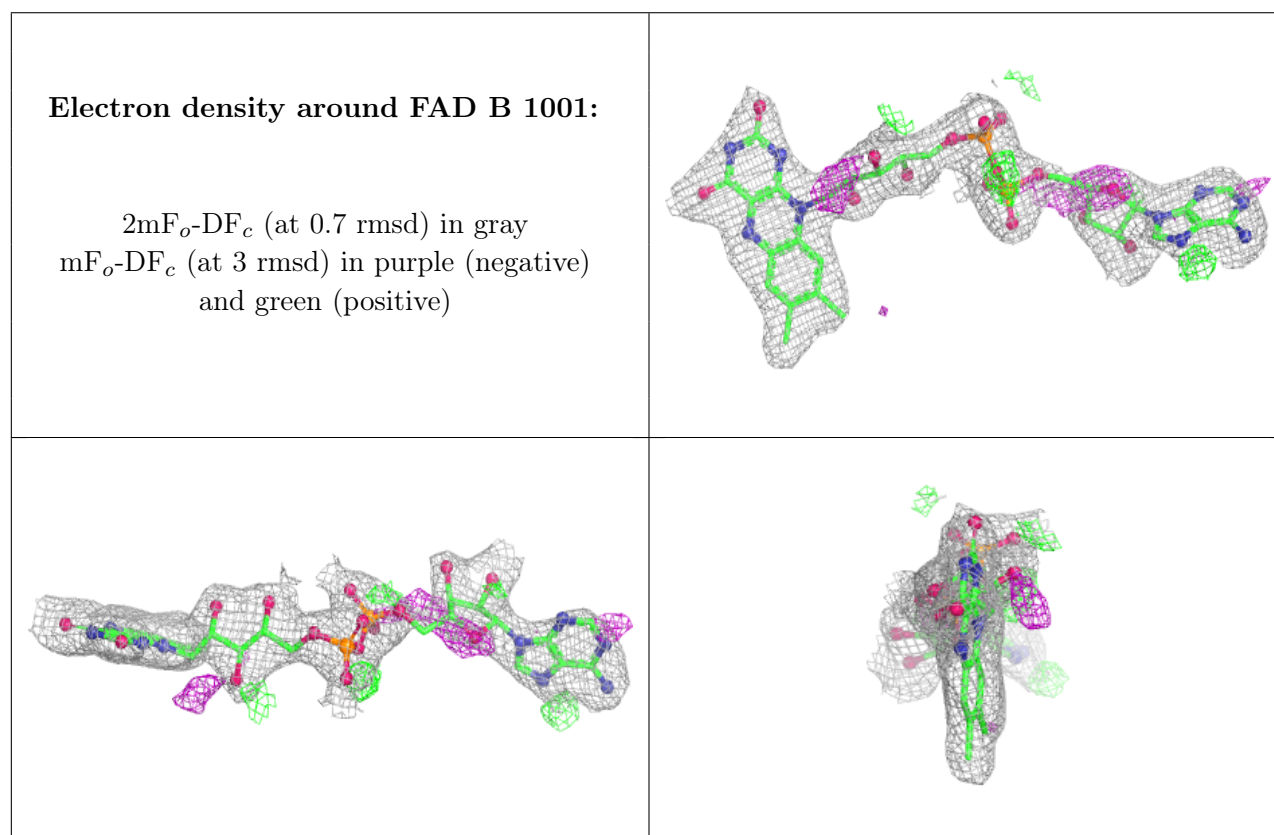
There are no oligosaccharides in this entry.

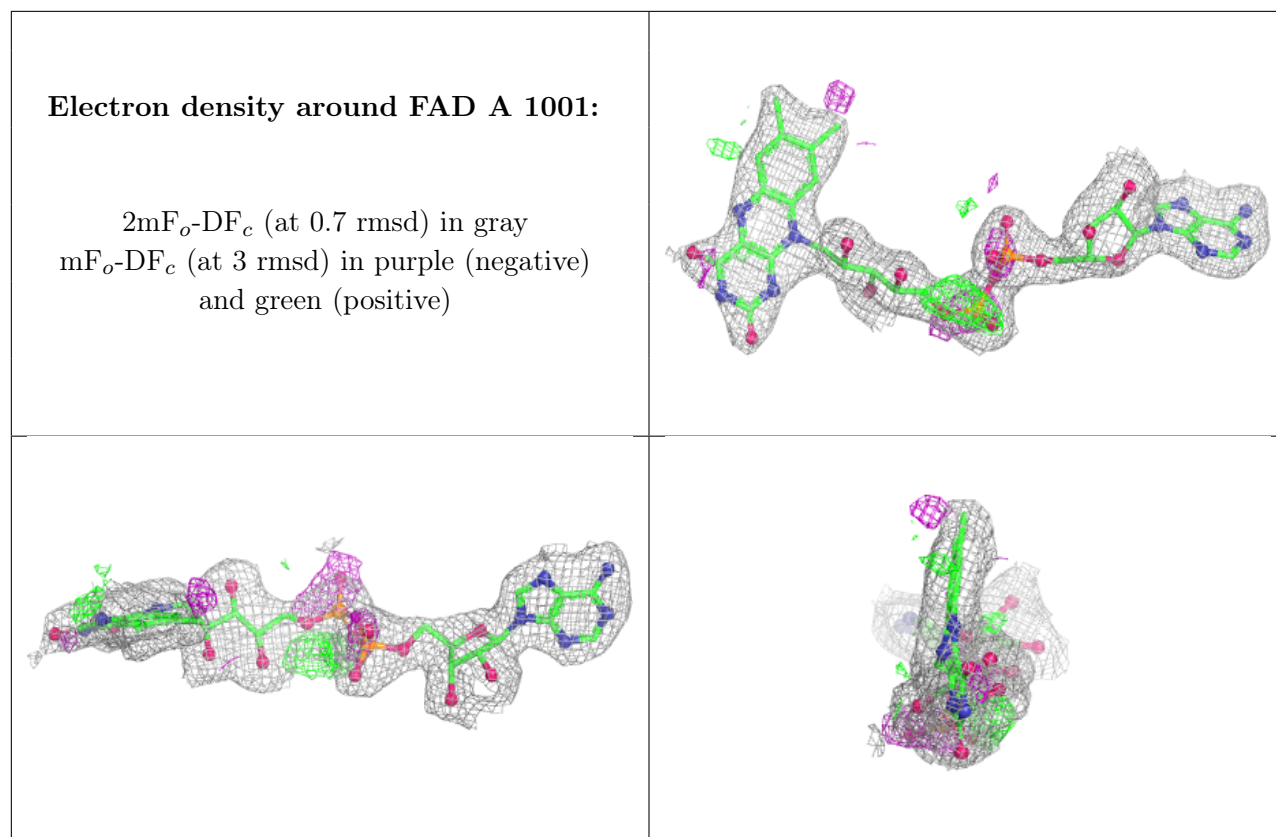
6.4 Ligands [i](#)

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($\text{\AA}^2$)	Q<0.9
2	FAD	B	1001	53/53	0.92	0.09	29,44,50,50	0
2	FAD	A	1001	53/53	0.96	0.06	12,22,29,33	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.





6.5 Other polymers [i](#)

There are no such residues in this entry.