



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 02:05 PM UTC

PDB ID : 4Z2U / pdb_00004z2u
Title : Crystal Structure of 2-hydroxybiphenyl 3-monooxygenase R242Q from *Pseudomonas azelaica*
Authors : Kanteev, M.; Bregman-Cohen, A.; Fishman, A.
Deposited on : 2015-03-30
Resolution : 2.50 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

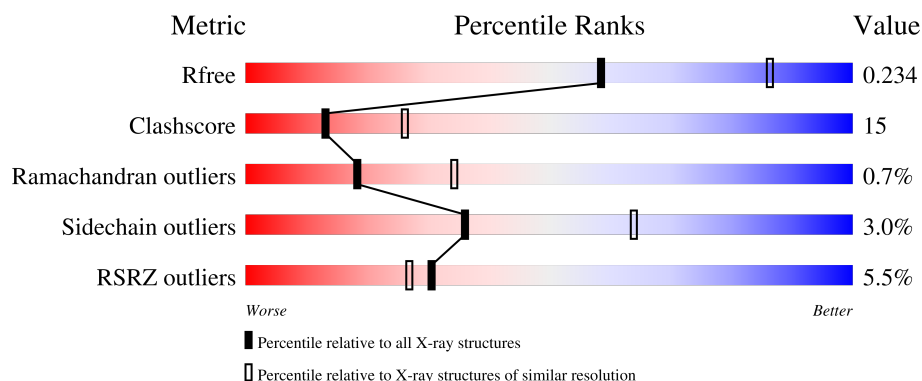
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	592	6% 69% 24% • 5%
1	B	592	5% 73% 19% • 5%

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
2	FAD	A	601	X	-	-	-
2	FAD	B	601	X	-	-	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9413 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 2-hydroxybiphenyl-3-monooxygenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	563	Total	C	N	O	S	0	0	0
			4330	2737	759	814	20			
1	B	563	Total	C	N	O	S	0	0	0
			4330	2737	759	814	20			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	MET	-	initiating methionine	UNP O06647
A	-4	HIS	-	expression tag	UNP O06647
A	-3	HIS	-	expression tag	UNP O06647
A	-2	HIS	-	expression tag	UNP O06647
A	-1	HIS	-	expression tag	UNP O06647
A	0	HIS	-	expression tag	UNP O06647
A	1	HIS	-	expression tag	UNP O06647
A	242	GLN	ARG	engineered mutation	UNP O06647
B	-5	MET	-	initiating methionine	UNP O06647
B	-4	HIS	-	expression tag	UNP O06647
B	-3	HIS	-	expression tag	UNP O06647
B	-2	HIS	-	expression tag	UNP O06647
B	-1	HIS	-	expression tag	UNP O06647
B	0	HIS	-	expression tag	UNP O06647
B	1	HIS	-	expression tag	UNP O06647
B	242	GLN	ARG	engineered mutation	UNP O06647

- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (CCD ID: FAD) (formula: C₂₇H₃₃N₉O₁₅P₂).



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

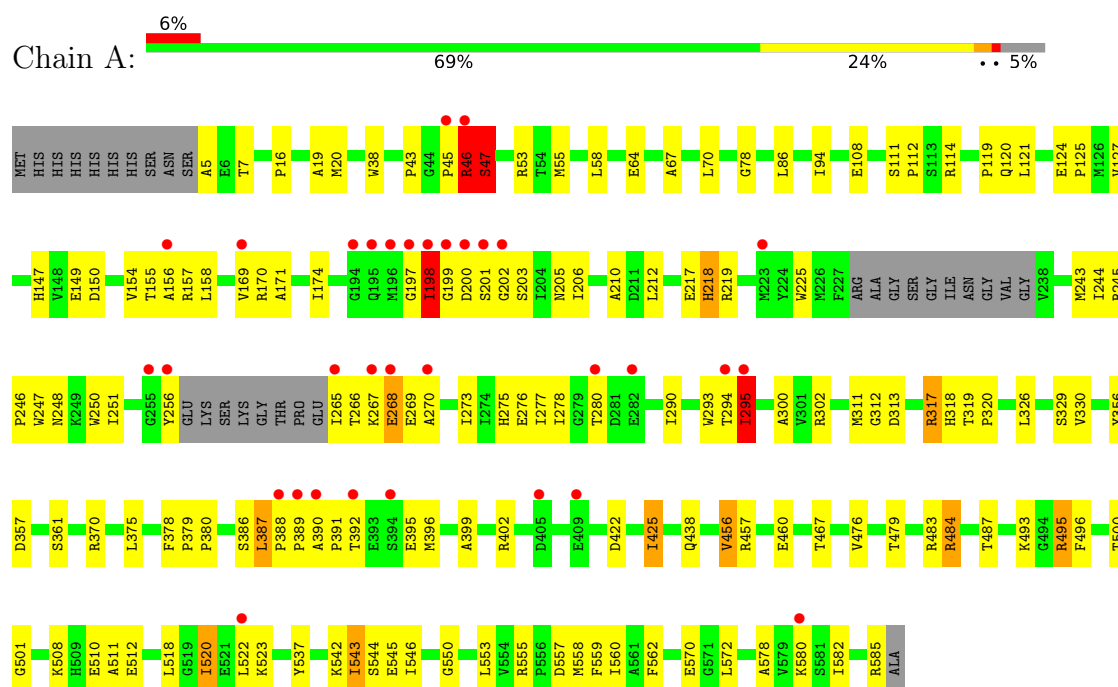
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	313	Total O 313 313	0	0
3	B	334	Total O 334 334	0	0

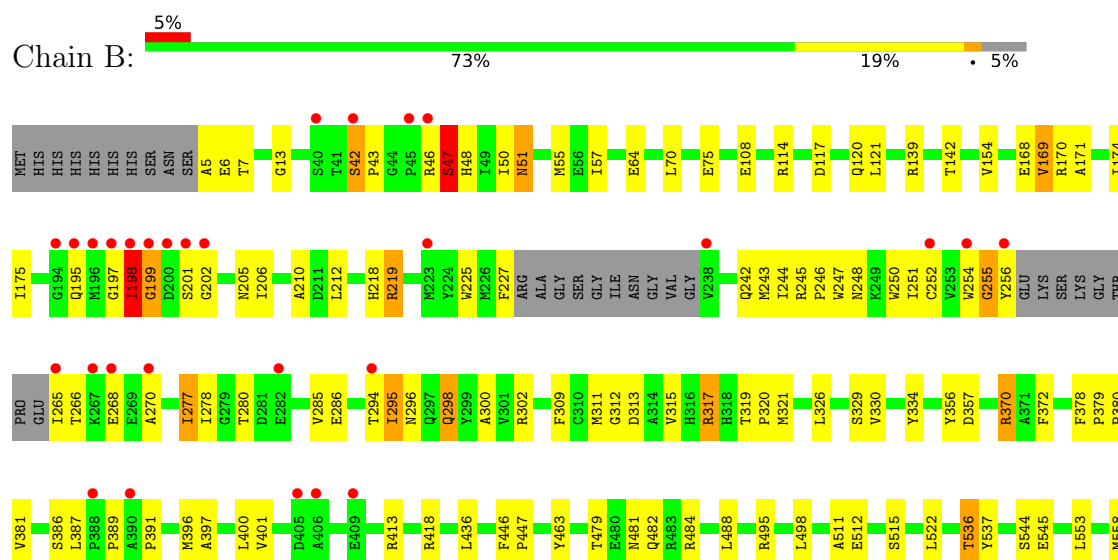
3 Residue-property plots [i](#)

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: 2-hydroxybiphenyl-3-monooxygenase



• Molecule 1: 2-hydroxybiphenyl-3-monooxygenase





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	154.14Å 131.54Å 79.00Å 90.00° 98.28° 90.00°	Depositor
Resolution (Å)	32.37 – 2.50 32.37 – 2.50	Depositor EDS
% Data completeness (in resolution range)	92.2 (32.37-2.50) 92.2 (32.37-2.50)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.67 (at 2.51Å)	Xtriage
Refinement program	PHENIX 1.7.3 _928	Depositor
R, R_{free}	0.204 , 0.237 0.203 , 0.234	Depositor DCC
R_{free} test set	2528 reflections (4.69%)	wwPDB-VP
Wilson B-factor (Å ²)	26.3	Xtriage
Anisotropy	0.500	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 40.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	9413	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.06% 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	0.80	7/4425 (0.2%)	1.01	19/5998 (0.3%)
1	B	0.69	5/4425 (0.1%)	0.89	9/5998 (0.2%)
All	All	0.75	12/8850 (0.1%)	0.95	28/11996 (0.2%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	5
1	B	0	1
All	All	0	6

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	149	GLU	C-N	-22.88	1.07	1.33
1	B	255	GLY	C-N	-22.33	1.02	1.33
1	A	46	ARG	C-N	-19.43	1.06	1.33
1	A	218	HIS	C-N	-12.56	1.17	1.33
1	B	169	VAL	C-N	-11.77	1.17	1.33
1	B	51	ASN	C-N	-11.07	1.17	1.33
1	A	217	GLU	C-N	-10.43	1.18	1.33
1	A	543	ILE	C-N	8.19	1.45	1.34
1	A	47	SER	C-N	-6.30	1.24	1.33
1	B	50	ILE	C-N	5.59	1.42	1.33
1	B	168	GLU	C-N	-5.57	1.26	1.33
1	A	150	ASP	C-N	5.50	1.42	1.33

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	46	ARG	O-C-N	-19.27	96.96	122.59
1	A	150	ASP	O-C-N	-15.82	102.32	122.73
1	A	46	ARG	CA-C-N	15.71	151.55	121.54
1	A	46	ARG	C-N-CA	15.71	151.55	121.54
1	A	295	ILE	N-CA-CB	-14.03	86.61	112.43
1	A	218	HIS	O-C-N	-12.64	104.41	122.36
1	B	169	VAL	O-C-N	-12.43	109.78	123.20
1	A	543	ILE	O-C-N	10.04	133.21	122.17
1	B	255	GLY	CA-C-N	9.33	138.49	121.70
1	B	255	GLY	C-N-CA	9.33	138.49	121.70
1	A	318	HIS	N-CA-CB	-8.06	96.36	110.87
1	A	47	SER	O-C-N	-7.06	113.19	122.59
1	A	276	GLU	N-CA-C	-6.53	105.44	113.41
1	A	318	HIS	N-CA-C	6.50	116.85	108.34
1	B	199	GLY	N-CA-C	-6.41	103.88	112.14
1	B	42	SER	N-CA-C	6.05	123.18	109.81
1	A	150	ASP	CA-C-N	6.00	131.64	121.14
1	A	150	ASP	C-N-CA	6.00	131.64	121.14
1	B	255	GLY	O-C-N	-5.69	117.77	122.77
1	A	218	HIS	CA-C-N	5.45	130.72	123.20
1	A	218	HIS	C-N-CA	5.45	130.72	123.20
1	B	168	GLU	CA-C-N	5.39	130.49	122.99
1	B	168	GLU	C-N-CA	5.39	130.49	122.99
1	A	38	TRP	N-CA-C	5.14	118.33	110.36
1	A	542	LYS	CA-C-N	5.10	127.92	121.85
1	A	542	LYS	C-N-CA	5.10	127.92	121.85
1	B	50	ILE	O-C-N	-5.05	116.30	122.97
1	A	78	GLY	N-CA-C	5.04	120.13	113.37

There are no chirality outliers.

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	218	HIS	Mainchain
1	A	46	ARG	Peptide,Mainchain
1	A	47	SER	Mainchain
1	A	495	ARG	Sidechain
1	B	255	GLY	Mainchain

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	4330	0	4275	148	0
1	B	4330	0	4276	110	0
2	A	53	0	30	8	0
2	B	53	0	30	8	0
3	A	313	0	0	3	4
3	B	334	0	0	1	5
All	All	9413	0	8611	259	6

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 15.

All (259) 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:201:SER:HB3	1:A:256:TYR:O	1.39	1.22
1:A:319:THR:CG2	1:A:320:PRO:HD2	1.71	1.21
1:A:319:THR:HG23	1:A:320:PRO:HD2	1.17	1.10
1:A:511:ALA:HB1	1:A:522:LEU:HD11	1.10	1.07
1:B:75:GLU:OE1	1:B:75:GLU:N	1.95	1.00
1:A:511:ALA:HB1	1:A:522:LEU:CD1	1.94	0.96
1:A:5:ALA:HB3	1:A:169:VAL:HG22	1.47	0.95
1:B:294:THR:O	1:B:295:ILE:HG22	1.69	0.92
1:A:511:ALA:CB	1:A:522:LEU:HD11	1.98	0.92
1:A:319:THR:CG2	1:A:320:PRO:CD	2.49	0.91
1:B:46:ARG:O	1:B:47:SER:HB3	1.70	0.90
1:B:108:GLU:HG3	1:B:114:ARG:NH2	1.86	0.90
1:B:46:ARG:HD3	2:B:601:FAD:C7	2.02	0.90
1:A:46:ARG:HD3	2:A:601:FAD:C8	2.04	0.87
1:B:370:ARG:O	1:B:370:ARG:HD3	1.74	0.87
1:A:206:ILE:HD13	1:A:270:ALA:HB1	1.57	0.86
1:B:511:ALA:HB1	1:B:522:LEU:HD11	1.58	0.85
1:A:370:ARG:O	1:A:370:ARG:HD3	1.75	0.85
1:B:205:ASN:HB3	1:B:251:ILE:HD11	1.59	0.84
1:A:154:VAL:HG21	1:A:174:ILE:HG13	1.60	0.84
1:A:202:GLY:O	1:A:256:TYR:HB3	1.78	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:511:ALA:HB1	1:B:522:LEU:CD1	2.08	0.82
1:A:225:TRP:HZ3	1:A:378:PHE:CE1	1.98	0.81
1:B:202:GLY:O	1:B:256:TYR:C	2.24	0.81
1:A:388:PRO:O	1:A:390:ALA:N	2.14	0.80
1:A:370:ARG:HD3	1:A:370:ARG:C	2.06	0.79
1:B:46:ARG:HD3	2:B:601:FAD:C8	2.13	0.78
1:B:370:ARG:HD3	1:B:370:ARG:C	2.08	0.78
1:A:46:ARG:HD3	2:A:601:FAD:C7	2.14	0.78
1:A:319:THR:HG22	1:A:320:PRO:CD	2.14	0.77
1:B:70:LEU:O	1:B:245:ARG:NH2	2.17	0.77
1:A:197:GLY:O	1:A:198:ILE:HB	1.85	0.76
1:A:154:VAL:O	1:A:170:ARG:HD2	1.86	0.76
1:A:379:PRO:HB2	1:A:380:PRO:HD3	1.64	0.76
1:A:70:LEU:O	1:A:245:ARG:NH2	2.19	0.73
1:A:94:ILE:HB	1:A:425:ILE:HD11	1.71	0.73
1:B:154:VAL:HG21	1:B:174:ILE:HG13	1.69	0.72
1:B:75:GLU:H	1:B:75:GLU:CD	1.95	0.72
1:B:225:TRP:CZ3	1:B:378:PHE:CE1	2.77	0.72
1:A:201:SER:CB	1:A:256:TYR:O	2.29	0.71
1:B:244:ILE:O	1:B:245:ARG:HD3	1.90	0.71
1:A:5:ALA:CB	1:A:169:VAL:HG22	2.20	0.71
1:A:546:ILE:HD12	1:A:550:GLY:HA3	1.72	0.71
1:A:244:ILE:O	1:A:245:ARG:HD3	1.90	0.71
1:A:270:ALA:HA	1:A:273:ILE:HD12	1.73	0.70
1:A:267:LYS:HG3	1:A:268:GLU:H	1.56	0.70
1:B:295:ILE:O	1:B:295:ILE:HG23	1.92	0.70
1:B:55:MET:HE3	1:B:64:GLU:HG3	1.74	0.69
1:A:225:TRP:CZ3	1:A:378:PHE:CE1	2.79	0.69
1:A:294:THR:HG23	1:A:294:THR:O	1.91	0.68
1:A:205:ASN:HB3	1:A:251:ILE:HD11	1.76	0.68
1:B:498:LEU:C	1:B:498:LEU:HD23	2.19	0.67
1:A:319:THR:HG22	1:A:320:PRO:HD2	1.68	0.67
1:B:48:HIS:CD2	1:B:242:GLN:NE2	2.63	0.67
1:B:387:LEU:HD13	1:B:391:PRO:HG3	1.77	0.67
1:A:108:GLU:HG3	1:A:114:ARG:NH2	2.09	0.67
1:A:278:ILE:HG22	1:A:280:THR:HG22	1.77	0.66
1:A:388:PRO:C	1:A:390:ALA:H	2.02	0.66
1:A:555:ARG:HG3	1:A:559:PHE:O	1.95	0.66
1:A:387:LEU:HG	1:A:391:PRO:HB3	1.78	0.66
1:B:43:PRO:O	1:B:121:LEU:HD13	1.96	0.66
1:B:46:ARG:O	1:B:120:GLN:OE1	2.14	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:569:ARG:O	1:B:573:GLU:HG2	1.97	0.65
1:A:94:ILE:HB	1:A:425:ILE:CD1	2.28	0.64
1:A:147:HIS:HB3	1:A:156:ALA:HB2	1.79	0.63
1:B:397:ALA:O	1:B:401:VAL:HG23	1.99	0.63
1:B:379:PRO:HB2	1:B:380:PRO:HD3	1.81	0.62
1:B:46:ARG:O	1:B:47:SER:CB	2.45	0.62
1:A:267:LYS:C	1:A:269:GLU:H	2.08	0.62
1:B:225:TRP:HZ3	1:B:378:PHE:CE1	2.16	0.62
1:A:312:GLY:HA2	1:A:329:SER:OG	2.00	0.62
1:A:495:ARG:NH2	1:A:523:LYS:HD3	2.15	0.61
1:A:154:VAL:HG21	1:A:174:ILE:CG1	2.29	0.61
1:A:422:ASP:O	1:A:425:ILE:HG12	2.00	0.61
1:A:370:ARG:HH21	1:A:460:GLU:CD	2.08	0.60
1:B:326:LEU:HD23	1:B:326:LEU:C	2.26	0.60
1:B:202:GLY:O	1:B:256:TYR:HB3	2.02	0.59
1:A:46:ARG:O	2:A:601:FAD:C4	2.50	0.59
1:A:391:PRO:HG3	1:A:396:MET:CE	2.33	0.59
1:B:202:GLY:O	1:B:256:TYR:CA	2.49	0.59
1:A:46:ARG:O	2:A:601:FAD:C4X	2.50	0.59
1:A:94:ILE:HD12	1:A:425:ILE:HD13	1.83	0.59
1:A:243:MET:HA	1:A:250:TRP:CE3	2.38	0.59
1:A:7:THR:O	1:A:171:ALA:HA	2.03	0.59
1:A:518:LEU:HB2	1:A:520:ILE:HG22	1.84	0.59
1:A:319:THR:HG22	1:A:320:PRO:N	2.16	0.58
1:B:225:TRP:HH2	1:B:321:MET:HE1	1.69	0.58
1:A:147:HIS:CB	1:A:156:ALA:HB2	2.34	0.58
1:B:225:TRP:CZ3	1:B:378:PHE:CZ	2.92	0.58
1:A:47:SER:HB3	1:A:120:GLN:OE1	2.04	0.57
1:A:147:HIS:HA	1:A:156:ALA:HA	1.87	0.57
1:B:46:ARG:CD	2:B:601:FAD:C8	2.82	0.57
1:A:557:ASP:O	1:A:558:MET:HB2	2.04	0.57
1:A:198:ILE:O	1:A:199:GLY:C	2.46	0.56
1:A:510:GLU:HG3	1:A:572:LEU:HD22	1.88	0.56
1:B:197:GLY:C	1:B:198:ILE:CG1	2.79	0.56
1:B:522:LEU:HD12	1:B:522:LEU:O	2.06	0.56
1:A:295:ILE:HG23	1:A:295:ILE:O	2.05	0.56
1:A:495:ARG:HH21	1:A:523:LYS:HD3	1.70	0.56
1:A:280:THR:HG23	1:A:280:THR:O	2.04	0.56
1:A:388:PRO:C	1:A:390:ALA:N	2.60	0.55
1:B:46:ARG:HD3	2:B:601:FAD:C7M	2.36	0.55
1:A:295:ILE:O	1:A:295:ILE:CG2	2.54	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:295:ILE:O	1:B:295:ILE:CG2	2.55	0.55
1:A:43:PRO:O	1:A:121:LEU:HD13	2.06	0.55
1:A:544:SER:O	1:A:545:GLU:HB2	2.07	0.55
1:A:267:LYS:HG3	1:A:268:GLU:N	2.21	0.55
1:A:157:ARG:O	1:A:158:LEU:HD23	2.06	0.55
1:A:580:LYS:HG2	1:A:585:ARG:O	2.06	0.55
1:A:202:GLY:O	1:A:256:TYR:CB	2.53	0.54
1:B:206:ILE:HD13	1:B:270:ALA:HB1	1.89	0.54
1:A:294:THR:O	1:A:294:THR:CG2	2.56	0.54
1:B:48:HIS:HB2	1:B:117:ASP:OD2	2.08	0.54
1:B:198:ILE:CG1	1:B:372:PHE:HE1	2.21	0.54
1:A:169:VAL:HG12	1:A:169:VAL:O	2.07	0.53
1:A:266:THR:CG2	1:A:290:ILE:HD13	2.39	0.53
1:A:311:MET:HB3	1:A:356:TYR:OH	2.08	0.53
1:A:391:PRO:CG	1:A:396:MET:HE2	2.39	0.53
1:B:296:ASN:O	1:B:298:GLN:HG3	2.08	0.53
1:A:45:PRO:O	1:A:120:GLN:HB2	2.08	0.53
1:A:200:ASP:CG	1:A:200:ASP:O	2.52	0.52
1:A:313:ASP:O	1:A:317:ARG:HG3	2.09	0.52
1:B:57:ILE:HG12	1:B:334:TYR:CD2	2.44	0.52
1:A:266:THR:HG23	1:A:290:ILE:HD13	1.92	0.52
1:B:212:LEU:HD21	1:B:278:ILE:HD13	1.90	0.52
1:A:570:GLU:CD	1:A:570:GLU:H	2.17	0.52
1:B:201:SER:O	1:B:294:THR:O	2.27	0.52
1:B:312:GLY:HA2	1:B:329:SER:OG	2.10	0.52
1:B:511:ALA:CB	1:B:522:LEU:HD11	2.37	0.52
1:B:294:THR:HG23	1:B:295:ILE:N	2.25	0.52
1:A:86:LEU:CD2	1:A:212:LEU:HD13	2.40	0.52
1:A:155:THR:HG22	1:A:155:THR:O	2.10	0.52
1:A:248:ASN:OD1	1:A:248:ASN:N	2.42	0.52
1:A:484:ARG:HD2	1:A:537:TYR:OH	2.10	0.51
1:A:387:LEU:HD21	1:A:396:MET:SD	2.50	0.51
1:B:47:SER:HB2	2:B:601:FAD:C2	2.40	0.51
1:B:5:ALA:HB3	1:B:169:VAL:HG22	1.91	0.51
1:A:543:ILE:O	3:A:703:HOH:O	2.19	0.51
1:B:294:THR:O	1:B:295:ILE:CG2	2.52	0.51
1:A:391:PRO:HG3	1:A:396:MET:HE1	1.92	0.51
1:B:319:THR:HB	1:B:320:PRO:HD2	1.92	0.51
1:B:391:PRO:HG2	1:B:396:MET:HE2	1.93	0.51
1:A:225:TRP:HZ3	1:A:378:PHE:CZ	2.29	0.50
2:A:601:FAD:O2A	3:A:704:HOH:O	2.19	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:496:PHE:O	1:A:522:LEU:HA	2.11	0.50
1:B:436:LEU:HB3	1:B:463:TYR:CD1	2.46	0.50
1:A:200:ASP:O	1:A:200:ASP:OD1	2.30	0.50
1:A:202:GLY:HA2	1:A:294:THR:HA	1.94	0.50
1:B:311:MET:HB3	1:B:356:TYR:OH	2.11	0.50
1:B:544:SER:O	1:B:545:GLU:HB2	2.12	0.49
1:A:265:ILE:HG22	1:A:266:THR:N	2.27	0.49
1:A:312:GLY:HA3	2:A:601:FAD:O1P	2.13	0.49
1:B:210:ALA:O	1:B:248:ASN:HA	2.11	0.49
1:A:225:TRP:CZ3	1:A:378:PHE:CZ	3.00	0.49
1:A:493:LYS:HD2	1:A:495:ARG:CZ	2.43	0.49
1:A:267:LYS:HB3	1:A:269:GLU:OE1	2.13	0.49
1:B:195:GLN:NE2	1:B:298:GLN:NE2	2.61	0.49
1:B:227:PHE:CE2	1:B:381:VAL:HG11	2.48	0.49
1:B:536:THR:HG22	1:B:537:TYR:CD2	2.48	0.49
1:B:198:ILE:HD11	1:B:372:PHE:HE1	1.78	0.49
1:B:202:GLY:O	1:B:256:TYR:CB	2.61	0.49
1:A:479:THR:HA	1:A:483:ARG:O	2.13	0.48
1:A:210:ALA:O	1:A:248:ASN:HA	2.14	0.48
1:A:319:THR:HG21	1:A:375:LEU:HD21	1.94	0.48
1:B:48:HIS:CD2	1:B:242:GLN:HE22	2.31	0.48
1:A:45:PRO:CB	1:A:119:PRO:HB3	2.44	0.48
1:B:285:VAL:HG12	1:B:286:GLU:N	2.28	0.48
1:B:218:HIS:NE2	1:B:219:ARG:HD2	2.28	0.48
1:A:246:PRO:HA	1:A:247:TRP:HA	1.56	0.47
1:A:379:PRO:CB	1:A:380:PRO:HD3	2.40	0.47
1:A:203:SER:N	1:A:293:TRP:O	2.46	0.47
1:A:19:ALA:HB1	1:A:127:VAL:HG12	1.97	0.47
1:B:48:HIS:HB2	1:B:117:ASP:CG	2.40	0.47
1:A:391:PRO:HG3	1:A:396:MET:HE2	1.96	0.47
1:B:302:ARG:HA	1:B:357:ASP:OD1	2.15	0.47
1:B:285:VAL:CG1	1:B:286:GLU:N	2.77	0.47
1:A:267:LYS:C	1:A:269:GLU:N	2.71	0.46
1:B:277:ILE:HG22	1:B:278:ILE:N	2.31	0.46
1:B:386:SER:OG	1:B:413:ARG:NH1	2.42	0.46
1:A:55:MET:HE3	1:A:64:GLU:HG3	1.96	0.46
1:A:370:ARG:C	1:A:370:ARG:CD	2.80	0.46
1:A:268:GLU:CG	1:A:268:GLU:O	2.64	0.46
1:A:326:LEU:C	1:A:326:LEU:HD23	2.40	0.46
1:A:399:ALA:HA	1:A:402:ARG:NH2	2.32	0.46
1:B:227:PHE:CZ	1:B:381:VAL:HG11	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:246:PRO:HA	1:B:247:TRP:HA	1.60	0.45
1:A:476:VAL:HG23	1:A:487:THR:HG23	1.99	0.45
1:B:278:ILE:HG22	1:B:280:THR:HG22	1.99	0.45
1:A:243:MET:HA	1:A:250:TRP:CZ3	2.52	0.45
1:A:267:LYS:CB	1:A:269:GLU:OE1	2.65	0.45
1:A:147:HIS:HB3	1:A:156:ALA:CB	2.47	0.45
1:A:124:GLU:N	1:A:125:PRO:HD2	2.32	0.44
1:A:295:ILE:HD13	3:A:924:HOH:O	2.17	0.44
1:B:198:ILE:HG22	1:B:199:GLY:H	1.82	0.44
1:B:387:LEU:HD13	1:B:391:PRO:CG	2.46	0.44
1:A:467:THR:HA	1:A:560:ILE:HD12	1.99	0.44
1:B:488:LEU:O	1:B:558:MET:HE1	2.16	0.44
1:A:58:LEU:HD13	1:A:67:ALA:HB2	1.98	0.44
1:B:312:GLY:HA3	2:B:601:FAD:O1P	2.17	0.44
1:B:244:ILE:HD11	1:B:251:ILE:CG2	2.47	0.44
1:A:300:ALA:HB1	1:A:302:ARG:O	2.18	0.43
1:A:379:PRO:HB2	1:A:380:PRO:CD	2.41	0.43
1:A:425:ILE:HG12	1:A:425:ILE:H	1.68	0.43
1:B:265:ILE:HG13	1:B:266:THR:N	2.34	0.43
1:B:391:PRO:HG2	1:B:396:MET:CE	2.48	0.43
1:B:446:PHE:HA	1:B:447:PRO:HD3	1.91	0.43
1:B:243:MET:HA	1:B:250:TRP:CE3	2.53	0.43
1:B:6:GLU:HG3	1:B:170:ARG:HB3	1.99	0.43
1:B:197:GLY:C	1:B:198:ILE:HG12	2.42	0.43
1:A:357:ASP:O	1:A:361:SER:HB2	2.19	0.43
1:A:202:GLY:O	1:A:256:TYR:N	2.52	0.43
1:A:422:ASP:O	1:A:425:ILE:CG1	2.67	0.43
1:B:46:ARG:CD	2:B:601:FAD:C7	2.87	0.43
1:B:511:ALA:HB1	1:B:522:LEU:HD12	1.98	0.43
1:A:280:THR:O	1:A:280:THR:CG2	2.67	0.43
1:B:5:ALA:O	1:B:169:VAL:HA	2.18	0.43
1:B:13:GLY:HA2	2:B:601:FAD:O4B	2.19	0.43
1:A:295:ILE:H	1:A:295:ILE:HG22	1.47	0.42
1:A:553:LEU:HB3	1:A:562:PHE:HB3	2.00	0.42
1:B:294:THR:CG2	1:B:295:ILE:N	2.82	0.42
1:A:46:ARG:O	2:A:601:FAD:N5	2.52	0.42
1:A:378:PHE:O	1:A:379:PRO:C	2.61	0.42
1:B:479:THR:HG23	1:B:536:THR:HB	2.01	0.42
1:A:391:PRO:CG	1:A:396:MET:CE	2.97	0.42
1:A:392:THR:H	1:A:395:GLU:HB2	1.84	0.42
1:B:379:PRO:CB	1:B:380:PRO:HD3	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:243:MET:CA	1:A:250:TRP:CZ3	3.02	0.42
1:A:16:PRO:HD2	2:A:601:FAD:O2P	2.20	0.42
1:A:20:MET:HG2	1:A:330:VAL:HG13	2.00	0.42
1:B:7:THR:O	1:B:171:ALA:HA	2.20	0.42
1:B:481:ASN:O	1:B:482:GLN:HB2	2.19	0.42
1:B:379:PRO:N	1:B:380:PRO:CD	2.83	0.42
1:A:111:SER:OG	1:A:112:PRO:HD2	2.19	0.41
1:A:510:GLU:CG	1:A:572:LEU:HD22	2.50	0.41
1:B:197:GLY:O	1:B:198:ILE:CG1	2.68	0.41
1:B:51:ASN:HB2	3:B:823:HOH:O	2.20	0.41
1:B:300:ALA:HB3	1:B:315:VAL:HB	2.01	0.41
1:B:197:GLY:O	1:B:198:ILE:HG13	2.21	0.41
1:B:553:LEU:HB3	1:B:562:PHE:HB3	2.02	0.41
1:B:570:GLU:H	1:B:570:GLU:CD	2.25	0.41
1:B:418:ARG:O	1:B:418:ARG:HD3	2.21	0.41
1:A:198:ILE:CG2	1:A:199:GLY:N	2.82	0.41
1:B:198:ILE:CD1	1:B:372:PHE:HE1	2.34	0.41
1:A:202:GLY:O	1:A:256:TYR:CA	2.69	0.41
1:A:268:GLU:O	1:A:268:GLU:CD	2.64	0.41
1:B:175:ILE:HA	1:B:309:PHE:O	2.20	0.41
1:B:205:ASN:HA	1:B:252:CYS:O	2.21	0.41
1:B:498:LEU:C	1:B:498:LEU:CD2	2.91	0.41
1:B:206:ILE:HD11	1:B:254:TRP:HH2	1.86	0.40
1:B:313:ASP:O	1:B:317:ARG:HG3	2.21	0.40
1:B:391:PRO:CG	1:B:396:MET:HE2	2.51	0.40
1:A:53:ARG:NH1	1:A:438:GLN:OE1	2.53	0.40
1:A:86:LEU:HD22	1:A:212:LEU:HD13	2.03	0.40
1:A:456:VAL:HG22	1:A:457:ARG:HG3	2.03	0.40
1:B:139:ARG:HG3	1:B:142:THR:OG1	2.21	0.40
1:A:243:MET:HE3	1:A:243:MET:HB3	1.98	0.40
1:A:500:THR:OG1	1:A:501:GLY:N	2.54	0.40
1:A:508:LYS:O	1:A:512:GLU:HG3	2.21	0.40
1:A:578:ALA:O	1:A:582:ILE:HG13	2.22	0.40

All (6) 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)
3:A:919:HOH:O	3:A:960:HOH:O[2_555]	1.92	0.28
3:A:935:HOH:O	3:B:849:HOH:O[4_445]	2.04	0.16

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:746:HOH:O	3:B:848:HOH:O[2_455]	2.06	0.14
3:A:1002:HOH:O	3:B:913:HOH:O[4_445]	2.08	0.12
3:A:975:HOH:O	3:B:911:HOH:O[4_445]	2.09	0.11
3:B:848:HOH:O	3:B:848:HOH:O[2_455]	2.19	0.01

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	557/592 (94%)	525 (94%)	29 (5%)	3 (0%)	24	43
1	B	557/592 (94%)	520 (93%)	32 (6%)	5 (1%)	14	27
All	All	1114/1184 (94%)	1045 (94%)	61 (6%)	8 (1%)	18	34

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	198	ILE
1	A	389	PRO
1	B	198	ILE
1	B	268	GLU
1	B	47	SER
1	B	295	ILE
1	A	268	GLU
1	B	389	PRO

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	451/473 (95%)	439 (97%)	12 (3%)	39	67
1	B	451/473 (95%)	436 (97%)	15 (3%)	33	61
All	All	902/946 (95%)	875 (97%)	27 (3%)	36	64

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	198	ILE
1	A	219	ARG
1	A	275	HIS
1	A	277	ILE
1	A	295	ILE
1	A	317	ARG
1	A	386	SER
1	A	387	LEU
1	A	425	ILE
1	A	456	VAL
1	A	484	ARG
1	A	520	ILE
1	B	42	SER
1	B	47	SER
1	B	198	ILE
1	B	219	ARG
1	B	277	ILE
1	B	298	GLN
1	B	317	ARG
1	B	330	VAL
1	B	370	ARG
1	B	400	LEU
1	B	484	ARG
1	B	495	ARG
1	B	512	GLU
1	B	515	SER
1	B	536	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	205	ASN

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Mol	Chain	Res	Type
1	A	275	HIS
1	A	531	GLN
1	B	48	HIS
1	B	51	ASN
1	B	120	GLN
1	B	242	GLN
1	B	296	ASN
1	B	298	GLN
1	B	331	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	B	601	-	58,58,58	1.98	15 (25%)	85,89,89	2.15	19 (22%)
2	FAD	A	601	-	58,58,58	2.10	14 (24%)	85,89,89	2.05	17 (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	B	601	-	1/1/9/9	13/34/50/50	0/6/6/6
2	FAD	A	601	-	1/1/9/9	13/34/50/50	0/6/6/6

All (29) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	601	FAD	C1'-C2'	7.46	1.63	1.52
2	B	601	FAD	C1'-C2'	6.90	1.62	1.52
2	A	601	FAD	C5X-N5	5.58	1.49	1.39
2	A	601	FAD	C2'-C3'	5.51	1.63	1.53
2	B	601	FAD	C5X-N5	5.31	1.49	1.39
2	A	601	FAD	C2-N1	5.06	1.48	1.36
2	B	601	FAD	C2-N1	4.97	1.47	1.36
2	B	601	FAD	C2'-C3'	4.67	1.61	1.53
2	A	601	FAD	C6A-N6A	3.47	1.43	1.34
2	B	601	FAD	C6A-N6A	3.30	1.42	1.34
2	B	601	FAD	C4X-C10	2.98	1.53	1.44
2	A	601	FAD	O2'-C2'	2.91	1.49	1.43
2	B	601	FAD	O2'-C2'	2.88	1.49	1.43
2	A	601	FAD	C5A-N7A	-2.86	1.33	1.39
2	A	601	FAD	C4X-C10	2.85	1.52	1.44
2	B	601	FAD	C5A-N7A	-2.58	1.34	1.39
2	A	601	FAD	O3'-C3'	-2.43	1.36	1.43
2	B	601	FAD	C2A-N3A	2.41	1.38	1.33
2	B	601	FAD	O3'-C3'	-2.33	1.37	1.43
2	A	601	FAD	C9A-N10	2.29	1.45	1.41
2	B	601	FAD	PA-O3P	-2.28	1.57	1.59
2	A	601	FAD	O2B-C2B	-2.24	1.37	1.43
2	B	601	FAD	O2B-C2B	-2.12	1.37	1.43
2	B	601	FAD	C2B-C3B	-2.12	1.47	1.53
2	B	601	FAD	P-O1P	2.12	1.58	1.50
2	A	601	FAD	PA-O1A	2.08	1.58	1.50
2	A	601	FAD	O3B-C3B	-2.08	1.37	1.43
2	A	601	FAD	C10-N10	2.07	1.41	1.37
2	B	601	FAD	PA-O1A	2.00	1.57	1.50

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	601	FAD	O2P-P-O3P	6.44	124.69	107.27
2	A	601	FAD	O2P-P-O3P	5.97	123.42	107.27
2	B	601	FAD	O5'-P-O1P	-5.94	85.38	108.94
2	A	601	FAD	N3A-C2A-N1A	-5.91	119.63	128.58
2	B	601	FAD	C5A-C4A-N3A	-5.84	118.67	126.72
2	A	601	FAD	O5'-P-O1P	-5.60	86.73	108.94
2	B	601	FAD	N3A-C2A-N1A	-5.55	120.18	128.58
2	A	601	FAD	O2P-P-O5'	-5.37	83.22	107.57
2	B	601	FAD	C4'-C3'-C2'	5.23	122.27	113.57
2	A	601	FAD	C5A-C4A-N3A	-4.92	119.94	126.72
2	B	601	FAD	O2P-P-O5'	-4.82	85.71	107.57
2	A	601	FAD	O2'-C2'-C3'	4.58	119.96	109.25
2	A	601	FAD	C1'-C2'-C3'	4.55	122.00	109.66
2	B	601	FAD	O2'-C2'-C3'	4.34	119.41	109.25
2	B	601	FAD	N3A-C4A-N9A	4.26	134.40	127.17
2	B	601	FAD	O3'-C3'-C4'	4.20	118.48	108.93
2	A	601	FAD	N3A-C4A-N9A	4.15	134.22	127.17
2	B	601	FAD	C5A-N7A-C8A	4.04	109.80	103.45
2	B	601	FAD	O3'-C3'-C2'	3.97	117.95	108.93
2	B	601	FAD	C2A-N3A-C4A	3.83	121.19	111.83
2	A	601	FAD	C5A-N7A-C8A	3.76	109.36	103.45
2	A	601	FAD	C5'-C4'-C3'	3.62	119.05	112.22
2	A	601	FAD	C2A-N3A-C4A	3.53	120.46	111.83
2	B	601	FAD	C4A-C5A-N7A	-3.40	106.69	110.58
2	B	601	FAD	C1'-C2'-C3'	3.30	118.60	109.66
2	A	601	FAD	N9A-C8A-N7A	-3.10	109.53	113.94
2	A	601	FAD	O3P-P-O1P	3.06	119.92	110.70
2	B	601	FAD	N9A-C8A-N7A	-2.93	109.78	113.94
2	B	601	FAD	O3P-P-O1P	2.73	118.91	110.70
2	A	601	FAD	O4-C4-C4X	-2.53	119.84	126.53
2	B	601	FAD	O3P-PA-O1A	-2.53	103.10	110.70
2	A	601	FAD	C4A-C5A-N7A	-2.46	107.77	110.58
2	A	601	FAD	O3P-PA-O1A	-2.38	103.56	110.70
2	B	601	FAD	O2'-C2'-C1'	2.26	119.49	110.20
2	B	601	FAD	O4B-C4B-C5B	2.25	116.55	109.33
2	A	601	FAD	O2A-PA-O3P	2.04	112.79	107.27

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	601	FAD	C3'
2	B	601	FAD	C3'

All (26) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	601	FAD	C1'-C2'-C3'-O3'
2	A	601	FAD	C1'-C2'-C3'-C4'
2	A	601	FAD	O2'-C2'-C3'-C4'
2	A	601	FAD	C2'-C3'-C4'-O4'
2	A	601	FAD	C2'-C3'-C4'-C5'
2	A	601	FAD	O3'-C3'-C4'-C5'
2	A	601	FAD	C5'-O5'-P-O1P
2	A	601	FAD	C5'-O5'-P-O3P
2	B	601	FAD	C5B-O5B-PA-O2A
2	B	601	FAD	C5B-O5B-PA-O3P
2	B	601	FAD	C1'-C2'-C3'-O3'
2	B	601	FAD	C1'-C2'-C3'-C4'
2	B	601	FAD	O2'-C2'-C3'-C4'
2	B	601	FAD	C2'-C3'-C4'-O4'
2	B	601	FAD	C5'-O5'-P-O1P
2	B	601	FAD	C5'-O5'-P-O3P
2	B	601	FAD	C2'-C3'-C4'-C5'
2	A	601	FAD	O3'-C3'-C4'-O4'
2	A	601	FAD	O2'-C2'-C3'-O3'
2	A	601	FAD	P-O3P-PA-O5B
2	B	601	FAD	P-O3P-PA-O5B
2	B	601	FAD	O4B-C4B-C5B-O5B
2	A	601	FAD	N10-C1'-C2'-C3'
2	B	601	FAD	C5B-O5B-PA-O1A
2	B	601	FAD	C3B-C4B-C5B-O5B
2	A	601	FAD	O4B-C4B-C5B-O5B

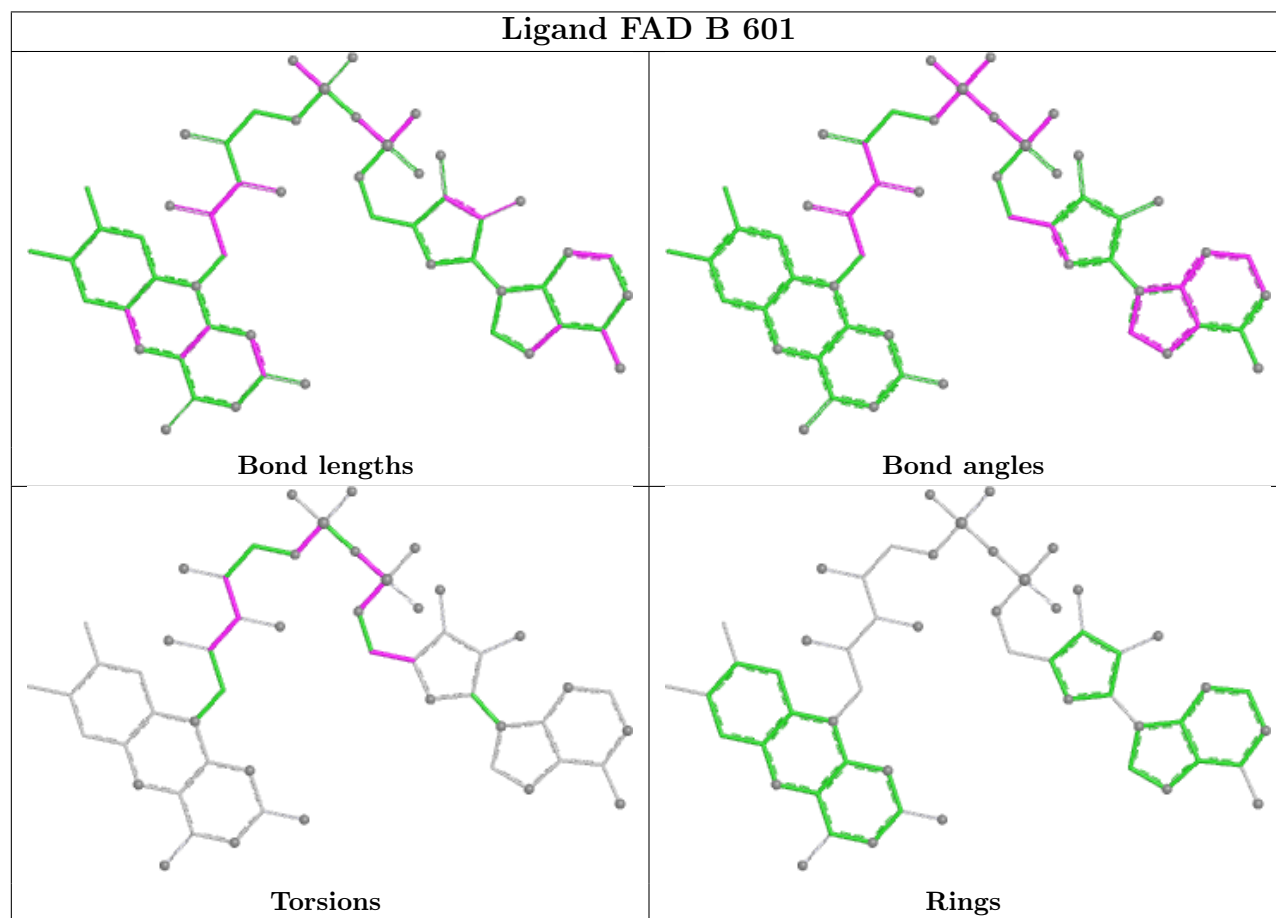
There are no ring outliers.

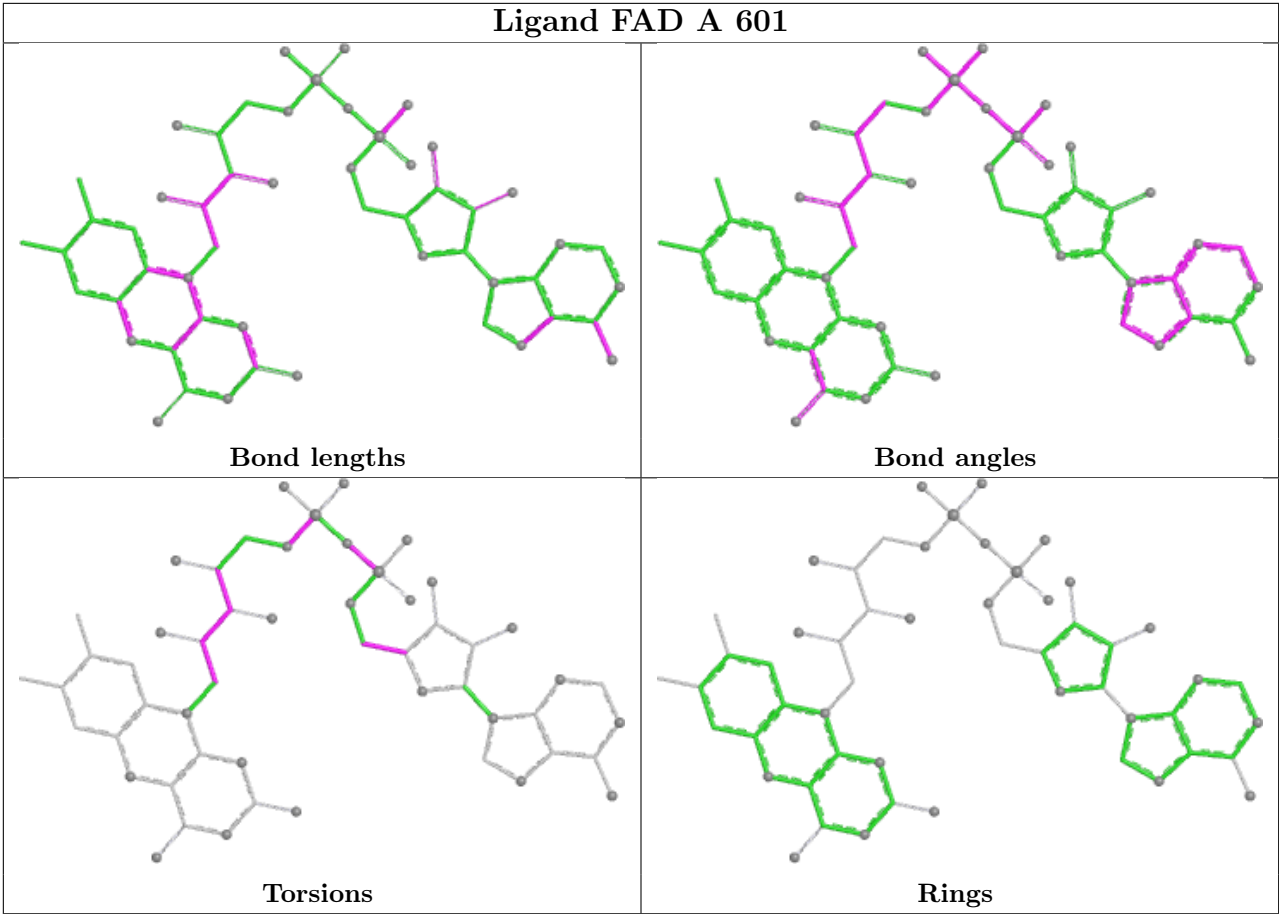
2 monomers are involved in 16 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	601	FAD	8	0
2	A	601	FAD	8	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 ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	4
1	B	3

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	217:GLU	C	218:HIS	N	1.18
1	A	218:HIS	C	219:ARG	N	1.17
1	B	51:ASN	C	52:GLN	N	1.17
1	B	169:VAL	C	170:ARG	N	1.17

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Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	149:GLU	C	150:ASP	N	1.07
1	A	46:ARG	C	47:SER	N	1.06
1	B	255:GLY	C	256:TYR	N	1.02

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	563/592 (95%)	0.16	33 (5%) 28 24	6, 22, 62, 157	0
1	B	563/592 (95%)	0.06	29 (5%) 33 29	6, 22, 62, 157	0
All	All	1126/1184 (95%)	0.11	62 (5%) 30 27	6, 22, 62, 157	0

All (62) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	198	ILE	8.4
1	A	202	GLY	7.0
1	B	198	ILE	6.3
1	B	268	GLU	5.5
1	A	197	GLY	5.4
1	A	223	MET	5.3
1	A	199	GLY	4.8
1	B	196	MET	4.7
1	A	294	THR	4.7
1	A	196	MET	4.6
1	B	223	MET	4.6
1	A	256	TYR	4.6
1	A	195	GLN	4.6
1	A	201	SER	4.6
1	B	270	ALA	4.1
1	A	265	ILE	4.0
1	B	195	GLN	4.0
1	B	256	TYR	3.9
1	A	270	ALA	3.9
1	B	197	GLY	3.7
1	A	169	VAL	3.6
1	A	268	GLU	3.6
1	B	199	GLY	3.6
1	B	265	ILE	3.5

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Mol	Chain	Res	Type	RSRZ
1	B	405	ASP	3.4
1	B	200	ASP	3.3
1	B	267	LYS	3.2
1	B	201	SER	3.2
1	A	282	GLU	3.2
1	B	202	GLY	3.2
1	B	282	GLU	3.1
1	A	200	ASP	3.1
1	A	388	PRO	3.1
1	A	580	LYS	3.0
1	A	46	ARG	2.9
1	B	254	TRP	2.9
1	A	194	GLY	2.8
1	A	390	ALA	2.7
1	A	156	ALA	2.7
1	A	405	ASP	2.5
1	B	40	SER	2.5
1	A	45	PRO	2.5
1	B	294	THR	2.5
1	A	255	GLY	2.5
1	A	392	THR	2.4
1	B	388	PRO	2.4
1	B	46	ARG	2.4
1	B	390	ALA	2.4
1	B	42	SER	2.3
1	B	252	CYS	2.3
1	B	45	PRO	2.3
1	A	280	THR	2.2
1	B	409	GLU	2.2
1	A	267	LYS	2.2
1	B	406	ALA	2.1
1	A	295	ILE	2.1
1	A	409	GLU	2.1
1	A	389	PRO	2.1
1	A	522	LEU	2.1
1	B	238	VAL	2.0
1	B	194	GLY	2.0
1	A	394	SER	2.0

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

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

6.3 Carbohydrates [i](#)

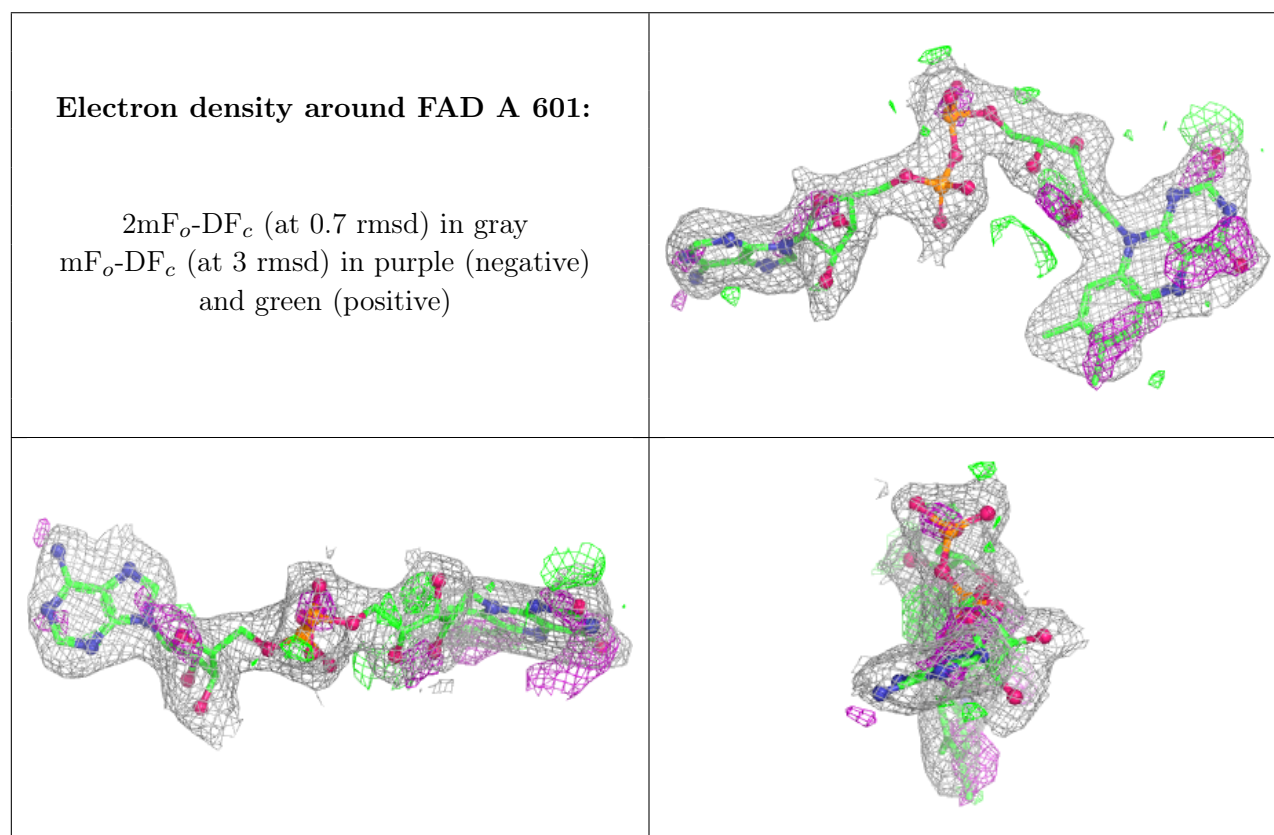
There are no oligosaccharides in this entry.

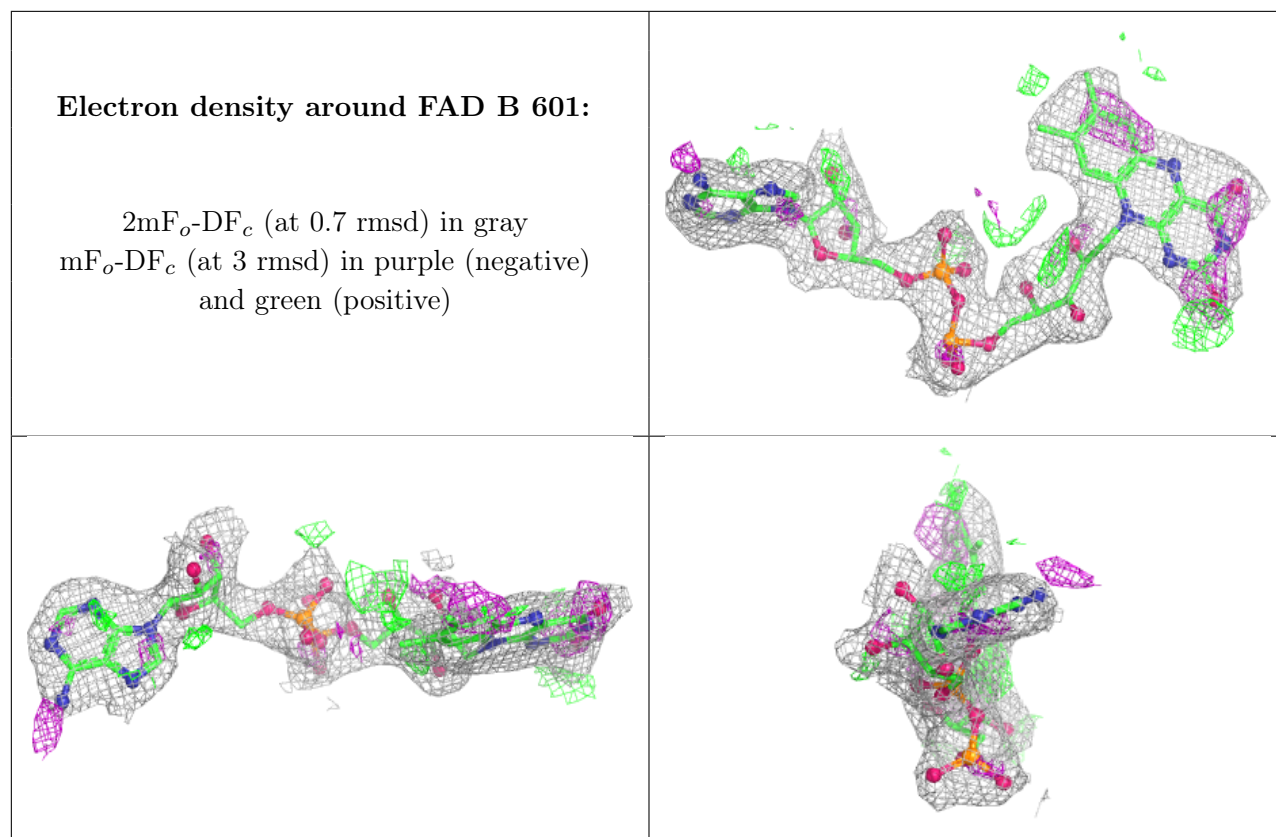
6.4 Ligands [i](#)

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	A	601	53/53	0.95	0.08	4,8,11,13	0
2	FAD	B	601	53/53	0.95	0.08	4,8,11,13	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.