



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 04:34 AM UTC

PDB ID : 4ML8 / pdb_00004ml8
Title : Structure of maize cytokinin oxidase/dehydrogenase 2 (ZmCKO2)
Authors : Morera, S.; Kopečný, D.; Briozzo, P.; Koncitikova, R.
Deposited on : 2013-09-06
Resolution : 2.70 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

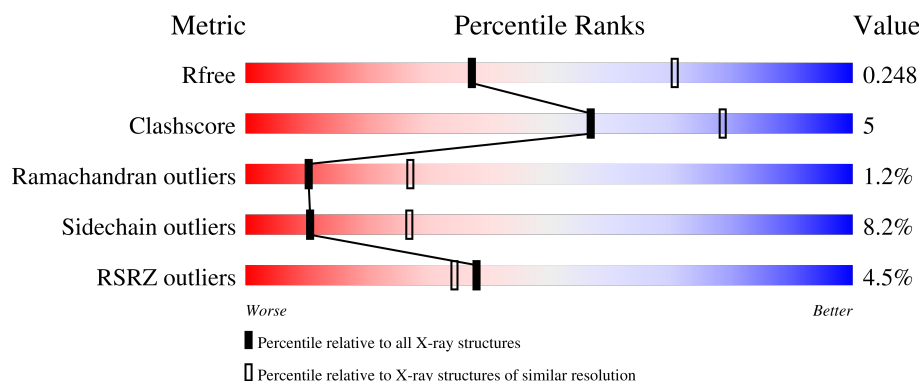
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	501	3% 76% 18% ...
1	B	501	5% 78% 17% ..
1	C	501	3% 79% 17% ..
1	D	501	6% 78% 15% ..

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	PEG	A	602	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 15588 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 Cytokinin oxidase 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	488	Total	C	N	O	S	0	0	0
			3842	2442	677	709	14			
1	B	487	Total	C	N	O	S	0	0	0
			3835	2437	676	708	14			
1	C	488	Total	C	N	O	S	0	0	0
			3842	2442	677	709	14			
1	D	479	Total	C	N	O	S	0	0	0
			3772	2398	665	695	14			

There are 16 discrepancies between the modelled and reference sequences:

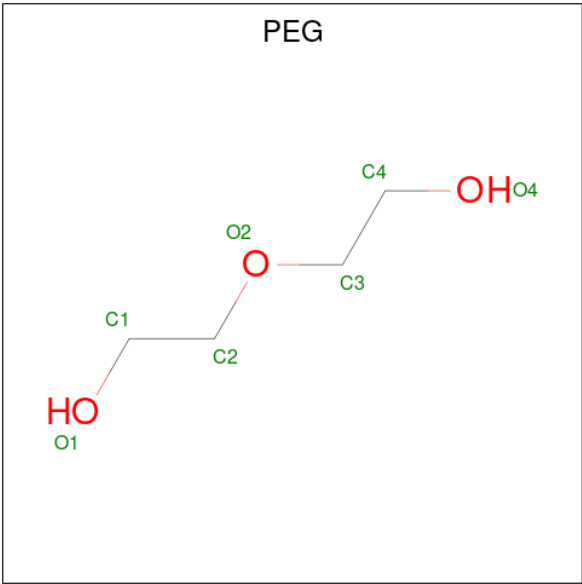
Chain	Residue	Modelled	Actual	Comment	Reference
A	16	ALA	-	expression tag	UNP Q709Q5
A	17	GLY	-	expression tag	UNP Q709Q5
A	18	HIS	-	expression tag	UNP Q709Q5
A	19	MET	-	expression tag	UNP Q709Q5
B	16	ALA	-	expression tag	UNP Q709Q5
B	17	GLY	-	expression tag	UNP Q709Q5
B	18	HIS	-	expression tag	UNP Q709Q5
B	19	MET	-	expression tag	UNP Q709Q5
C	16	ALA	-	expression tag	UNP Q709Q5
C	17	GLY	-	expression tag	UNP Q709Q5
C	18	HIS	-	expression tag	UNP Q709Q5
C	19	MET	-	expression tag	UNP Q709Q5
D	16	ALA	-	expression tag	UNP Q709Q5
D	17	GLY	-	expression tag	UNP Q709Q5
D	18	HIS	-	expression tag	UNP Q709Q5
D	19	MET	-	expression tag	UNP Q709Q5

- 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		
2	C	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	D	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 3 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: C₄H₁₀O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			7	4	3		

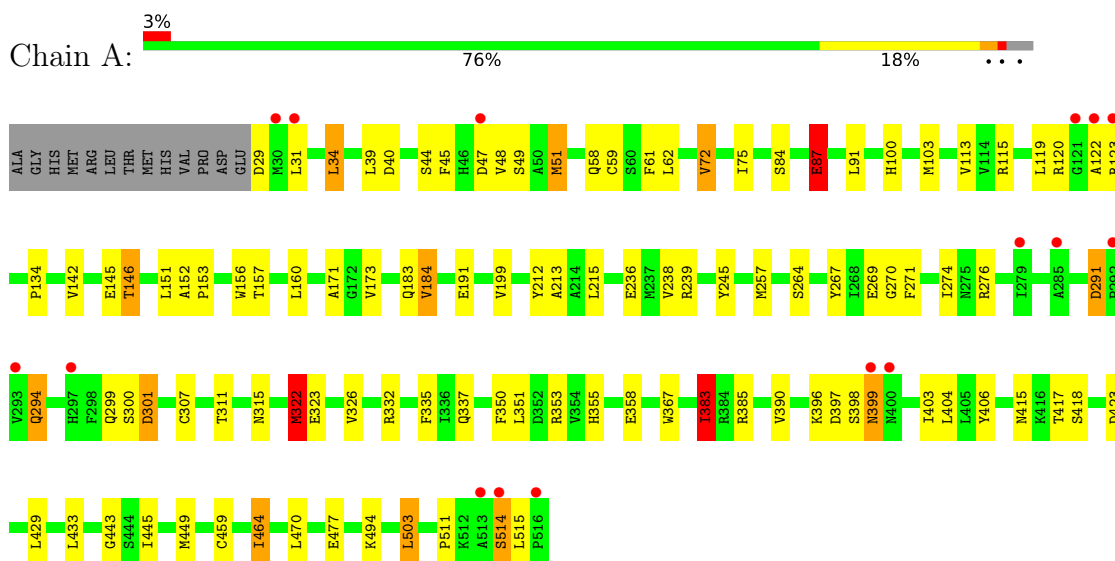
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	28	Total	O	0	0
			28	28		
4	B	22	Total	O	0	0
			22	22		
4	C	19	Total	O	0	0
			19	19		
4	D	9	Total	O	0	0
			9	9		

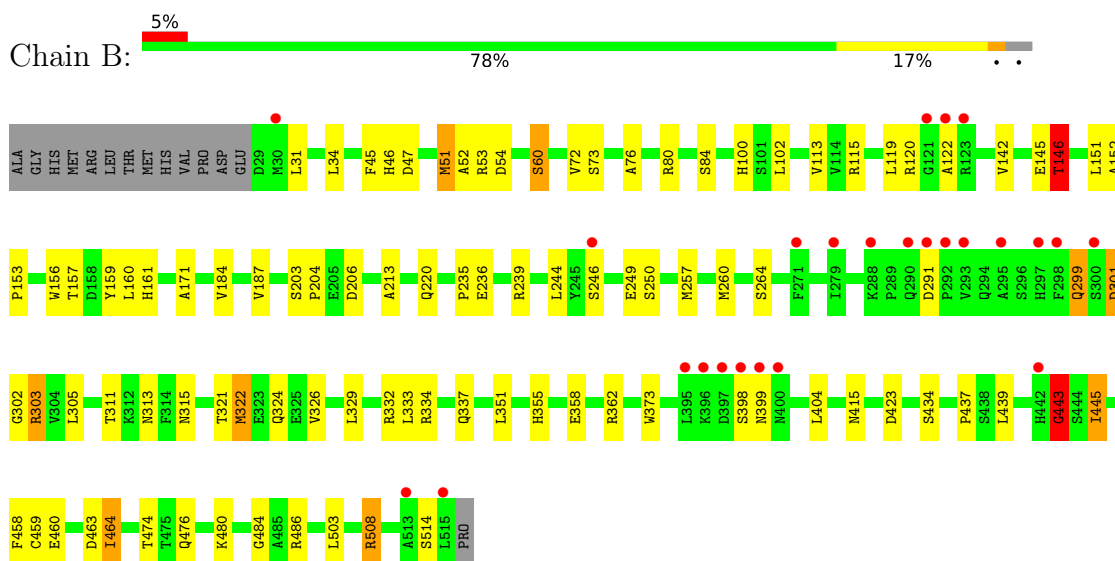
3 Residue-property plots [i](#)

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.

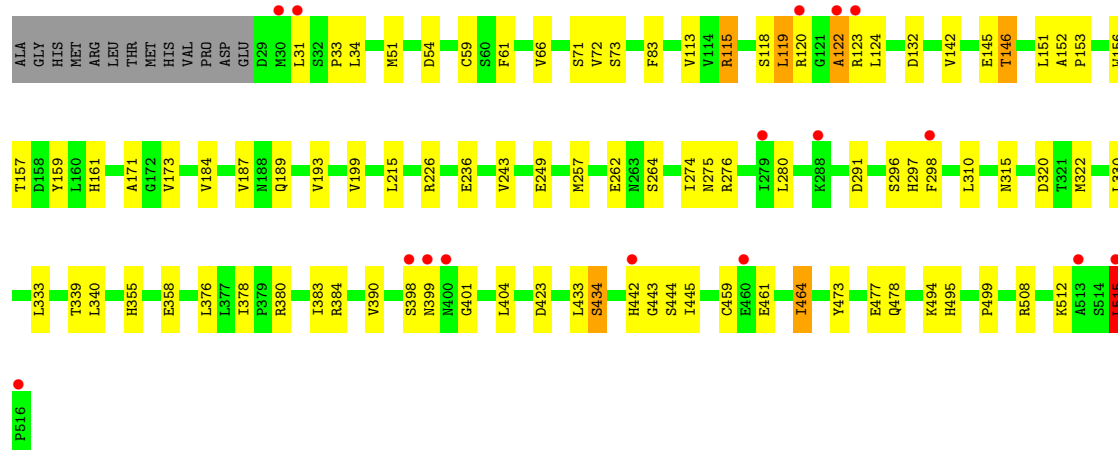
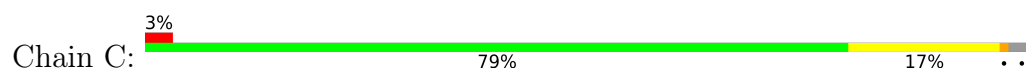
• Molecule 1: Cytokinin oxidase 2



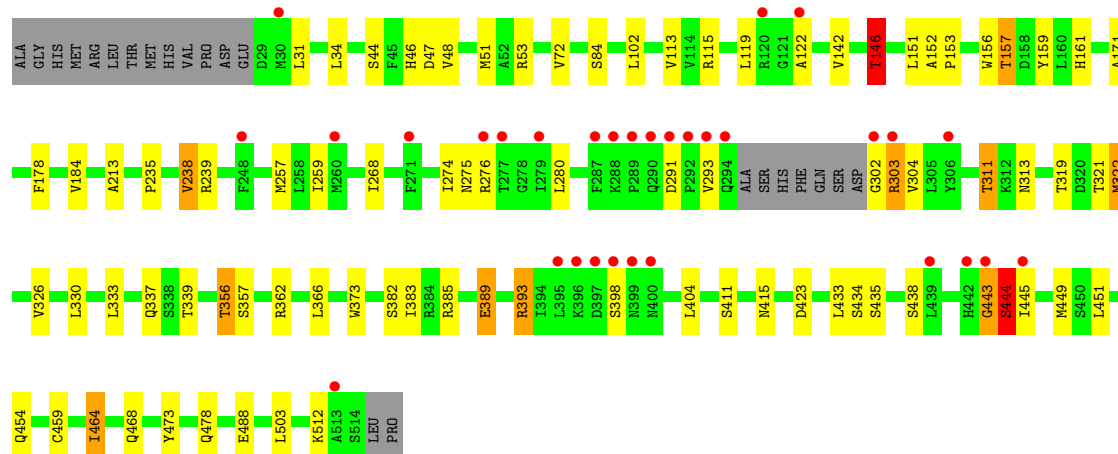
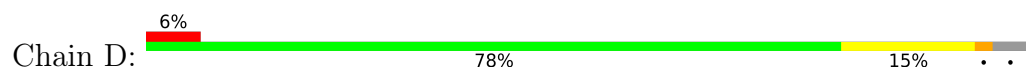
• Molecule 1: Cytokinin oxidase 2



• Molecule 1: Cytokinin oxidase 2



• Molecule 1: Cytokinin oxidase 2



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	183.21Å 76.67Å 201.04Å 90.00° 92.08° 90.00°	Depositor
Resolution (Å)	38.37 – 2.70 38.37 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.8 (38.37-2.70) 99.7 (38.37-2.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.79 (at 2.69Å)	Xtriage
Refinement program	BUSTER 2.10.0	Depositor
R, R_{free}	0.187 , 0.237 0.197 , 0.248	Depositor DCC
R_{free} test set	3859 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	46.3	Xtriage
Anisotropy	0.487	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 46.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.115 for -h,-k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	15588	wwPDB-VP
Average B, all atoms (Å ²)	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 19.73% 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: PEG, 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	0.92	2/3939 (0.1%)	1.39	27/5348 (0.5%)
1	B	0.90	1/3931 (0.0%)	1.39	27/5336 (0.5%)
1	C	0.87	0/3939	1.38	17/5348 (0.3%)
1	D	0.87	0/3865	1.40	24/5245 (0.5%)
All	All	0.89	3/15674 (0.0%)	1.39	95/21277 (0.4%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	322	MET	SD-CE	-5.72	1.65	1.79
1	A	51	MET	SD-CE	-5.50	1.65	1.79
1	B	187	VAL	CA-C	5.40	1.58	1.52

All (95) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	514	SER	N-CA-C	9.11	122.89	111.69
1	C	383	ILE	N-CA-C	7.68	118.43	110.36
1	B	514	SER	N-CA-C	7.68	119.56	111.03
1	A	443	GLY	CA-C-N	7.35	134.94	121.70
1	A	443	GLY	C-N-CA	7.35	134.94	121.70
1	A	383	ILE	N-CA-C	7.21	117.35	110.42
1	C	157	THR	N-CA-C	-7.01	99.84	110.14
1	C	443	GLY	CA-C-N	6.82	133.98	121.70
1	C	443	GLY	C-N-CA	6.82	133.98	121.70
1	C	33	PRO	CA-C-N	6.81	129.70	120.44
1	C	33	PRO	C-N-CA	6.81	129.70	120.44
1	A	399	ASN	CA-C-N	6.77	130.03	120.28
1	A	399	ASN	C-N-CA	6.77	130.03	120.28
1	B	246	SER	N-CA-C	-6.75	105.99	114.56

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	142	VAL	N-CA-C	-6.68	104.01	110.42
1	C	54	ASP	CA-CB-CG	6.59	119.19	112.60
1	A	142	VAL	N-CA-C	-6.59	104.09	110.42
1	B	301	ASP	CA-CB-CG	6.55	119.15	112.60
1	D	423	ASP	N-CA-C	6.41	120.36	112.54
1	A	300	SER	CA-C-N	6.31	131.29	121.26
1	A	300	SER	C-N-CA	6.31	131.29	121.26
1	B	443	GLY	CA-C-N	6.27	132.99	121.70
1	B	443	GLY	C-N-CA	6.27	132.99	121.70
1	C	145	GLU	CA-C-N	6.23	129.25	120.28
1	C	145	GLU	C-N-CA	6.23	129.25	120.28
1	B	333	LEU	N-CA-C	6.22	119.39	109.24
1	A	145	GLU	CA-C-N	6.20	129.21	120.28
1	A	145	GLU	C-N-CA	6.20	129.21	120.28
1	B	220	GLN	N-CA-C	6.13	118.93	111.82
1	B	54	ASP	CA-CB-CG	6.07	118.67	112.60
1	D	321	THR	CA-C-N	6.05	128.68	120.44
1	D	321	THR	C-N-CA	6.05	128.68	120.44
1	C	142	VAL	N-CA-C	-5.89	104.77	110.42
1	D	443	GLY	CA-C-N	5.83	132.20	121.70
1	D	443	GLY	C-N-CA	5.83	132.20	121.70
1	D	459	CYS	CA-C-N	5.81	128.34	120.44
1	D	459	CYS	C-N-CA	5.81	128.34	120.44
1	D	444	SER	CA-C-N	5.79	127.93	120.70
1	D	444	SER	C-N-CA	5.79	127.93	120.70
1	D	503	LEU	N-CA-C	5.76	119.30	110.20
1	B	157	THR	N-CA-C	-5.62	101.43	110.14
1	C	390	VAL	N-CA-C	5.58	115.67	110.42
1	D	157	THR	N-CA-C	-5.55	100.83	109.72
1	A	423	ASP	N-CA-C	5.55	118.09	111.71
1	D	213	ALA	CA-C-N	5.55	127.72	120.28
1	D	213	ALA	C-N-CA	5.55	127.72	120.28
1	A	87	GLU	N-CA-C	5.54	118.85	111.75
1	D	383	ILE	N-CA-C	5.52	116.30	110.72
1	B	145	GLU	CA-C-N	5.51	127.66	120.28
1	B	145	GLU	C-N-CA	5.51	127.66	120.28
1	A	390	VAL	N-CA-C	5.50	115.70	110.53
1	B	503	LEU	CA-C-N	5.50	130.86	124.21
1	B	503	LEU	C-N-CA	5.50	130.86	124.21
1	B	100	HIS	N-CA-C	5.49	119.85	113.20
1	C	122	ALA	CA-C-N	5.48	132.01	121.54
1	C	122	ALA	C-N-CA	5.48	132.01	121.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	146	THR	CB-CA-C	5.47	121.16	110.67
1	A	385	ARG	CA-C-N	5.33	127.69	120.65
1	A	385	ARG	C-N-CA	5.33	127.69	120.65
1	C	512	LYS	CA-C-N	5.32	130.29	122.63
1	C	512	LYS	C-N-CA	5.32	130.29	122.63
1	D	46	HIS	N-CA-C	5.31	120.00	112.93
1	D	356	THR	CB-CA-C	5.31	119.60	110.79
1	B	445	ILE	N-CA-CB	5.28	118.50	110.58
1	B	484	GLY	CA-C-N	5.28	127.88	120.28
1	B	484	GLY	C-N-CA	5.28	127.88	120.28
1	C	515	LEU	N-CA-C	5.27	118.99	109.44
1	B	213	ALA	CA-C-N	5.27	128.32	120.31
1	B	213	ALA	C-N-CA	5.27	128.32	120.31
1	B	458	PHE	CA-C-N	5.27	127.60	120.38
1	B	458	PHE	C-N-CA	5.27	127.60	120.38
1	B	146	THR	CB-CA-C	5.25	119.51	110.79
1	B	398	SER	N-CA-C	-5.21	106.23	112.59
1	C	378	ILE	N-CA-C	5.21	114.20	108.05
1	A	301	ASP	CA-CB-CG	5.19	117.79	112.60
1	B	423	ASP	N-CA-C	5.18	118.70	112.38
1	A	301	ASP	N-CA-C	5.16	118.04	110.30
1	A	40	ASP	N-CA-C	-5.15	105.75	111.36
1	B	362	ARG	CA-C-N	5.14	127.12	120.44
1	B	362	ARG	C-N-CA	5.14	127.12	120.44
1	A	48	VAL	CA-C-N	5.13	127.41	120.38
1	A	48	VAL	C-N-CA	5.13	127.41	120.38
1	A	183	GLN	CA-C-N	5.12	128.56	120.47
1	A	183	GLN	C-N-CA	5.12	128.56	120.47
1	D	385	ARG	CA-C-N	5.10	127.37	120.44
1	D	385	ARG	C-N-CA	5.10	127.37	120.44
1	D	362	ARG	CA-C-N	5.06	127.06	120.28
1	D	362	ARG	C-N-CA	5.06	127.06	120.28
1	A	503	LEU	CA-C-N	5.03	130.30	124.21
1	A	503	LEU	C-N-CA	5.03	130.30	124.21
1	D	48	VAL	CA-C-N	5.03	127.33	120.54
1	D	48	VAL	C-N-CA	5.03	127.33	120.54
1	B	46	HIS	N-CA-C	5.01	119.13	112.72
1	A	157	THR	N-CA-C	-5.00	101.94	109.85
1	A	503	LEU	N-CA-C	5.00	118.10	110.20

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	3842	0	3758	49	0
1	B	3835	0	3751	37	0
1	C	3842	0	3758	31	0
1	D	3772	0	3696	26	0
2	A	53	0	29	0	0
2	B	53	0	29	0	0
2	C	53	0	29	0	0
2	D	53	0	29	0	0
3	A	7	0	10	4	0
4	A	28	0	0	0	0
4	B	22	0	0	0	0
4	C	19	0	0	0	0
4	D	9	0	0	0	0
All	All	15588	0	15089	141	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 5.

All (141) 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:51:MET:HE2	1:B:115:ARG:HG3	1.54	0.89
1:B:51:MET:CE	1:B:115:ARG:HG3	2.07	0.85
1:D:239:ARG:HD2	1:D:322:MET:HE1	1.58	0.85
1:A:367:TRP:HE1	3:A:602:PEG:H41	1.44	0.82
1:B:146:THR:HG23	1:B:151:LEU:O	1.82	0.79
1:A:274:ILE:HB	1:A:398:SER:HB3	1.71	0.72
1:B:302:GLY:HA2	1:B:303:ARG:HB3	1.70	0.71
1:D:382:SER:HB3	1:D:464:ILE:HG13	1.72	0.70
1:A:100:HIS:HE2	3:A:602:PEG:H42	1.59	0.68
1:B:146:THR:CG2	1:B:152:ALA:HA	2.26	0.65
1:A:51:MET:HE3	1:A:113:VAL:HG11	1.80	0.63
1:D:146:THR:HG23	1:D:151:LEU:O	1.99	0.62
1:A:236:GLU:HG2	1:A:315:ASN:HA	1.81	0.62
1:D:274:ILE:HB	1:D:398:SER:HB3	1.81	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:51:MET:HE3	1:C:113:VAL:HG11	1.81	0.62
1:C:146:THR:CG2	1:C:152:ALA:HA	2.30	0.62
1:B:146:THR:HG21	1:B:152:ALA:HA	1.83	0.61
1:A:58:GLN:CD	1:A:103:MET:HE1	2.27	0.59
1:A:184:VAL:HG22	1:A:215:LEU:HB3	1.84	0.59
1:C:132:ASP:HB3	1:C:226:ARG:HD2	1.85	0.59
1:A:100:HIS:NE2	3:A:602:PEG:H42	2.17	0.58
1:D:330:LEU:HA	1:D:333:LEU:HD12	1.86	0.58
1:C:146:THR:HG21	1:C:152:ALA:HA	1.86	0.57
1:C:59:CYS:HB3	1:C:61:PHE:CE2	2.39	0.57
1:B:236:GLU:HG2	1:B:315:ASN:HA	1.87	0.56
1:C:494:LYS:HG3	1:C:515:LEU:HD11	1.88	0.56
1:C:274:ILE:HB	1:C:398:SER:HB3	1.89	0.55
1:C:189:GLN:HE21	1:C:226:ARG:HH22	1.55	0.55
1:C:146:THR:HG23	1:C:151:LEU:O	2.07	0.54
1:D:389:GLU:HG3	1:D:393:ARG:HH11	1.70	0.54
1:D:303:ARG:HG2	1:D:304:VAL:H	1.71	0.54
1:C:193:VAL:HG22	1:C:199:VAL:HG22	1.90	0.54
1:C:495:HIS:HB2	1:C:515:LEU:HD22	1.90	0.54
1:A:146:THR:HG23	1:A:151:LEU:O	2.08	0.54
1:D:146:THR:HG21	1:D:153:PRO:HD3	1.90	0.53
1:B:355:HIS:HD2	1:B:358:GLU:OE1	1.91	0.53
1:A:123:ARG:HB3	1:A:134:PRO:HG3	1.90	0.53
1:A:58:GLN:NE2	1:A:103:MET:HE1	2.23	0.53
1:A:59:CYS:HB3	1:A:61:PHE:CE2	2.45	0.52
1:C:51:MET:HE3	1:C:113:VAL:CG1	2.39	0.52
1:D:178:PHE:HB2	1:D:259:ILE:HB	1.92	0.51
1:A:212:TYR:O	1:A:417:THR:HA	2.11	0.51
1:D:51:MET:HE3	1:D:113:VAL:HG11	1.93	0.51
1:B:302:GLY:CA	1:B:303:ARG:HB3	2.40	0.51
1:A:269:GLU:HG2	1:A:406:TYR:CD1	2.46	0.51
1:B:153:PRO:HD2	1:B:156:TRP:CH2	2.45	0.51
1:C:122:ALA:O	1:C:124:LEU:N	2.44	0.50
1:A:160:LEU:HD11	1:A:351:LEU:HB3	1.92	0.50
1:D:280:LEU:HD11	1:D:339:THR:HG22	1.92	0.50
1:D:159:TYR:CZ	1:D:161:HIS:HB2	2.47	0.50
1:A:123:ARG:NH2	1:A:191:GLU:OE2	2.44	0.50
1:C:236:GLU:HG2	1:C:315:ASN:HA	1.93	0.49
1:B:51:MET:CE	1:B:115:ARG:CG	2.85	0.49
1:B:146:THR:HG21	1:B:153:PRO:HD3	1.94	0.49
1:B:146:THR:HG23	1:B:151:LEU:C	2.37	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:153:PRO:HD2	1:C:156:TRP:CH2	2.48	0.48
1:A:367:TRP:NE1	3:A:602:PEG:H41	2.21	0.48
1:B:235:PRO:HG3	1:B:313:ASN:HB3	1.96	0.48
1:C:187:VAL:HG11	1:C:215:LEU:HD21	1.96	0.48
1:B:322:MET:HE3	1:B:326:VAL:HG21	1.96	0.47
1:D:153:PRO:HD2	1:D:156:TRP:CH2	2.49	0.47
1:B:52:ALA:O	1:B:60:SER:HA	2.15	0.47
1:D:146:THR:CG2	1:D:152:ALA:HA	2.44	0.47
1:D:451:LEU:HA	1:D:454:GLN:HE21	1.80	0.47
1:B:142:VAL:O	1:B:146:THR:HB	2.14	0.47
1:B:160:LEU:HD11	1:B:351:LEU:HB3	1.96	0.47
1:A:350:PHE:O	1:A:353:ARG:HD3	2.15	0.47
1:B:159:TYR:CZ	1:B:161:HIS:HB2	2.49	0.47
1:C:71:SER:HA	1:C:119:LEU:HD21	1.96	0.47
1:C:459:CYS:HA	1:C:464:ILE:HD13	1.97	0.47
1:A:270:GLY:HA2	1:A:307:CYS:O	2.15	0.47
1:D:473:TYR:HB3	1:D:478:GLN:HB3	1.95	0.47
1:C:330:LEU:HA	1:C:333:LEU:HD12	1.97	0.46
1:B:146:THR:HG23	1:B:152:ALA:HA	1.96	0.46
1:B:51:MET:HE1	1:B:115:ARG:CG	2.46	0.46
1:C:159:TYR:CZ	1:C:161:HIS:HB2	2.51	0.46
1:A:39:LEU:O	1:B:355:HIS:HE1	1.99	0.46
1:B:239:ARG:HG2	1:B:322:MET:HE3	1.97	0.46
1:C:51:MET:HG3	1:C:66:VAL:HG21	1.98	0.46
1:C:115:ARG:O	1:C:118:SER:HB2	2.16	0.45
1:A:213:ALA:O	1:A:418:SER:OG	2.33	0.45
1:A:239:ARG:HD2	1:A:322:MET:HE1	1.98	0.45
1:A:404:LEU:HD13	1:A:433:LEU:HD22	1.98	0.45
1:C:404:LEU:HD13	1:C:433:LEU:HD22	1.98	0.45
1:D:438:SER:O	1:D:444:SER:HA	2.16	0.45
1:A:146:THR:CG2	1:A:152:ALA:HA	2.46	0.45
1:A:267:TYR:HB3	1:A:311:THR:OG1	2.17	0.45
1:B:244:LEU:HD22	1:B:305:LEU:HB3	1.97	0.45
1:B:459:CYS:HA	1:B:464:ILE:HD13	1.98	0.45
1:B:476:GLN:HG2	1:B:480:LYS:HE3	1.98	0.45
1:C:473:TYR:HB3	1:C:478:GLN:HB3	1.99	0.45
1:D:404:LEU:HD13	1:D:433:LEU:HD22	1.99	0.45
1:C:243:VAL:HG12	1:C:340:LEU:HA	1.99	0.45
1:B:31:LEU:HB2	1:B:45:PHE:CE1	2.51	0.45
1:A:239:ARG:NH1	1:A:322:MET:HE1	2.32	0.45
1:B:51:MET:HE1	1:B:115:ARG:HG3	1.97	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:494:LYS:HG3	1:A:515:LEU:HD11	1.99	0.44
1:B:474:THR:HA	1:B:508:ARG:HD3	1.98	0.44
1:D:51:MET:HE2	1:D:115:ARG:HG3	1.98	0.44
1:D:302:GLY:HA2	1:D:303:ARG:HA	1.79	0.44
1:A:153:PRO:HD2	1:A:156:TRP:CH2	2.52	0.44
1:C:249:GLU:HG3	1:C:384:ARG:CZ	2.48	0.44
1:A:146:THR:HG21	1:A:153:PRO:HD3	1.99	0.44
1:A:51:MET:HE3	1:A:113:VAL:CG1	2.47	0.44
1:A:72:VAL:HA	1:A:75:ILE:HD12	2.00	0.43
1:C:59:CYS:HB3	1:C:61:PHE:HE2	1.80	0.43
1:D:102:LEU:HD13	1:D:373:TRP:HB2	2.00	0.43
1:B:299:GLN:O	1:B:303:ARG:NH2	2.52	0.43
1:A:383:ILE:HD12	1:A:383:ILE:HA	1.91	0.43
1:A:87:GLU:H	1:A:87:GLU:HG2	1.58	0.43
1:A:153:PRO:HD2	1:A:156:TRP:CZ2	2.54	0.43
1:A:31:LEU:HA	1:A:34:LEU:HD22	2.00	0.43
1:B:203:SER:HB2	1:B:204:PRO:CD	2.49	0.43
1:B:264:SER:HA	1:B:329:LEU:HD21	2.00	0.43
1:A:355:HIS:HD2	1:A:358:GLU:OE1	2.01	0.43
1:B:51:MET:HG3	1:B:113:VAL:HG11	2.02	0.42
1:A:91:LEU:HD23	1:A:91:LEU:HA	1.87	0.42
1:C:401:GLY:HA3	1:C:434:SER:HA	2.02	0.42
1:A:459:CYS:HA	1:A:464:ILE:HD13	2.00	0.42
1:B:76:ALA:O	1:B:80:ARG:HB2	2.20	0.42
1:A:511:PRO:O	1:A:514:SER:HB3	2.20	0.42
1:B:437:PRO:HD2	1:B:443:GLY:O	2.20	0.42
1:D:238:VAL:HG13	1:D:311:THR:HB	2.02	0.42
1:A:355:HIS:CD2	1:A:358:GLU:OE1	2.73	0.42
1:B:102:LEU:HD13	1:B:373:TRP:HB2	2.02	0.42
1:A:31:LEU:HB2	1:A:45:PHE:CE1	2.55	0.41
1:D:274:ILE:HG12	1:D:304:VAL:HG22	2.01	0.41
1:C:83:PHE:HE1	1:C:499:PRO:HG2	1.85	0.41
1:D:51:MET:CE	1:D:115:ARG:HG3	2.50	0.41
1:C:280:LEU:HD11	1:C:339:THR:HG22	2.03	0.41
1:A:322:MET:HE3	1:A:326:VAL:HG21	2.02	0.41
1:A:239:ARG:HH12	1:A:323:GLU:HG2	1.85	0.41
1:A:291:ASP:HB3	1:A:294:GLN:HE21	1.85	0.41
1:D:178:PHE:CZ	1:D:268:ILE:HD12	2.56	0.41
1:A:120:ARG:HH12	1:B:120:ARG:HB3	1.86	0.41
1:A:238:VAL:HG13	1:A:311:THR:HB	2.03	0.41
1:A:245:TYR:CE1	1:A:335:PHE:HB3	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:355:HIS:HD2	1:C:358:GLU:OE1	2.05	0.40
1:A:406:TYR:CZ	1:A:429:LEU:HD23	2.57	0.40
1:A:271:PHE:HA	1:A:403:ILE:O	2.22	0.40
1:D:235:PRO:HG3	1:D:313:ASN:HB3	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	486/501 (97%)	457 (94%)	25 (5%)	4 (1%)	16	37
1	B	485/501 (97%)	457 (94%)	23 (5%)	5 (1%)	12	32
1	C	486/501 (97%)	456 (94%)	23 (5%)	7 (1%)	9	23
1	D	475/501 (95%)	443 (93%)	25 (5%)	7 (2%)	8	22
All	All	1932/2004 (96%)	1813 (94%)	96 (5%)	23 (1%)	10	27

All (23) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	122	ALA
1	C	123	ARG
1	C	298	PHE
1	D	122	ALA
1	A	171	ALA
1	A	264	SER
1	B	171	ALA
1	D	171	ALA
1	D	276	ARG
1	C	264	SER
1	D	275	ASN

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Mol	Chain	Res	Type
1	B	122	ALA
1	C	171	ALA
1	C	291	ASP
1	C	297	HIS
1	B	303	ARG
1	C	444	SER
1	D	291	ASP
1	D	443	GLY
1	A	291	ASP
1	B	291	ASP
1	D	444	SER
1	B	443	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	414/425 (97%)	380 (92%)	34 (8%)	10	27
1	B	413/425 (97%)	377 (91%)	36 (9%)	9	24
1	C	414/425 (97%)	384 (93%)	30 (7%)	13	32
1	D	406/425 (96%)	371 (91%)	35 (9%)	10	25
All	All	1647/1700 (97%)	1512 (92%)	135 (8%)	10	27

All (135) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	29	ASP
1	A	34	LEU
1	A	44	SER
1	A	47	ASP
1	A	49	SER
1	A	62	LEU
1	A	72	VAL
1	A	84	SER
1	A	87	GLU

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Mol	Chain	Res	Type
1	A	115	ARG
1	A	119	LEU
1	A	146	THR
1	A	173	VAL
1	A	184	VAL
1	A	199	VAL
1	A	257	MET
1	A	276	ARG
1	A	294	GLN
1	A	299	GLN
1	A	301	ASP
1	A	322	MET
1	A	332	ARG
1	A	337	GLN
1	A	383	ILE
1	A	396	LYS
1	A	397	ASP
1	A	399	ASN
1	A	415	ASN
1	A	445	ILE
1	A	449	MET
1	A	464	ILE
1	A	470	LEU
1	A	477	GLU
1	A	503	LEU
1	B	34	LEU
1	B	47	ASP
1	B	51	MET
1	B	53	ARG
1	B	60	SER
1	B	72	VAL
1	B	73	SER
1	B	84	SER
1	B	119	LEU
1	B	146	THR
1	B	184	VAL
1	B	206	ASP
1	B	249	GLU
1	B	250	SER
1	B	257	MET
1	B	260	MET
1	B	299	GLN

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Mol	Chain	Res	Type
1	B	301	ASP
1	B	311	THR
1	B	321	THR
1	B	322	MET
1	B	324	GLN
1	B	332	ARG
1	B	334	ARG
1	B	337	GLN
1	B	399	ASN
1	B	404	LEU
1	B	415	ASN
1	B	434	SER
1	B	439	LEU
1	B	445	ILE
1	B	460	GLU
1	B	463	ASP
1	B	464	ILE
1	B	486	ARG
1	B	508	ARG
1	C	31	LEU
1	C	34	LEU
1	C	72	VAL
1	C	73	SER
1	C	115	ARG
1	C	119	LEU
1	C	120	ARG
1	C	146	THR
1	C	173	VAL
1	C	184	VAL
1	C	257	MET
1	C	262	GLU
1	C	275	ASN
1	C	276	ARG
1	C	296	SER
1	C	310	LEU
1	C	320	ASP
1	C	322	MET
1	C	376	LEU
1	C	380	ARG
1	C	399	ASN
1	C	423	ASP
1	C	434	SER

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Mol	Chain	Res	Type
1	C	442	HIS
1	C	445	ILE
1	C	461	GLU
1	C	464	ILE
1	C	477	GLU
1	C	508	ARG
1	C	515	LEU
1	D	31	LEU
1	D	34	LEU
1	D	44	SER
1	D	47	ASP
1	D	53	ARG
1	D	72	VAL
1	D	84	SER
1	D	119	LEU
1	D	146	THR
1	D	157	THR
1	D	184	VAL
1	D	238	VAL
1	D	257	MET
1	D	293	VAL
1	D	303	ARG
1	D	311	THR
1	D	319	THR
1	D	322	MET
1	D	326	VAL
1	D	337	GLN
1	D	356	THR
1	D	357	SER
1	D	366	LEU
1	D	389	GLU
1	D	393	ARG
1	D	411	SER
1	D	415	ASN
1	D	434	SER
1	D	435	SER
1	D	445	ILE
1	D	449	MET
1	D	464	ILE
1	D	468	GLN
1	D	488	GLU
1	D	512	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (39) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	42	HIS
1	A	58	GLN
1	A	188	ASN
1	A	207	ASN
1	A	275	ASN
1	A	282	ASN
1	A	355	HIS
1	A	364	GLN
1	A	399	ASN
1	A	415	ASN
1	A	447	HIS
1	B	42	HIS
1	B	81	HIS
1	B	161	HIS
1	B	207	ASN
1	B	275	ASN
1	B	297	HIS
1	B	355	HIS
1	B	400	ASN
1	C	58	GLN
1	C	68	HIS
1	C	188	ASN
1	C	189	GLN
1	C	207	ASN
1	C	282	ASN
1	C	355	HIS
1	C	400	ASN
1	C	447	HIS
1	C	476	GLN
1	D	81	HIS
1	D	186	ASN
1	D	188	ASN
1	D	207	ASN
1	D	364	GLN
1	D	409	ASN
1	D	454	GLN
1	D	468	GLN
1	D	476	GLN
1	D	478	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

5 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	PEG	A	602	-	6,6,6	0.26	0	5,5,5	0.17	0
2	FAD	A	601	1	58,58,58	0.92	1 (1%)	85,89,89	1.02	6 (7%)
2	FAD	C	601	1	58,58,58	0.86	1 (1%)	85,89,89	0.89	5 (5%)
2	FAD	D	601	1	58,58,58	0.61	1 (1%)	85,89,89	1.10	7 (8%)
2	FAD	B	601	1	58,58,58	0.54	1 (1%)	85,89,89	0.96	5 (5%)

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
3	PEG	A	602	-	-	4/4/4/4	-
2	FAD	A	601	1	-	6/34/50/50	0/6/6/6
2	FAD	C	601	1	-	1/34/50/50	0/6/6/6
2	FAD	D	601	1	-	3/34/50/50	0/6/6/6

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	FAD	B	601	1	-	4/34/50/50	0/6/6/6

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	601	FAD	PA-O3P	6.07	1.66	1.59
2	C	601	FAD	PA-O3P	5.78	1.65	1.59
2	D	601	FAD	PA-O3P	3.73	1.63	1.59
2	B	601	FAD	PA-O3P	3.04	1.62	1.59

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	601	FAD	C4'-C3'-C2'	5.11	122.08	113.57
2	B	601	FAD	O3P-P-O1P	-5.10	95.36	110.70
2	C	601	FAD	O3P-P-O1P	-4.31	97.74	110.70
2	D	601	FAD	O3P-P-O1P	-4.27	97.86	110.70
2	D	601	FAD	C4'-C3'-C2'	4.19	120.55	113.57
2	D	601	FAD	O3P-PA-O1A	-4.08	98.42	110.70
2	A	601	FAD	O3P-P-O1P	-3.52	100.11	110.70
2	B	601	FAD	O3P-PA-O1A	-3.19	101.10	110.70
2	B	601	FAD	C4'-C3'-C2'	2.86	118.33	113.57
2	D	601	FAD	O2P-P-O5'	-2.75	95.09	107.57
2	A	601	FAD	C5'-C4'-C3'	2.65	117.22	112.22
2	B	601	FAD	O2P-P-O5'	-2.49	96.28	107.57
2	B	601	FAD	O5'-P-O1P	2.49	118.78	108.94
2	D	601	FAD	O5'-C5'-C4'	2.45	115.89	109.36
2	A	601	FAD	O2P-P-O5'	-2.44	96.50	107.57
2	A	601	FAD	O5'-P-O1P	2.43	118.56	108.94
2	A	601	FAD	O3P-PA-O1A	-2.42	103.42	110.70
2	C	601	FAD	C4'-C3'-C2'	2.30	117.40	113.57
2	C	601	FAD	O2P-P-O5'	-2.29	97.16	107.57
2	D	601	FAD	O5'-P-O1P	2.24	117.81	108.94
2	C	601	FAD	C5'-C4'-C3'	2.20	116.36	112.22
2	C	601	FAD	O5'-P-O1P	2.15	117.44	108.94
2	D	601	FAD	C5'-C4'-C3'	2.11	116.20	112.22

There are no chirality outliers.

All (18) torsion outliers are listed below:

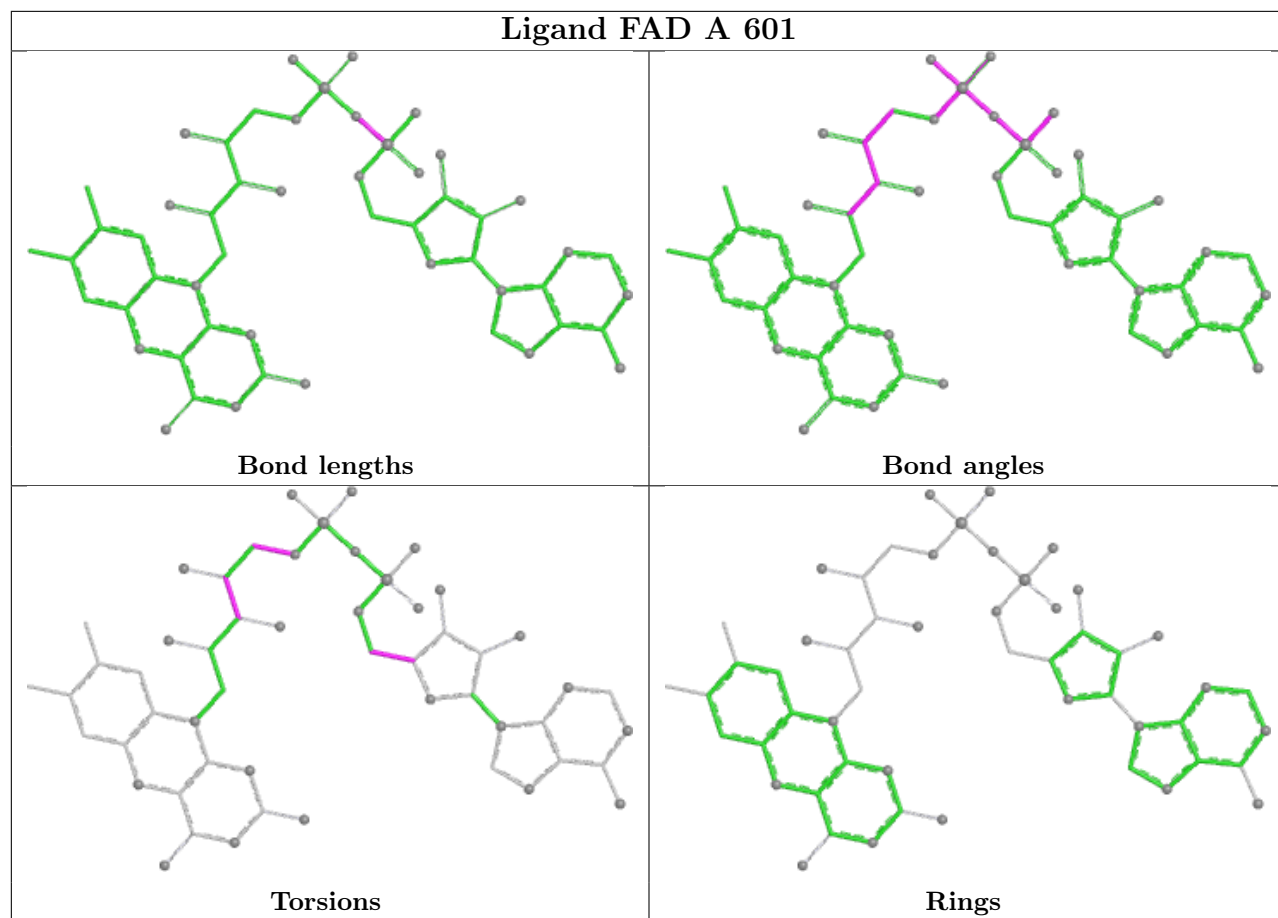
Mol	Chain	Res	Type	Atoms
2	C	601	FAD	PA-O3P-P-O5'
2	A	601	FAD	O3'-C3'-C4'-C5'
2	B	601	FAD	O3'-C3'-C4'-C5'
2	D	601	FAD	O3'-C3'-C4'-C5'
2	A	601	FAD	C2'-C3'-C4'-C5'
3	A	602	PEG	O2-C3-C4-O4
2	B	601	FAD	C2'-C3'-C4'-C5'
3	A	602	PEG	C1-C2-O2-C3
2	A	601	FAD	O3'-C3'-C4'-O4'
2	D	601	FAD	C2'-C3'-C4'-C5'
2	B	601	FAD	O3'-C3'-C4'-O4'
3	A	602	PEG	O1-C1-C2-O2
2	D	601	FAD	C4'-C5'-O5'-P
2	B	601	FAD	C4'-C5'-O5'-P
3	A	602	PEG	C4-C3-O2-C2
2	A	601	FAD	O4B-C4B-C5B-O5B
2	A	601	FAD	C4'-C5'-O5'-P
2	A	601	FAD	C3B-C4B-C5B-O5B

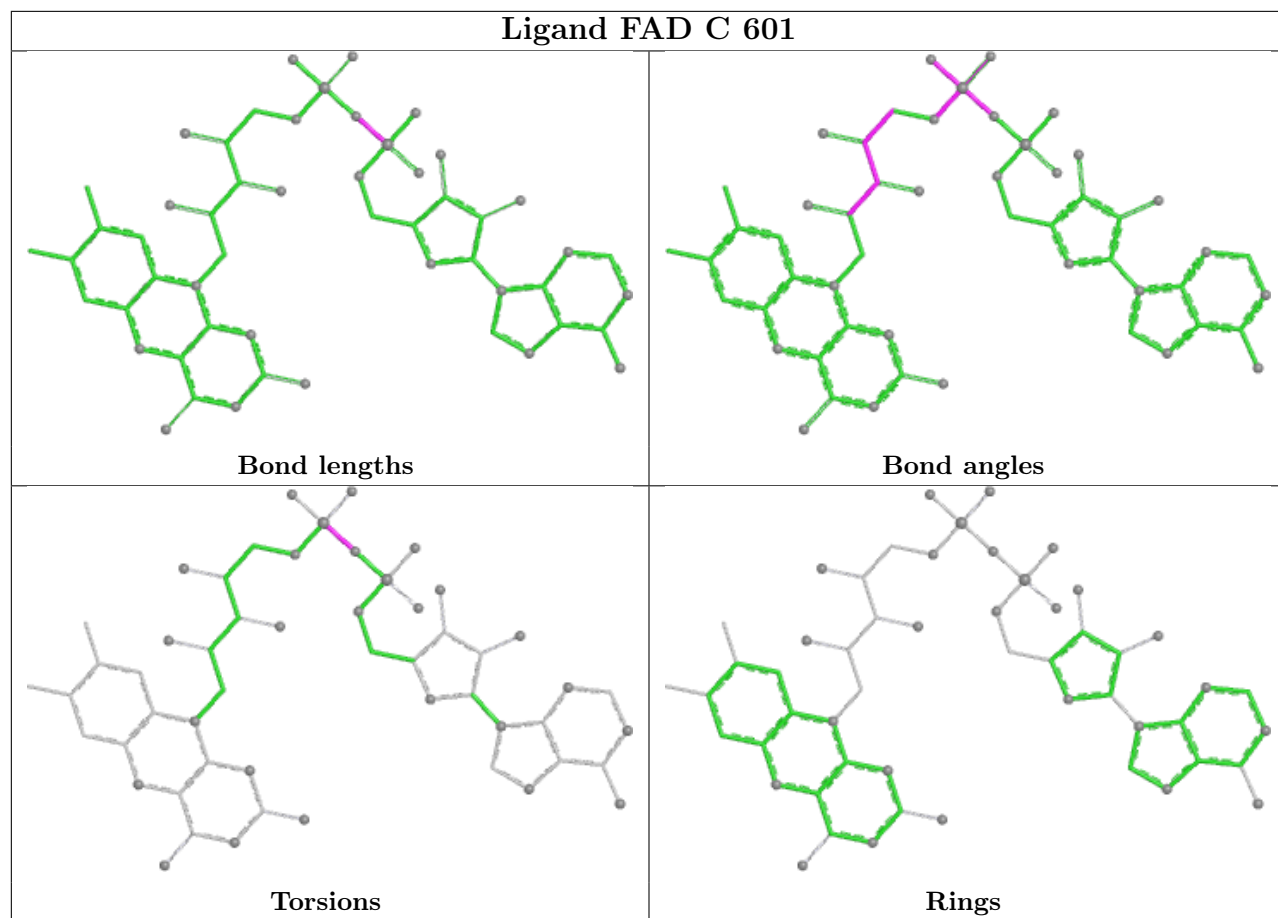
There are no ring outliers.

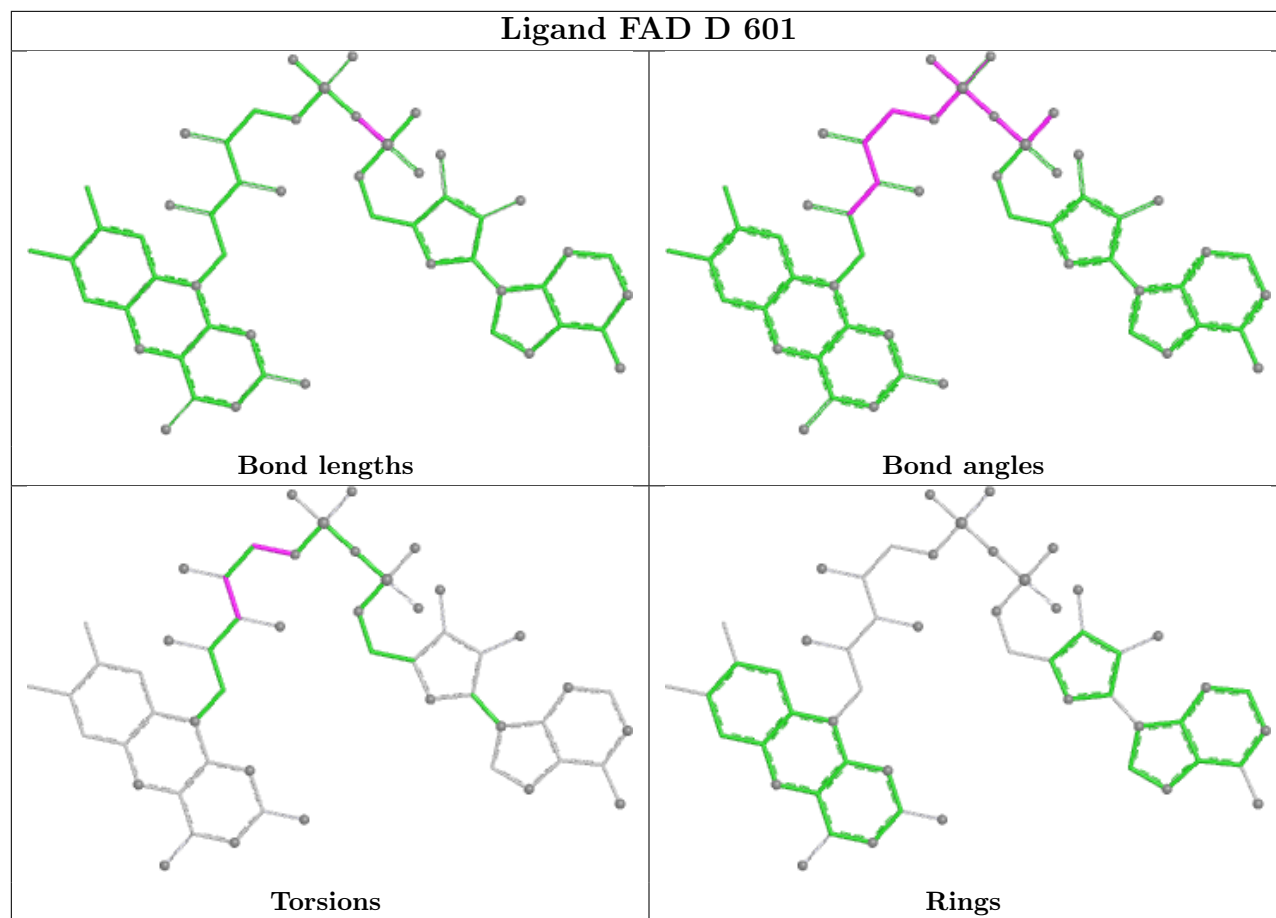
1 monomer is involved in 4 short contacts:

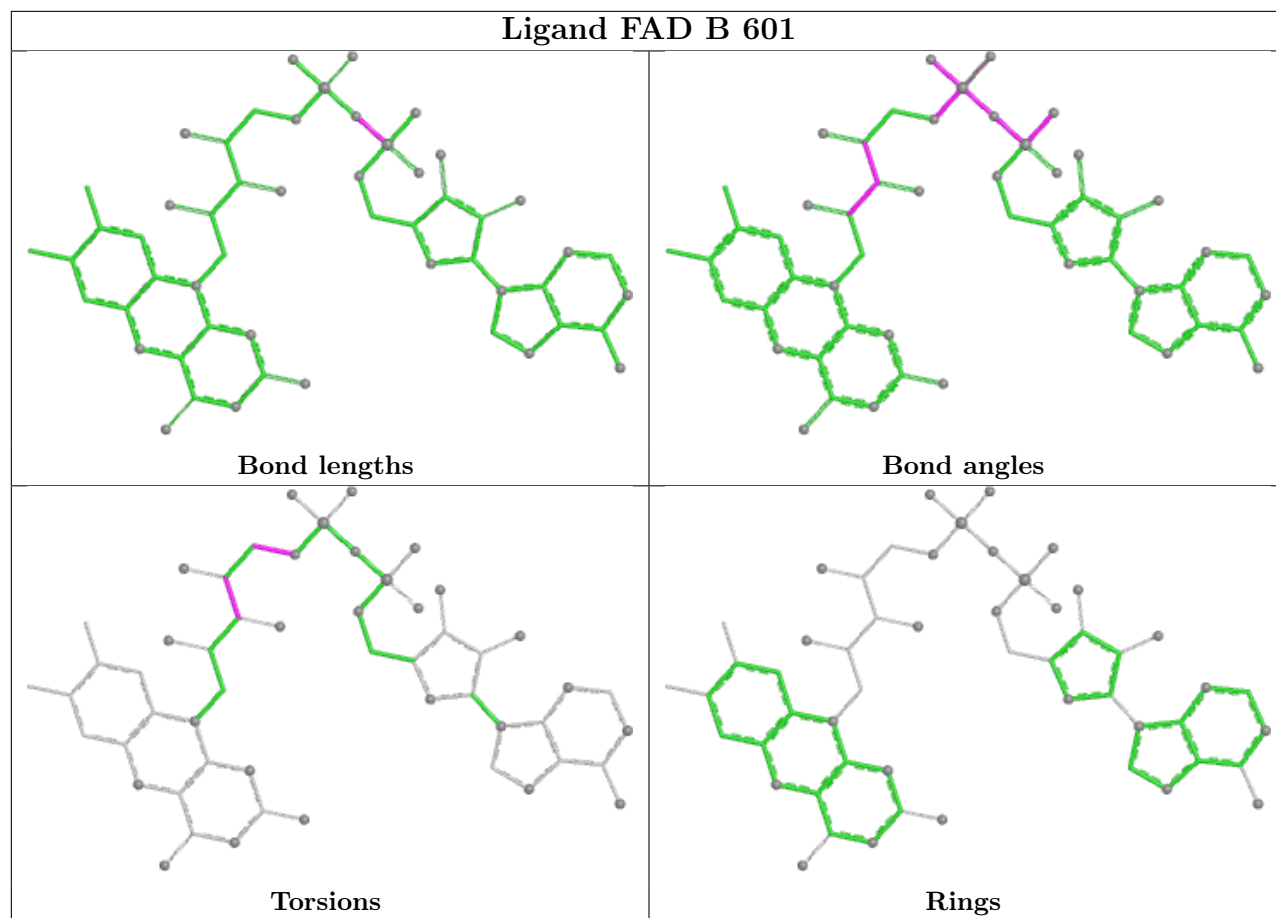
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	602	PEG	4	0

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	488/501 (97%)	-0.08	16 (3%)	49 45	21, 43, 85, 113	0
1	B	487/501 (97%)	0.03	25 (5%)	33 29	21, 44, 92, 118	0
1	C	488/501 (97%)	0.06	16 (3%)	49 45	27, 49, 88, 112	0
1	D	479/501 (95%)	0.36	31 (6%)	25 22	27, 56, 105, 123	0
All	All	1942/2004 (96%)	0.09	88 (4%)	38 34	21, 48, 92, 123	0

All (88) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	122	ALA	5.3
1	D	293	VAL	5.2
1	B	398	SER	5.0
1	D	279	ILE	4.5
1	D	445	ILE	4.3
1	B	279	ILE	4.0
1	C	120	ARG	4.0
1	A	122	ALA	4.0
1	B	396	LYS	3.9
1	C	399	ASN	3.9
1	C	30	MET	3.8
1	B	293	VAL	3.7
1	C	123	ARG	3.7
1	C	516	PRO	3.7
1	B	122	ALA	3.7
1	B	397	ASP	3.6
1	C	398	SER	3.6
1	D	288	LYS	3.6
1	C	279	ILE	3.5
1	C	400	ASN	3.5
1	D	398	SER	3.4

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Mol	Chain	Res	Type	RSRZ
1	D	271	PHE	3.4
1	A	516	PRO	3.4
1	D	294	GLN	3.4
1	B	292	PRO	3.3
1	B	515	LEU	3.3
1	D	400	ASN	3.2
1	D	396	LYS	3.2
1	B	400	ASN	3.1
1	B	297	HIS	3.1
1	A	293	VAL	3.0
1	A	30	MET	3.0
1	D	290	GLN	3.0
1	D	260	MET	3.0
1	A	279	ILE	3.0
1	A	513	ALA	3.0
1	D	292	PRO	2.9
1	C	298	PHE	2.9
1	B	291	ASP	2.9
1	D	397	ASP	2.9
1	D	439	LEU	2.8
1	B	288	LYS	2.8
1	D	399	ASN	2.8
1	B	399	ASN	2.7
1	D	442	HIS	2.7
1	A	399	ASN	2.7
1	C	513	ALA	2.7
1	B	298	PHE	2.7
1	D	30	MET	2.6
1	D	513	ALA	2.6
1	A	47	ASP	2.6
1	A	123	ARG	2.6
1	C	442	HIS	2.5
1	D	302	GLY	2.5
1	D	277	THR	2.5
1	D	120	ARG	2.5
1	B	123	ARG	2.5
1	A	292	PRO	2.5
1	C	460	GLU	2.4
1	B	121	GLY	2.4
1	B	290	GLN	2.4
1	D	122	ALA	2.4
1	D	291	ASP	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	271	PHE	2.4
1	D	248	PHE	2.4
1	A	297	HIS	2.4
1	D	443	GLY	2.3
1	D	303	ARG	2.3
1	A	285	ALA	2.2
1	A	400	ASN	2.2
1	C	515	LEU	2.2
1	D	276	ARG	2.2
1	A	514	SER	2.2
1	B	513	ALA	2.2
1	D	306	TYR	2.2
1	C	288	LYS	2.1
1	B	300	SER	2.1
1	B	295	ALA	2.1
1	C	31	LEU	2.1
1	B	246	SER	2.1
1	D	395	LEU	2.1
1	A	121	GLY	2.1
1	D	289	PRO	2.0
1	B	442	HIS	2.0
1	D	287	PHE	2.0
1	B	30	MET	2.0
1	A	31	LEU	2.0
1	B	395	LEU	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 ⓘ

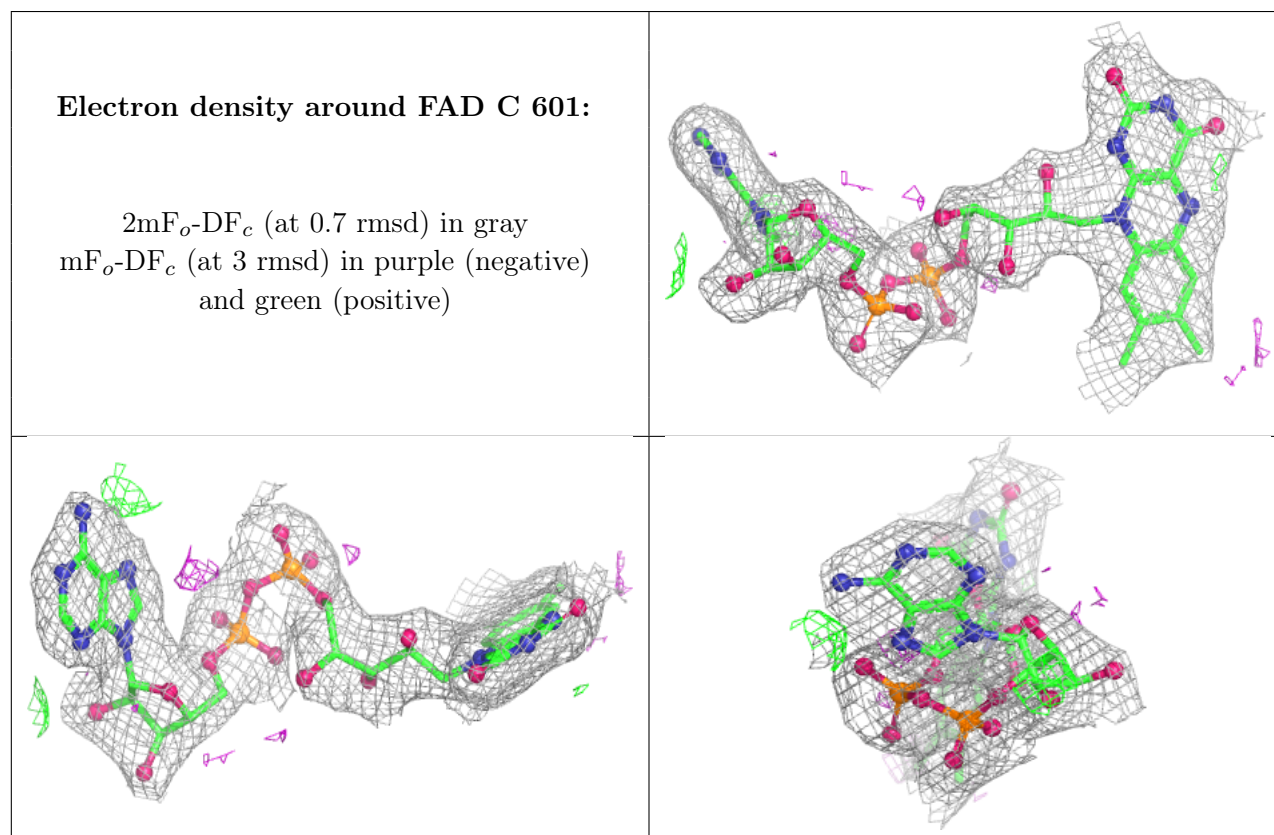
There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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.

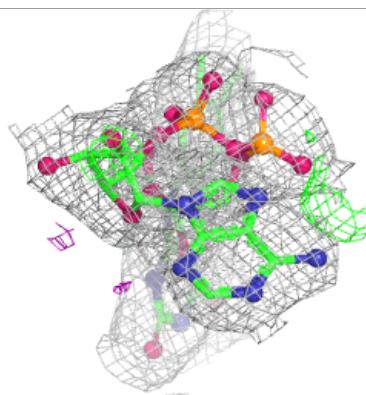
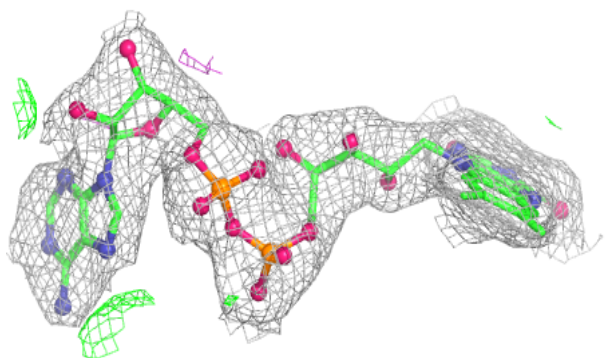
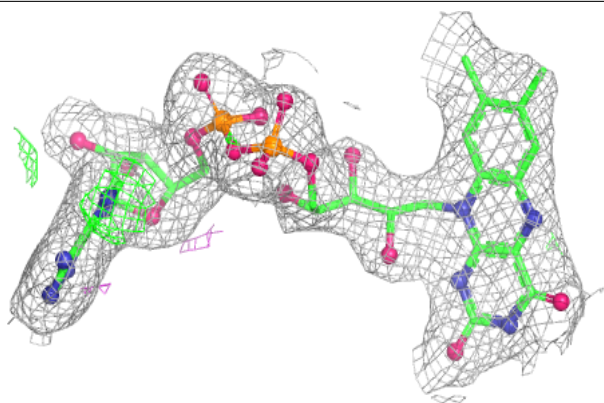
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	PEG	A	602	7/7	0.90	0.16	58,64,66,67	0
2	FAD	C	601	53/53	0.96	0.07	32,37,42,43	0
2	FAD	D	601	53/53	0.97	0.07	37,41,48,49	0
2	FAD	A	601	53/53	0.97	0.06	19,32,39,41	0
2	FAD	B	601	53/53	0.98	0.06	22,33,38,40	0

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.

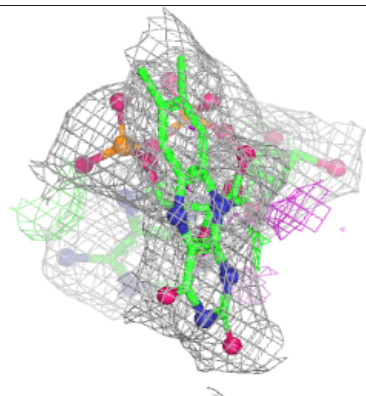
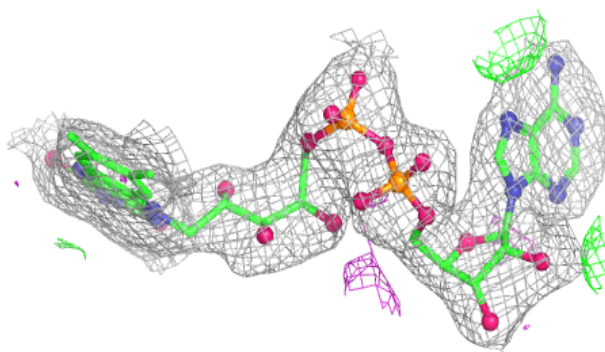
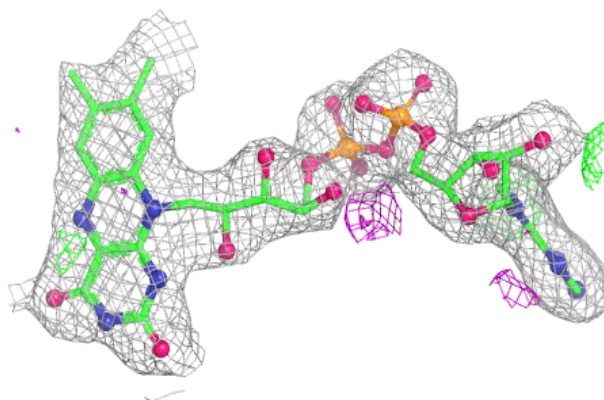


Electron density around FAD D 601:

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

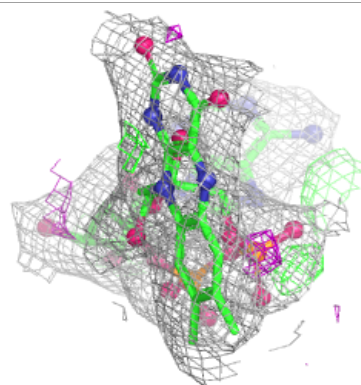
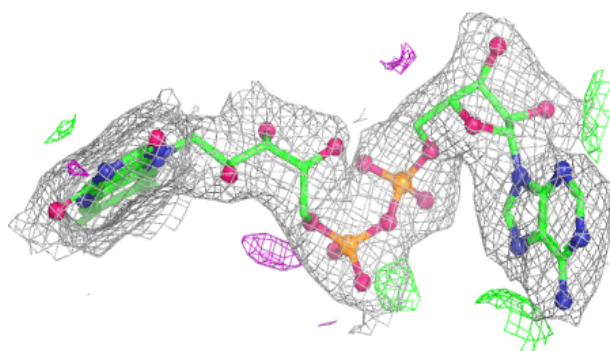
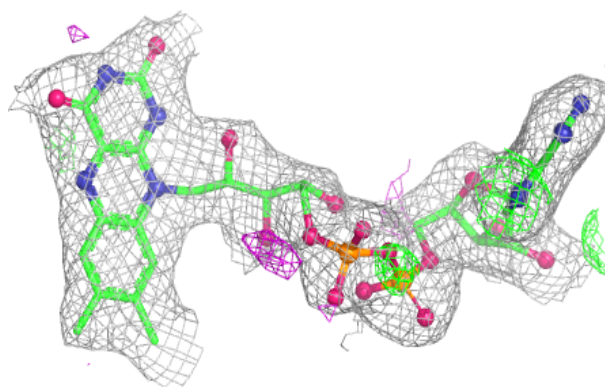
**Electron density around FAD A 601:**

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



Electron density around FAD B 601:

$2mF_o-DF_c$ (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.