



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 17, 2025 – 12:14 PM JST

PDB ID : 9U3D / pdb_00009u3d
Title : Monomeric sarcosine oxidase from Bacillus sp. (SoxB) complexed with D-Proline
Authors : Zhang, Y.; Nakajima, Y.; Kurobe, M.; Nakamura, T.; Himiyama, T.; Nishiya, Y.
Deposited on : 2025-03-18
Resolution : 1.35 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 Phenix2.0rc1
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	2.0rc1
EDS	:	3.0
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.006 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.45.1

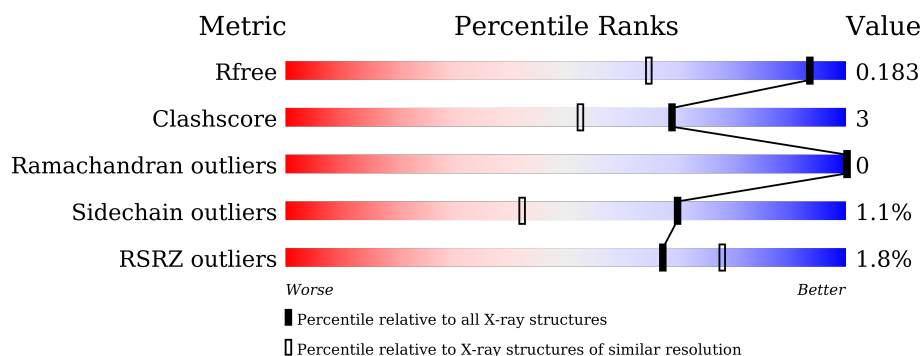
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.35 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}	164625	1089 (1.36-1.36)
Clashscore	180529	1157 (1.36-1.36)
Ramachandran outliers	177936	1146 (1.36-1.36)
Sidechain outliers	177891	1146 (1.36-1.36)
RSRZ outliers	164620	1088 (1.36-1.36)

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	397	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 1%, green 98%, grey 2%);"></div> <div style="display: flex; justify-content: space-between; margin-top: 2px;"> % 90% 5% . . </div> </div>
1	B	397	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 3%, orange 1%, yellow 1%, green 95%, grey 2%);"></div> <div style="display: flex; justify-content: space-between; margin-top: 2px;"> 3% 87% 7% . 5% </div> </div>

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	DPR	A	402	-	-	X	-
5	GOL	A	405	-	-	X	-

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 6764 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.

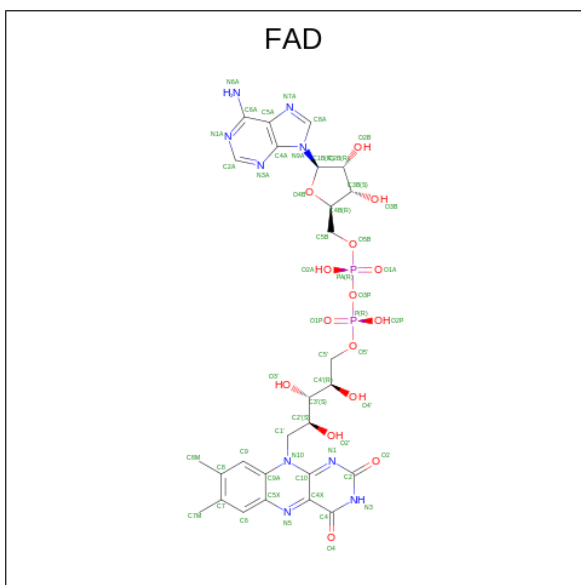
- Molecule 1 is a protein called Monomeric sarcosine oxidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	381	Total	C	N	O	S	0	21	0
			3095	1984	512	589	10			
1	B	378	Total	C	N	O	S	0	16	0
			3043	1949	507	577	10			

There are 16 discrepancies between the modelled and reference sequences:

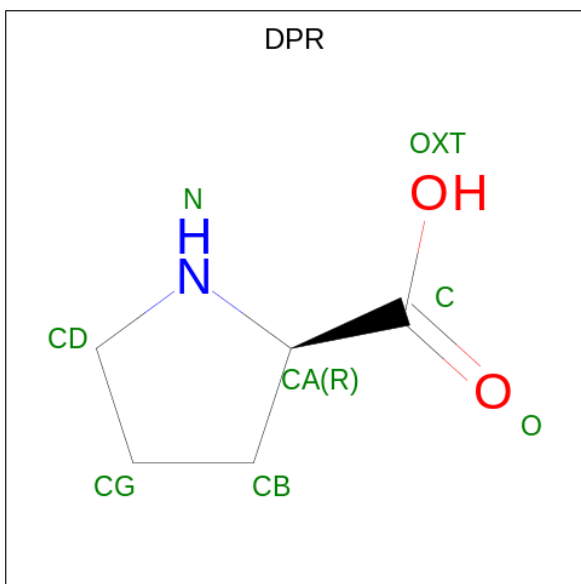
Chain	Residue	Modelled	Actual	Comment	Reference
A	390	GLY	-	expression tag	UNP P40859
A	391	SER	-	expression tag	UNP P40859
A	392	HIS	-	expression tag	UNP P40859
A	393	HIS	-	expression tag	UNP P40859
A	394	HIS	-	expression tag	UNP P40859
A	395	HIS	-	expression tag	UNP P40859
A	396	HIS	-	expression tag	UNP P40859
A	397	HIS	-	expression tag	UNP P40859
B	390	GLY	-	expression tag	UNP P40859
B	391	SER	-	expression tag	UNP P40859
B	392	HIS	-	expression tag	UNP P40859
B	393	HIS	-	expression tag	UNP P40859
B	394	HIS	-	expression tag	UNP P40859
B	395	HIS	-	expression tag	UNP P40859
B	396	HIS	-	expression tag	UNP P40859
B	397	HIS	-	expression tag	UNP P40859

- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (CCD ID: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$) (labeled as "Ligand of Interest" by depositor).



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 D-PROLINE (CCD ID: DPR) (formula: $\text{C}_5\text{H}_9\text{NO}_2$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			8	5	1	2		

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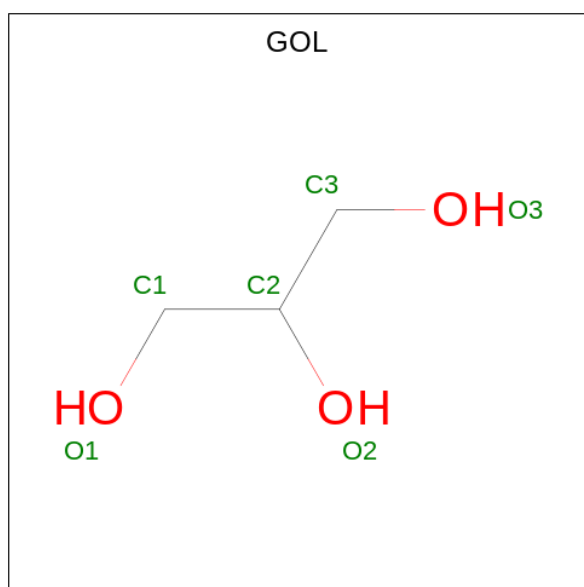
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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	1	Total	C	N	O	0	0
			8	5	1	2		

- Molecule 4 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

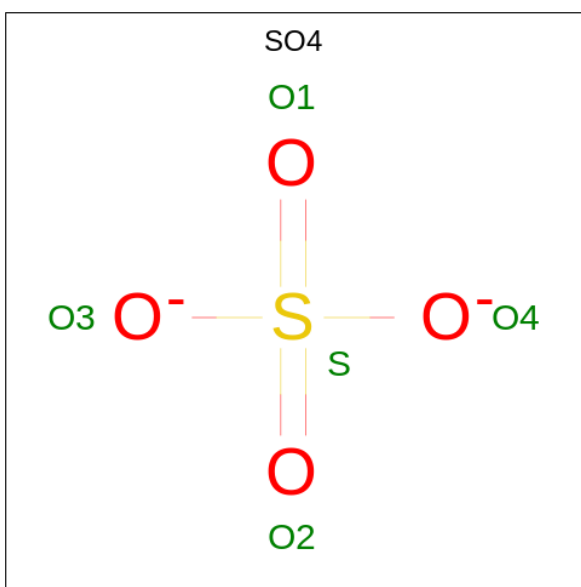
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Cl	0	0
			1	1		
4	B	1	Total	Cl	0	0
			1	1		

- Molecule 5 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			6	3	3		
5	A	1	Total	C	O	0	0
			6	3	3		
5	A	1	Total	C	O	0	0
			6	3	3		
5	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 6 is SULFATE ION (CCD ID: SO4) (formula: O₄S).

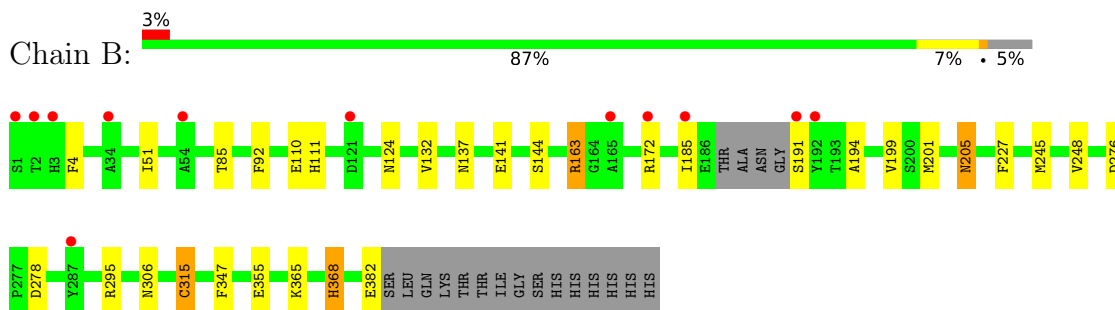


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	O	S	0	0
			5	4	1		
6	B	1	Total	O	S	0	0
			5	4	1		
6	B	1	Total	O	S	0	0
			5	4	1		
6	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	245	Total	O	0	0
			245	245		
7	B	213	Total	O	0	0
			213	213		

- Molecule 1: Monomeric sarcosine oxidase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	71.37Å 69.69Å 73.09Å 90.00° 92.10° 90.00°	Depositor
Resolution (Å)	41.67 – 1.35 41.67 – 1.35	Depositor EDS
% Data completeness (in resolution range)	96.8 (41.67-1.35) 96.8 (41.67-1.35)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.65 (at 1.35Å)	Xtriage
Refinement program	REFMAC 5.8.0425	Depositor
R, R_{free}	0.158 , 0.183 0.158 , 0.183	Depositor DCC
R_{free} test set	8068 reflections (5.14%)	wwPDB-VP
Wilson B-factor (Å ²)	16.7	Xtriage
Anisotropy	0.045	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 31.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.000 for l,k,-h 0.019 for h,-k,-l 0.014 for l,-k,h	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	6764	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.27% 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: SO4, CL, FAD, GOL, DPR

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	0.75	0/3237	1.12	6/4378 (0.1%)
1	B	0.73	0/3165	1.13	8/4280 (0.2%)
All	All	0.74	0/6402	1.12	14/8658 (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	2
1	B	0	2
All	All	0	4

There are no bond length outliers.

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	170	HIS	CB-CG-CD2	9.73	143.85	131.20
1	A	170	HIS	CB-CG-ND1	-8.80	109.51	122.70
1	B	315	CYS	N-CA-CB	-7.56	99.37	111.62
1	A	92	PHE	CA-CB-CG	6.67	120.47	113.80
1	A	328	ASP	CA-CB-CG	6.38	118.98	112.60
1	B	276	ASP	CA-CB-CG	5.97	118.57	112.60
1	B	227	PHE	CA-CB-CG	5.52	119.32	113.80
1	B	368	HIS	CB-CG-CD2	-5.52	124.03	131.20
1	A	322	ASP	CA-CB-CG	5.49	118.09	112.60
1	A	52	ARG	NE-CZ-NH2	5.39	124.05	119.20
1	B	205	ASN	CA-CB-CG	5.34	117.94	112.60
1	B	347	PHE	CA-CB-CG	-5.16	108.64	113.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	248	VAL	O-C-N	-5.14	117.13	121.40
1	B	92	PHE	CA-CB-CG	5.14	118.94	113.80

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	163[A]	ARG	Sidechain
1	A	163[B]	ARG	Sidechain
1	B	163[A]	ARG	Sidechain
1	B	163[B]	ARG	Sidechain

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	3095	0	3060	21	0
1	B	3043	0	3002	17	0
2	A	53	0	30	2	0
2	B	53	0	30	3	0
3	A	8	0	8	6	0
3	B	8	0	8	3	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	18	0	24	10	0
5	B	6	0	8	0	0
6	A	5	0	0	0	0
6	B	15	0	0	0	0
7	A	245	0	0	1	0
7	B	213	0	0	4	0
All	All	6764	0	6170	42	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 3.

All (42) 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:172[B]:ARG:HD2	5:A:405:GOL:O2	1.36	1.26
1:A:172[B]:ARG:CD	5:A:405:GOL:O2	2.17	0.93
1:B:124[B]:ASN:OD1	7:B:501:HOH:O	1.93	0.86
2:B:401:FAD:N5	3:B:402:DPR:HG2	1.92	0.84
2:A:401:FAD:H61A	5:A:405:GOL:H31	1.48	0.79
1:A:172[B]:ARG:HD2	5:A:405:GOL:HO2	1.48	0.76
1:B:124[A]:ASN:HD21	1:B:132:VAL:H	1.33	0.76
1:A:124:ASN:HD21	1:A:132:VAL:H	1.34	0.74
1:A:20:TYR:CZ	1:A:24:LYS:HD2	2.23	0.74
1:A:204:TRP:CD1	5:A:405:GOL:H32	2.23	0.74
2:A:401:FAD:N6A	5:A:405:GOL:H31	2.06	0.70
1:A:60:GLU:OE1	1:A:372:ILE:HG21	1.94	0.66
2:B:401:FAD:C4X	3:B:402:DPR:HG2	2.26	0.66
1:A:254:TYR:CD1	3:A:402:DPR:HG3	2.31	0.64
1:A:208:LEU:HD11	5:A:405:GOL:H2	1.81	0.61
1:A:163[B]:ARG:NH1	7:A:504:HOH:O	2.36	0.58
1:B:368:HIS:HD2	7:B:691:HOH:O	1.88	0.57
1:B:355:GLU:OE2	1:B:368:HIS:HE1	1.89	0.56
1:A:269:HIS:ND1	3:A:402:DPR:HD2	2.23	0.53
1:B:141:GLU:OE1	1:B:144[B]:SER:OG	2.18	0.53
1:A:172[B]:ARG:CG	5:A:405:GOL:O2	2.58	0.51
1:B:85[B]:THR:HG22	7:B:589:HOH:O	2.08	0.51
1:A:204:TRP:CD1	5:A:405:GOL:H11	2.46	0.51
1:A:172[A]:ARG:NH1	5:A:405:GOL:O3	2.45	0.50
1:B:51:ILE:O	1:B:144[A]:SER:HB2	2.11	0.50
1:B:245:MET:SD	3:B:402:DPR:HA	2.52	0.49
1:A:269:HIS:CE1	3:A:402:DPR:HD2	2.48	0.49
1:B:295:ARG:HH11	1:B:306:ASN:ND2	2.11	0.48
1:B:185:ILE:O	1:B:191:SER:HB2	2.14	0.48
1:A:275:ILE:HD12	1:A:275:ILE:C	2.39	0.47
1:B:199:VAL:CG1	1:B:201:MET:HE2	2.45	0.47
1:A:295:ARG:HH11	1:A:306:ASN:ND2	2.14	0.46
1:A:20:TYR:CZ	1:A:24:LYS:CD	2.97	0.46
1:B:172:ARG:NH1	1:B:278:ASP:OD1	2.49	0.45
1:A:245:MET:SD	3:A:402:DPR:HA	2.56	0.45
1:B:111:HIS:HD2	7:B:697:HOH:O	2.02	0.43
1:B:4:PHE:O	1:B:194:ALA:HA	2.20	0.42
1:A:52:ARG:HH21	3:A:402:DPR:C	2.33	0.42
1:B:315:CYS:HB3	2:B:401:FAD:HM71	2.02	0.41
1:B:141:GLU:CD	1:B:144[B]:SER:OG	2.63	0.41
1:B:201:MET:H	1:B:205:ASN:HD21	1.69	0.41
1:A:254:TYR:CG	3:A:402:DPR:HG3	2.57	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	401/397 (101%)	394 (98%)	7 (2%)	0	100	100
1	B	390/397 (98%)	382 (98%)	8 (2%)	0	100	100
All	All	791/794 (100%)	776 (98%)	15 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	340/333 (102%)	337 (99%)	3 (1%)	75	51
1	B	333/333 (100%)	327 (98%)	6 (2%)	54	23
All	All	673/666 (101%)	664 (99%)	9 (1%)	70	35

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	137	ASN
1	A	365[A]	LYS
1	A	365[B]	LYS
1	B	110	GLU
1	B	137	ASN

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Mol	Chain	Res	Type
1	B	163[A]	ARG
1	B	163[B]	ARG
1	B	365	LYS
1	B	382	GLU

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	3	HIS
1	A	21	GLN
1	A	25	GLN
1	A	41	ASN
1	A	124	ASN
1	A	143	ASN
1	A	189	ASN
1	A	205	ASN
1	A	293	ASN
1	A	306	ASN
1	B	21	GLN
1	B	41	ASN
1	B	111	HIS
1	B	137	ASN
1	B	170	HIS
1	B	205	ASN
1	B	293	ASN
1	B	306	ASN
1	B	368	HIS

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

Of 14 ligands modelled in this entry, 2 are monoatomic - leaving 12 for Mogul analysis.

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
5	GOL	A	404	-	5,5,5	0.17	0	5,5,5	0.43	0
5	GOL	B	404	-	5,5,5	0.13	0	5,5,5	0.35	0
3	DPR	B	402	-	8,8,8	0.63	0	10,10,10	1.69	3 (30%)
5	GOL	A	406	-	5,5,5	0.19	0	5,5,5	0.42	0
5	GOL	A	405	-	5,5,5	0.21	0	5,5,5	0.89	0
2	FAD	A	401	1	53,58,58	0.92	4 (7%)	68,89,89	0.90	3 (4%)
2	FAD	B	401	1	53,58,58	0.90	1 (1%)	68,89,89	0.85	3 (4%)
6	SO4	A	407	-	4,4,4	0.33	0	6,6,6	0.08	0
3	DPR	A	402	-	8,8,8	0.47	0	10,10,10	1.61	2 (20%)
6	SO4	B	407	-	4,4,4	0.24	0	6,6,6	0.29	0
6	SO4	B	405	-	4,4,4	0.37	0	6,6,6	0.08	0
6	SO4	B	406	-	4,4,4	0.36	0	6,6,6	0.09	0

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
5	GOL	A	404	-	-	2/4/4/4	-
5	GOL	B	404	-	-	3/4/4/4	-
3	DPR	B	402	-	-	0/4/11/11	0/1/1/1
5	GOL	A	406	-	-	2/4/4/4	-
5	GOL	A	405	-	-	2/4/4/4	-
2	FAD	A	401	1	-	1/30/50/50	0/6/6/6
2	FAD	B	401	1	-	1/30/50/50	0/6/6/6
3	DPR	A	402	-	-	0/4/11/11	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	401	FAD	C4X-N5	3.07	1.36	1.30
2	A	401	FAD	C4X-N5	2.72	1.36	1.30
2	A	401	FAD	O4B-C1B	2.40	1.44	1.41
2	A	401	FAD	P-O2P	-2.03	1.45	1.55
2	A	401	FAD	C4A-N3A	-2.00	1.32	1.35

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	402	DPR	O-C-CA	-3.48	110.84	122.26
3	A	402	DPR	O-C-CA	-3.38	111.16	122.26
2	A	401	FAD	O4B-C1B-C2B	-2.54	103.22	106.93
3	B	402	DPR	OXT-C-O	2.52	129.80	124.09
2	B	401	FAD	C4A-C5A-N7A	2.50	112.01	109.40
3	A	402	DPR	OXT-C-O	2.47	129.69	124.09
2	A	401	FAD	C5A-C6A-N6A	2.38	123.97	120.35
2	B	401	FAD	O4B-C1B-C2B	-2.25	103.64	106.93
3	B	402	DPR	CB-CA-C	-2.12	107.22	112.75
2	A	401	FAD	O2'-C2'-C3'	2.04	114.05	109.10
2	B	401	FAD	O2P-P-O1P	2.02	122.24	112.24

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	404	GOL	C1-C2-C3-O3
5	A	404	GOL	O2-C2-C3-O3
5	A	405	GOL	O1-C1-C2-O2
5	A	405	GOL	O1-C1-C2-C3
5	A	406	GOL	O1-C1-C2-C3
5	B	404	GOL	O1-C1-C2-C3
5	B	404	GOL	O1-C1-C2-O2
5	A	406	GOL	C1-C2-C3-O3
5	B	404	GOL	C1-C2-C3-O3
2	A	401	FAD	O4B-C4B-C5B-O5B
2	B	401	FAD	O4B-C4B-C5B-O5B

There are no ring outliers.

5 monomers are involved in 20 short contacts:

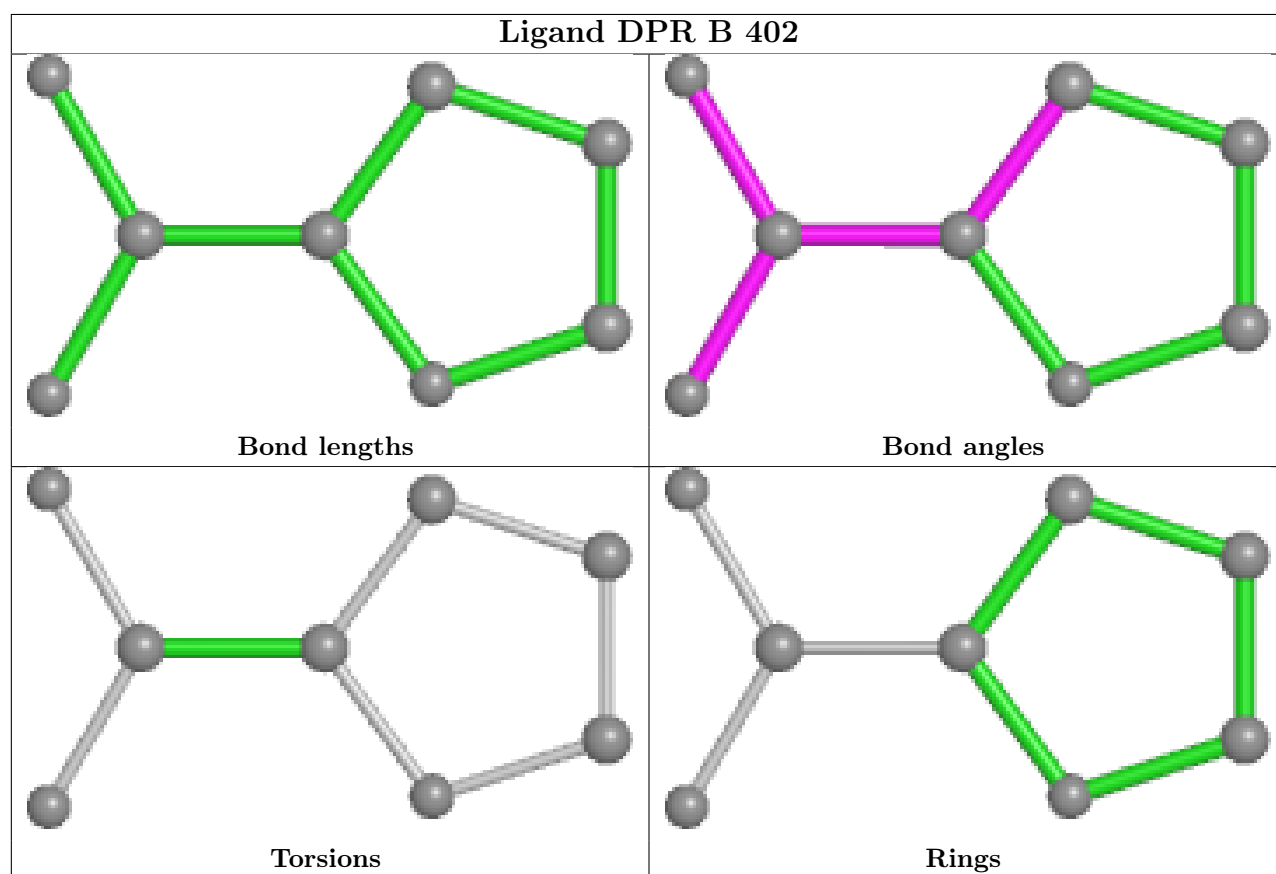
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	402	DPR	3	0

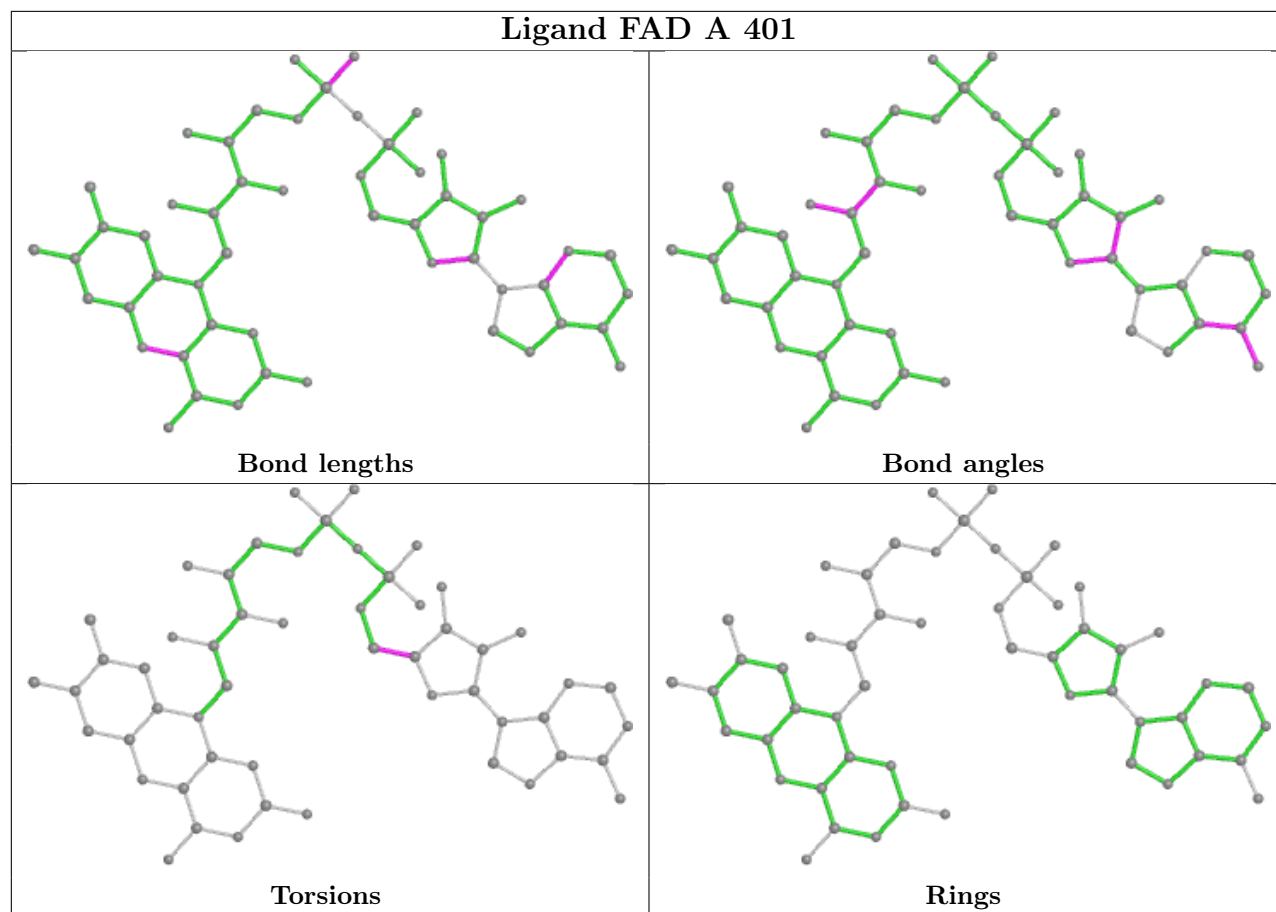
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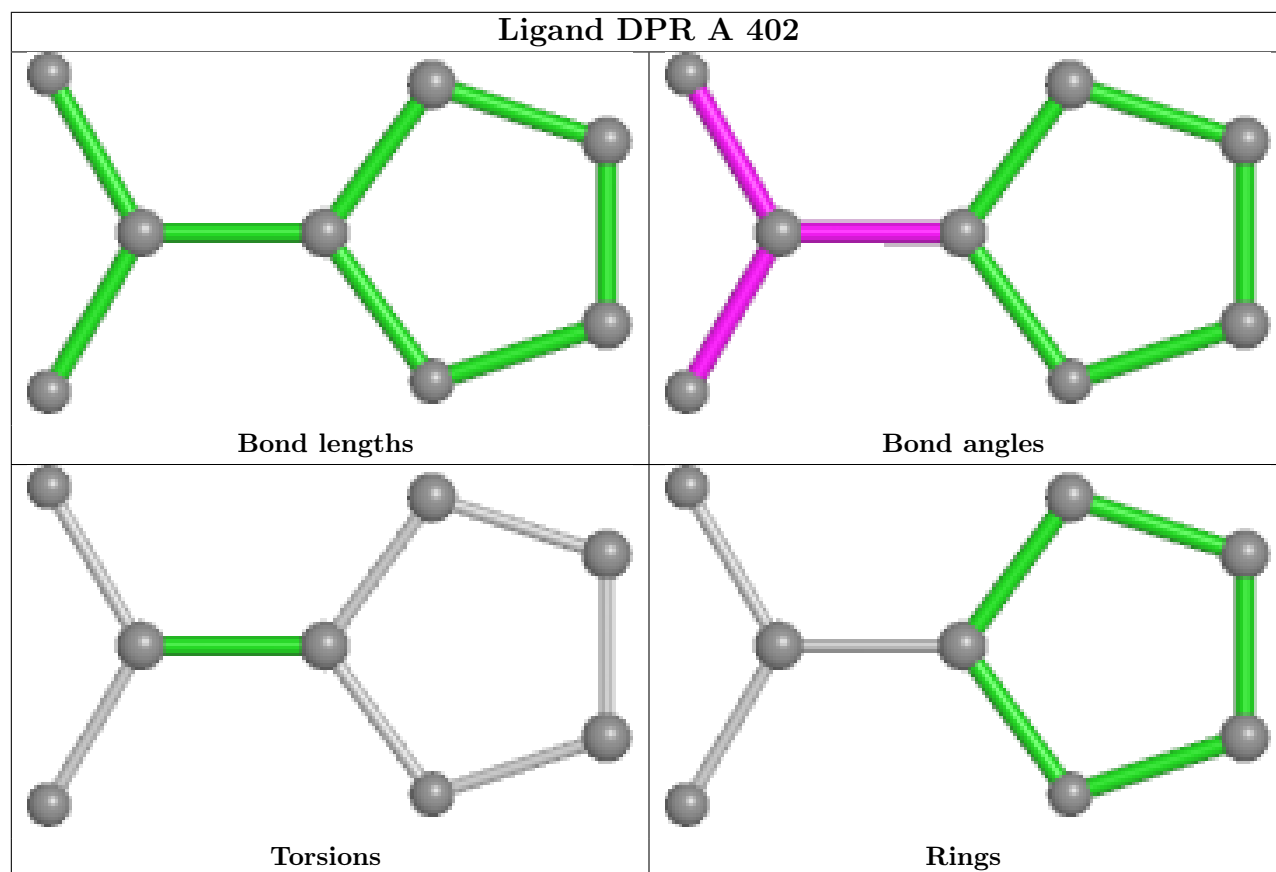
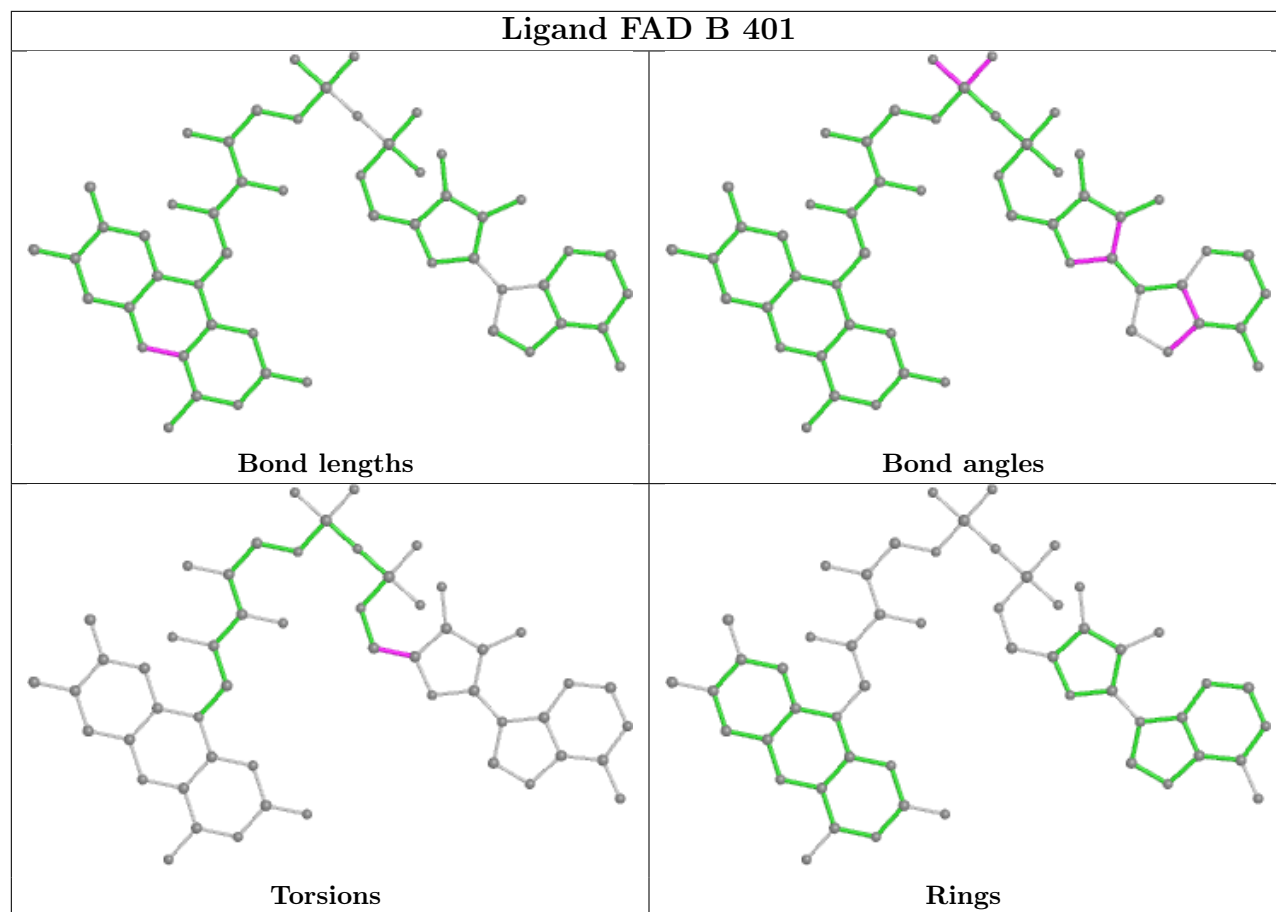
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	405	GOL	10	0
2	A	401	FAD	2	0
2	B	401	FAD	3	0
3	A	402	DPR	6	0

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	381/397 (95%)	0.09	2 (0%) 87 93	8, 17, 29, 47	21 (5%)
1	B	378/397 (95%)	0.36	12 (3%) 50 61	8, 18, 34, 50	16 (4%)
All	All	759/794 (95%)	0.22	14 (1%) 67 78	8, 17, 32, 50	37 (4%)

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	192	TYR	3.9
1	B	1	SER	3.8
1	B	191	SER	3.0
1	B	185	ILE	2.8
1	B	2	THR	2.7
1	B	172	ARG	2.6
1	B	3	HIS	2.6
1	B	54	ALA	2.5
1	B	121	ASP	2.3
1	A	2	THR	2.3
1	A	3	HIS	2.2
1	B	287	TYR	2.1
1	B	34	ALA	2.0
1	B	165	ALA	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

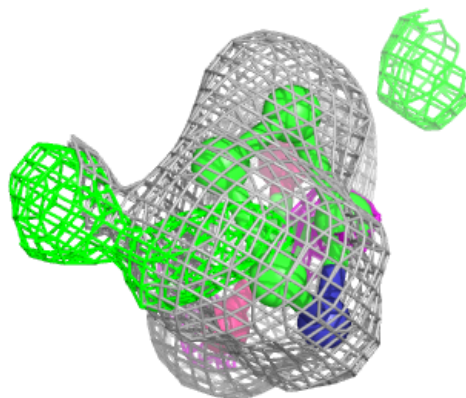
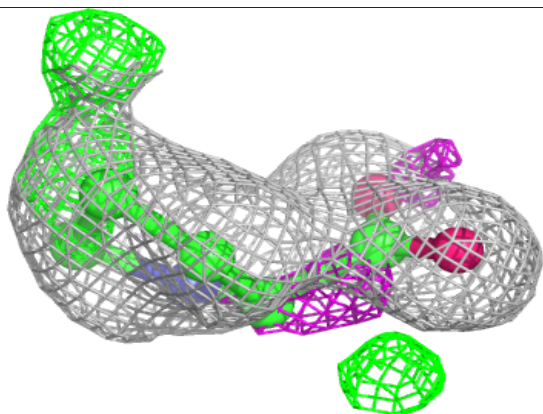
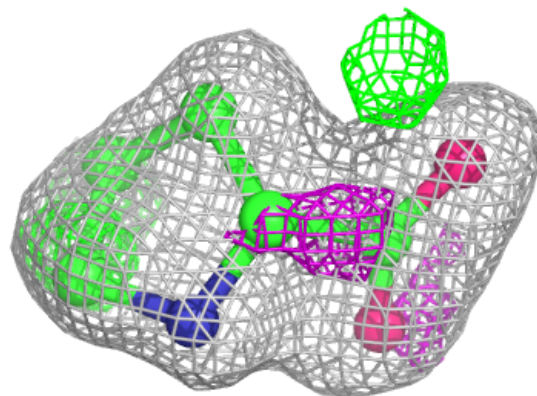
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
6	SO4	B	405	5/5	0.66	0.13	71,78,78,83	0
6	SO4	A	407	5/5	0.68	0.14	82,86,91,100	0
6	SO4	B	406	5/5	0.78	0.15	102,105,110,113	0
5	GOL	A	406	6/6	0.82	0.13	32,34,36,41	0
5	GOL	A	404	6/6	0.83	0.16	29,37,51,65	0
5	GOL	B	404	6/6	0.84	0.15	33,39,54,62	0
6	SO4	B	407	5/5	0.84	0.12	35,43,58,63	0
5	GOL	A	405	6/6	0.85	0.12	30,36,37,48	0
3	DPR	B	402	8/8	0.86	0.16	22,36,38,40	0
3	DPR	A	402	8/8	0.90	0.11	23,31,32,34	0
2	FAD	B	401	53/53	0.98	0.05	12,14,24,27	0
2	FAD	A	401	53/53	0.99	0.04	10,12,14,17	0
4	CL	B	403	1/1	0.99	0.02	14,14,14,14	0
4	CL	A	403	1/1	1.00	0.02	13,13,13,13	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.

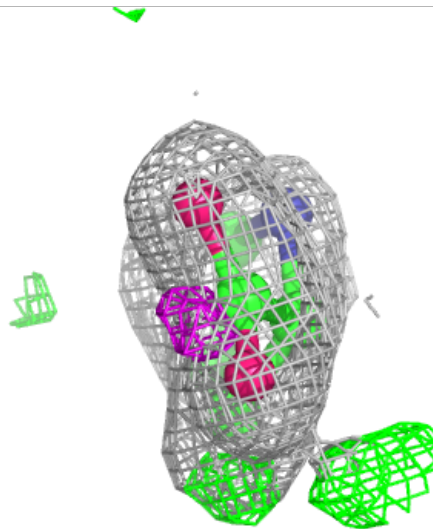
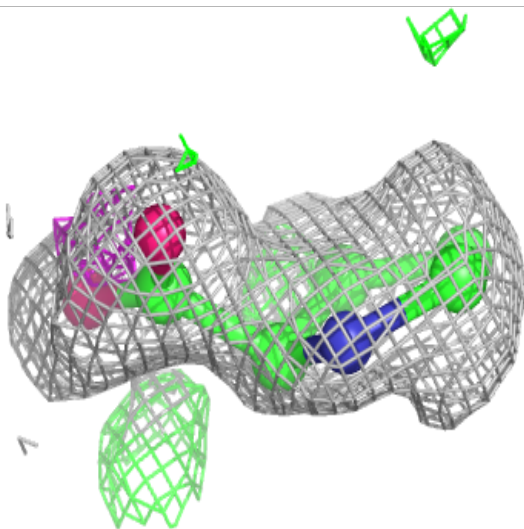
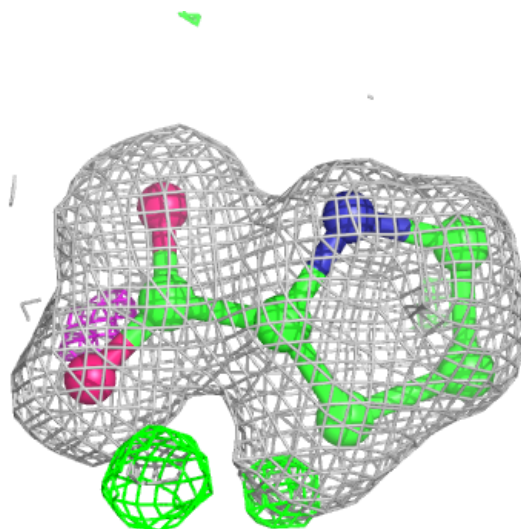
Electron density around DPR B 402:

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



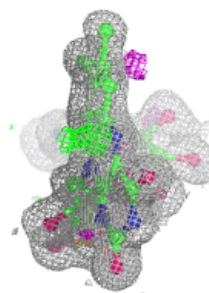
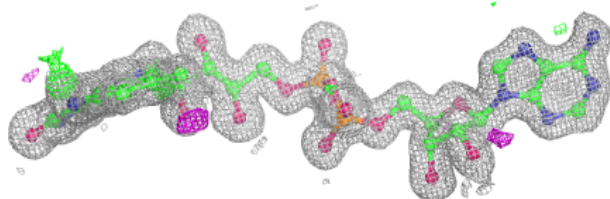
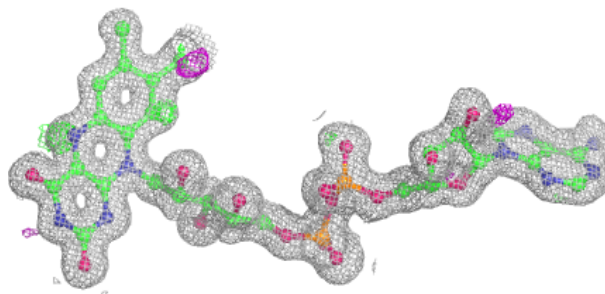
Electron density around DPR A 402:

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

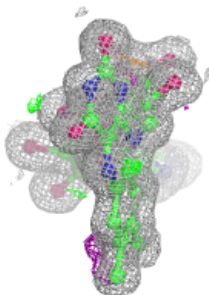
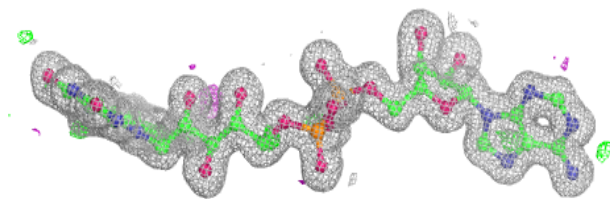
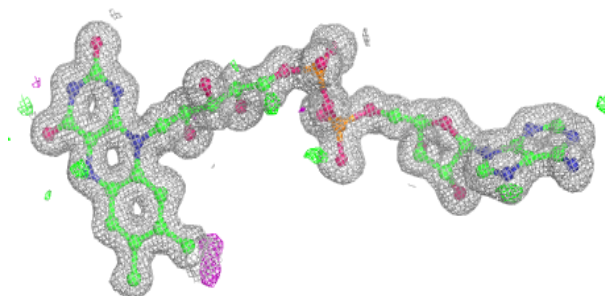


Electron density around FAD B 401:

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

**Electron density around FAD A 401:**

$2mF_o - DF_c$ (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.