



wwPDB X-ray Structure Validation Summary Report ⓘ

May 22, 2020 – 06:54 am BST

PDB ID : 1OFE
Title : Glutamate Synthase from Synechocystis sp in complex with 2-Oxoglutarate and L-DON at 2.45 Angstrom resolution
Authors : Van Den heuvel, R.H.H.; Mattevi, A.
Deposited on : 2003-04-14
Resolution : 2.45 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

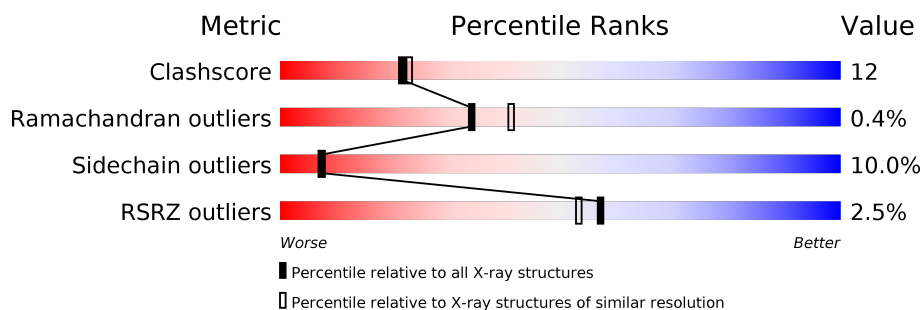
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.45 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))
Clashscore	141614	1613 (2.48-2.44)
Ramachandran outliers	138981	1598 (2.48-2.44)
Sidechain outliers	138945	1598 (2.48-2.44)
RSRZ outliers	127900	1523 (2.48-2.44)

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 on 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	1520	<div> <div></div> <div>70%</div> <div>23%</div> <div>...</div> </div>
1	B	1520	<div> <div>4%</div> <div>73%</div> <div>21%</div> <div>...</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
5	ONL	B	2511	-	-	X	-

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 23181 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.

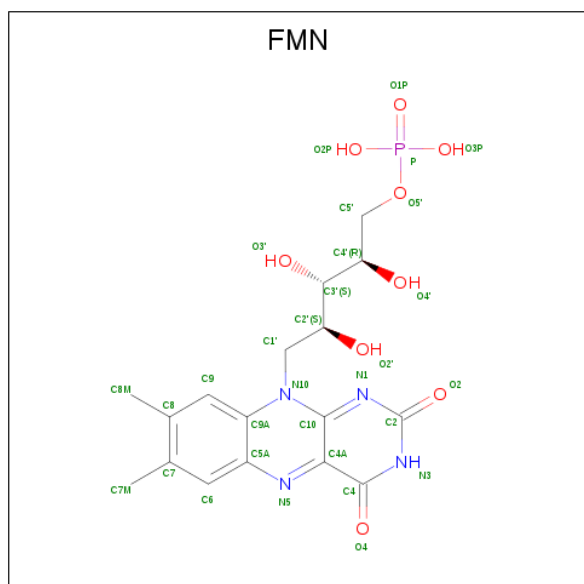
- Molecule 1 is a protein called FERREDONIN-DEPENDENT GLUTAMATE SYNTHASE 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	1485	Total	C	N	O	S	0	0	0
			11388	7184	1986	2163	55			
1	B	1485	Total	C	N	O	S	0	0	0
			11388	7184	1986	2163	55			

There are 6 discrepancies between the modelled and reference sequences:

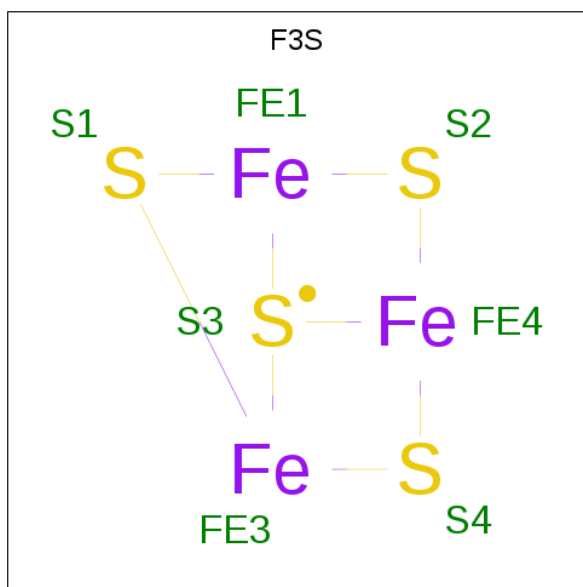
Chain	Residue	Modelled	Actual	Comment	Reference
A	578	ASP	THR	conflict	UNP P55038
A	581	THR	ASP	conflict	UNP P55038
A	1507	ASN	GLY	conflict	UNP P55038
B	578	ASP	THR	conflict	UNP P55038
B	581	THR	ASP	conflict	UNP P55038
B	1507	ASN	GLY	conflict	UNP P55038

- Molecule 2 is FLAVIN MONONUCLEOTIDE (three-letter code: FMN) (formula: C₁₇H₂₁N₄O₉P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			31	17	4	9	1		
2	B	1	Total	C	N	O	P	0	0
			31	17	4	9	1		

- Molecule 3 is FE3-S4 CLUSTER (three-letter code: F3S) (formula: Fe_3S_4).



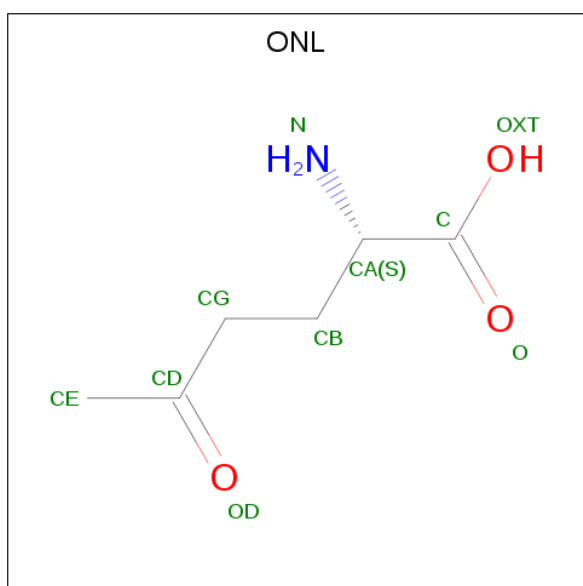
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	Fe	S	0	0
			7	3	4		
3	B	1	Total	Fe	S	0	0
			7	3	4		

- Molecule 4 is 2-OXOGLUTARIC ACID (three-letter code: AKG) (formula: $\text{C}_5\text{H}_6\text{O}_5$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			10	5	5		
4	B	1	Total	C	O	0	0
			10	5	5		

- Molecule 5 is 5-OXO-L-NORLEUCINE (three-letter code: ONL) (formula: $C_6H_{11}NO_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			10	6	1	3		
5	B	1	Total	C	N	O	0	0
			10	6	1	3		

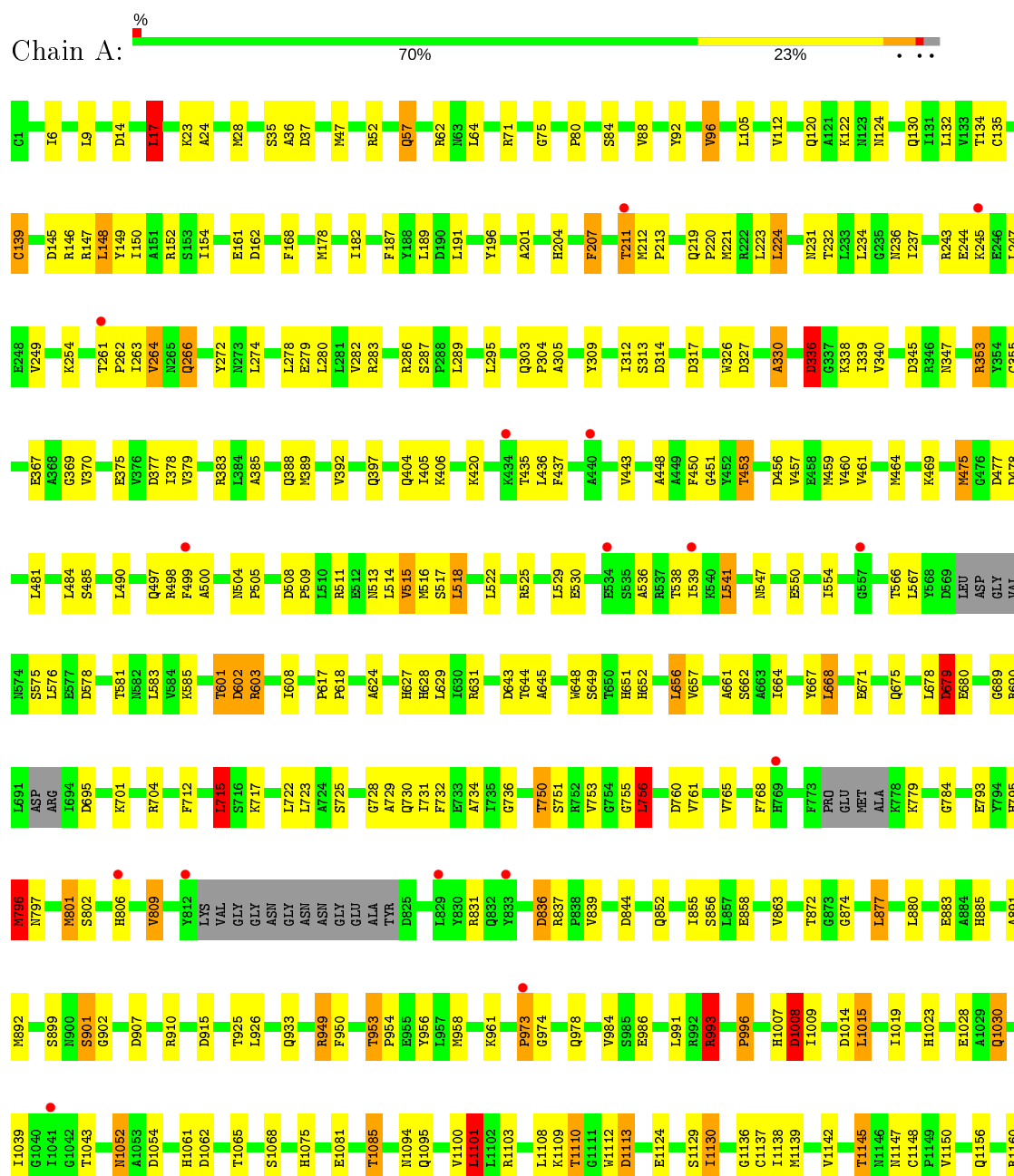
- Molecule 6 is water.

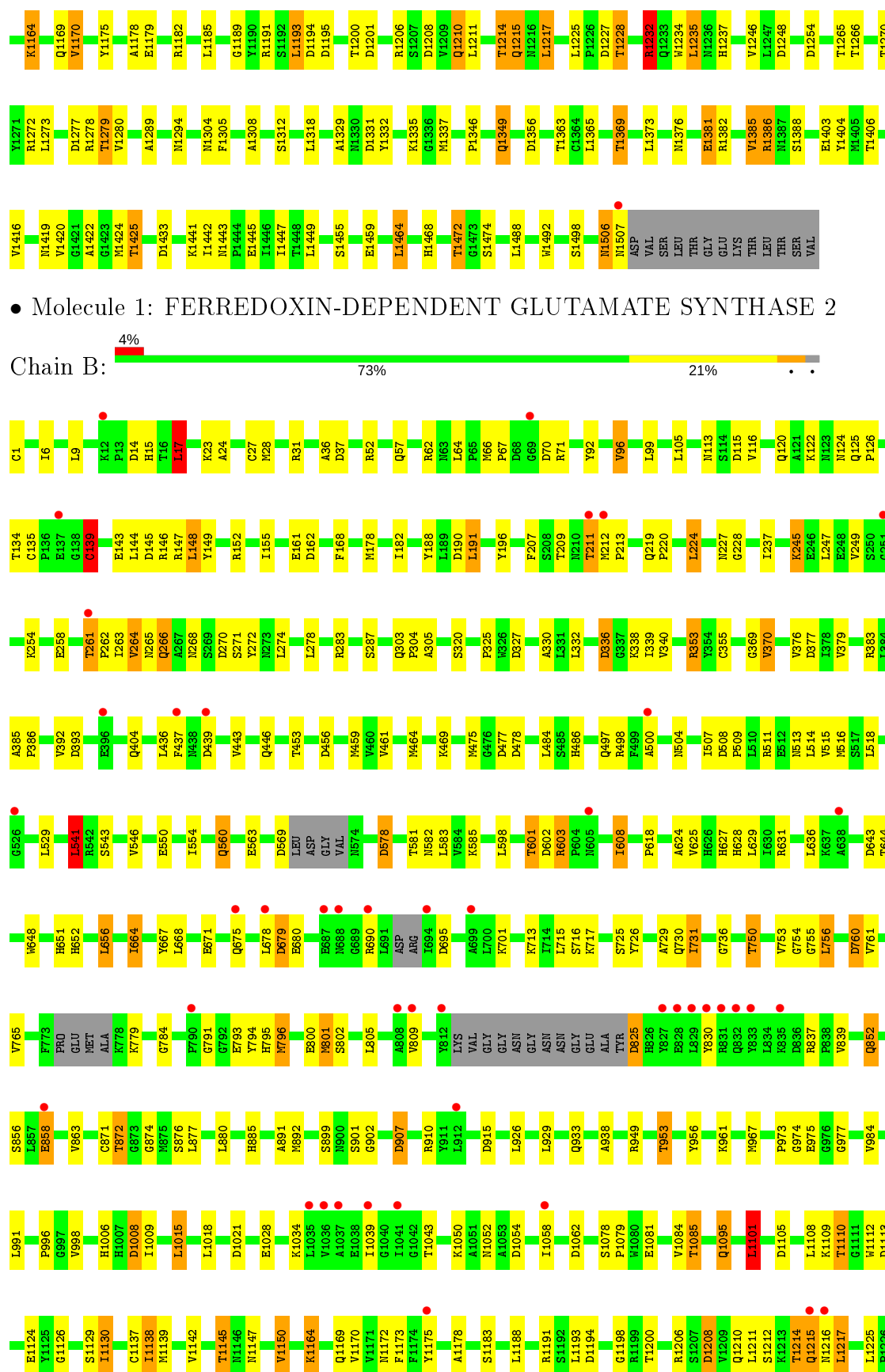
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	208	Total 208	O 208	0	0
6	B	81	Total 81	O 81	0	0

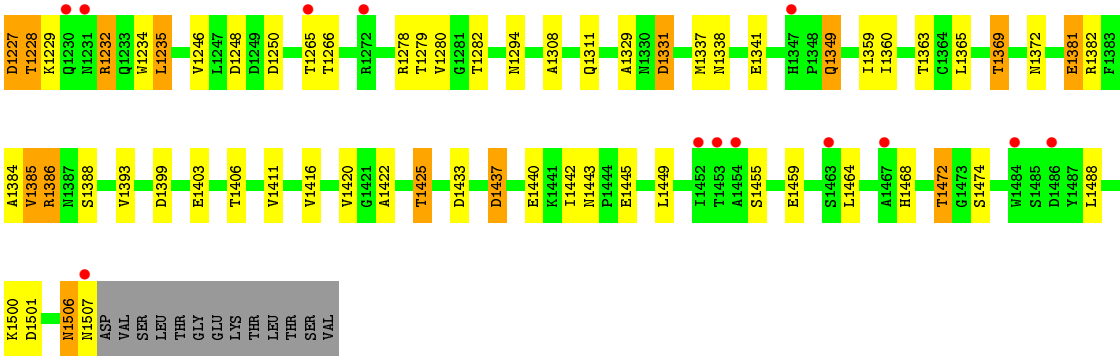
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA 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: FERREDOXIN-DEPENDENT GLUTAMATE SYNTHASE 2







4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	122.67Å 134.38Å 198.77Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.45 46.61 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.3 (20.00-2.45) 99.4 (46.61-2.40)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.15 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5	Depositor
R, R_{free}	0.207 , 0.257 0.205 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	37.8	Xtriage
Anisotropy	0.389	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 37.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	23181	wwPDB-VP
Average B, all atoms (Å ²)	17.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.84% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: FMN, F3S, AKG, ONL

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	4/11609 (0.0%)	1.02	56/15739 (0.4%)
1	B	0.67	0/11609	0.92	46/15739 (0.3%)
All	All	0.74	4/23218 (0.0%)	0.97	102/31478 (0.3%)

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	3
1	B	0	1
All	All	0	4

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	950	PHE	C-O	6.56	1.35	1.23
1	A	355	CYS	CB-SG	-5.72	1.72	1.81
1	A	475	MET	SD-CE	5.70	2.09	1.77
1	A	796	MET	SD-CE	5.09	2.06	1.77

The worst 5 of 102 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	353	ARG	NE-CZ-NH2	-12.44	114.08	120.30
1	A	949	ARG	NE-CZ-NH2	-11.31	114.65	120.30
1	A	1062	ASP	CB-CG-OD2	8.98	126.38	118.30
1	B	695	ASP	CB-CG-OD2	8.69	126.12	118.30
1	A	17	LEU	CA-CB-CG	8.49	134.83	115.30

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	330	ALA	Peptide
1	A	689	GLY	Peptide
1	A	728	GLY	Peptide
1	B	330	ALA	Peptide

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	11388	0	11335	296	0
1	B	11388	0	11336	245	0
2	A	31	0	19	1	0
2	B	31	0	19	3	0
3	A	7	0	0	0	0
3	B	7	0	0	0	0
4	A	10	0	4	1	0
4	B	10	0	4	0	0
5	A	10	0	9	0	0
5	B	10	0	10	7	0
6	A	208	0	0	20	0
6	B	81	0	0	12	0
All	All	23181	0	22736	543	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 12.

The worst 5 of 543 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:1130:ILE:CG1	1:A:1130:ILE:CD1	1.74	1.56
1:A:796:MET:CE	1:A:796:MET:SD	2.06	1.43
1:A:475:MET:CE	1:A:475:MET:SD	2.09	1.40
1:B:1:CYS:SG	5:B:2511:ONL:CE	2.20	1.29
1:B:1:CYS:SG	5:B:2511:ONL:CD	2.27	1.22

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	1475/1520 (97%)	1398 (95%)	72 (5%)	5 (0%)	41	49
1	B	1475/1520 (97%)	1394 (94%)	73 (5%)	8 (0%)	29	34
All	All	2950/3040 (97%)	2792 (95%)	145 (5%)	13 (0%)	34	41

5 of 13 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	679	ASP
1	B	679	ASP
1	B	1385	VAL
1	B	902	GLY
1	B	977	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	1209/1236 (98%)	1085 (90%)	124 (10%)	7	7
1	B	1209/1236 (98%)	1092 (90%)	117 (10%)	8	8
All	All	2418/2472 (98%)	2177 (90%)	241 (10%)	7	7

5 of 241 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1248	ASP
1	B	139	CYS
1	B	1228	THR
1	A	1280	VAL
1	A	1488	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 63 such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1376	ASN
1	B	120	GLN
1	B	1372	ASN
1	A	1419	ASN
1	A	1493	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

8 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	F3S	A	2509	1	0,9,9	0.00	-	-		
4	AKG	A	2510	-	3,9,9	5.25	1 (33%)	4,11,11	3.40	1 (25%)
2	FMN	A	2508	-	31,33,33	1.50	6 (19%)	40,50,50	2.50	9 (22%)
2	FMN	B	2508	-	31,33,33	1.83	6 (19%)	40,50,50	2.39	13 (32%)
3	F3S	B	2509	1	0,9,9	0.00	-	-		
4	AKG	B	2510	-	3,9,9	4.49	1 (33%)	4,11,11	4.89	3 (75%)
5	ONL	A	2511	1	5,9,9	0.81	0	5,11,11	2.97	4 (80%)
5	ONL	B	2511	-	5,9,9	0.47	0	5,11,11	2.84	1 (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
3	F3S	A	2509	1	-	-	0/3/3/3
4	AKG	A	2510	-	-	0/3/9/9	-
3	F3S	B	2509	1	-	-	0/3/3/3
2	FMN	A	2508	-	-	7/18/18/18	0/3/3/3
2	FMN	B	2508	-	-	2/18/18/18	0/3/3/3
5	ONL	B	2511	-	-	0/5/9/9	-
4	AKG	B	2510	-	-	1/3/9/9	-
5	ONL	A	2511	1	-	2/5/9/9	-

The worst 5 of 14 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	2510	AKG	O5-C2	9.02	1.36	1.22
4	B	2510	AKG	O5-C2	7.76	1.34	1.22
2	B	2508	FMN	C1'-N10	5.39	1.53	1.48
2	B	2508	FMN	C4A-N5	4.61	1.39	1.33
2	A	2508	FMN	C10-N1	3.90	1.38	1.33

The worst 5 of 31 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	2510	AKG	C4-C3-C2	-9.20	93.30	113.14
2	A	2508	FMN	C4-N3-C2	7.70	121.65	115.14
2	B	2508	FMN	C4-N3-C2	7.26	121.27	115.14
2	B	2508	FMN	C1'-N10-C10	6.83	124.53	118.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	2510	AKG	C4-C3-C2	-6.64	98.81	113.14

There are no chirality outliers.

5 of 12 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	2508	FMN	C2'-C1'-N10-C9A
2	A	2508	FMN	O3'-C3'-C4'-C5'
2	A	2508	FMN	C4'-C5'-O5'-P
2	B	2508	FMN	C4'-C5'-O5'-P
2	A	2508	FMN	O3'-C3'-C4'-O4'

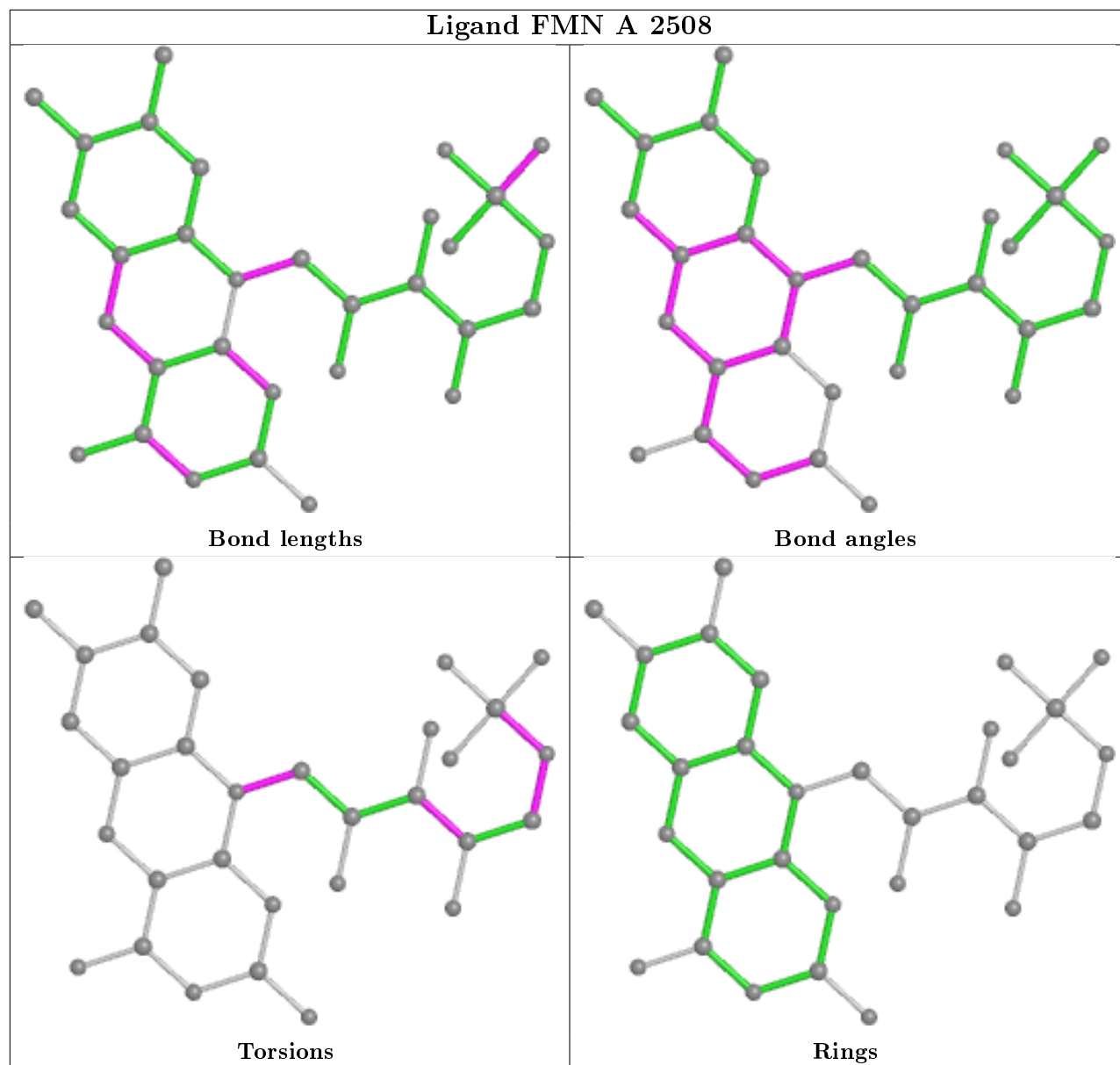
There are no ring outliers.

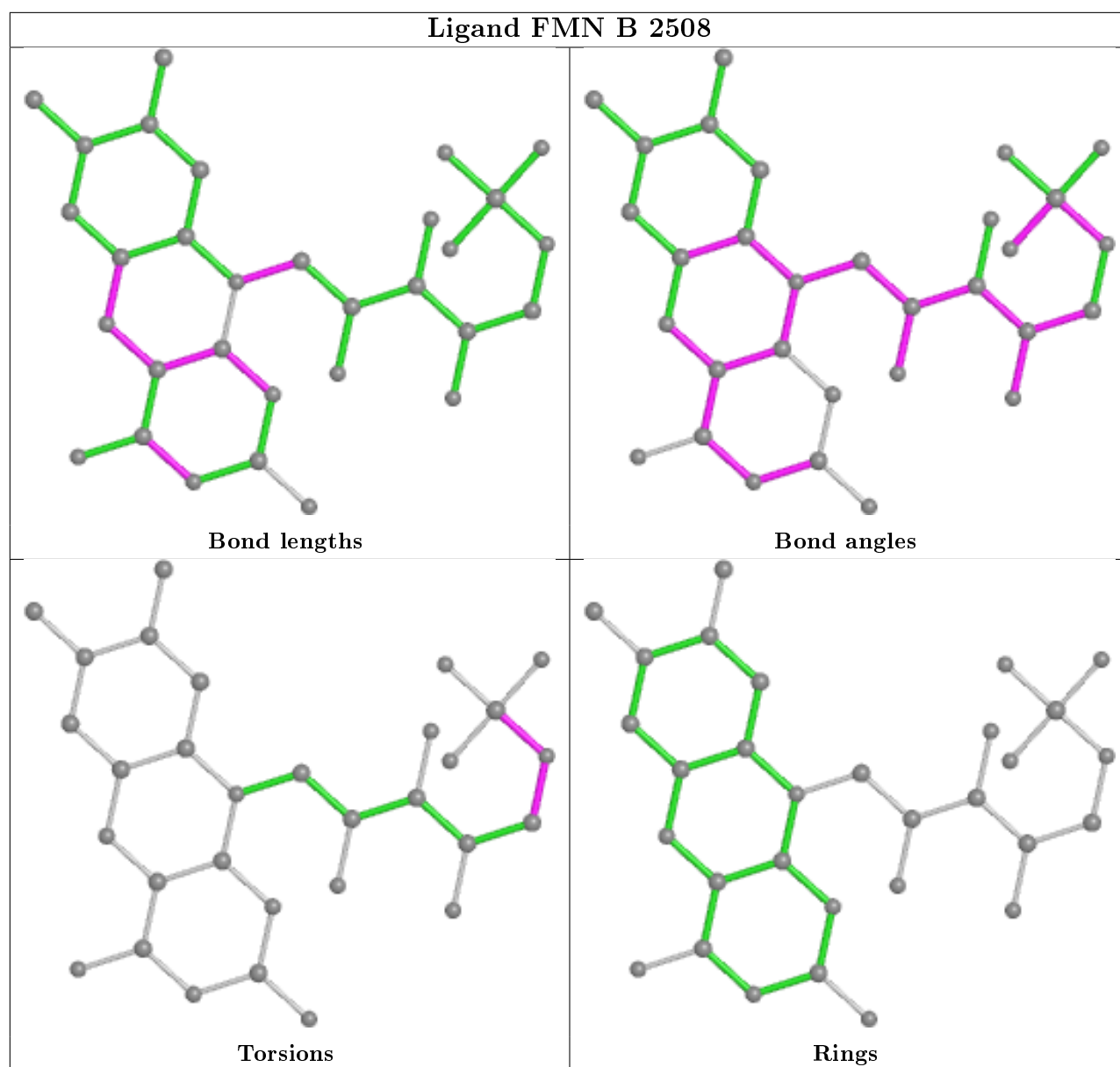
4 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	2510	AKG	1	0
2	A	2508	FMN	1	0
2	B	2508	FMN	3	0
5	B	2511	ONL	7	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.

Ligand FMN A 2508





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	1485/1520 (97%)	-0.21	17 (1%) 80 80	8, 16, 23, 29	0
1	B	1485/1520 (97%)	0.08	57 (3%) 40 37	7, 17, 23, 31	0
All	All	2970/3040 (97%)	-0.07	74 (2%) 57 53	7, 16, 23, 31	0

The worst 5 of 74 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	1507	ASN	5.1
1	B	812	TYR	5.0
1	B	1453	THR	4.0
1	B	833	TYR	4.0
1	B	261	THR	3.6

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 carbohydrates in this entry.

6.4 Ligands [i](#)

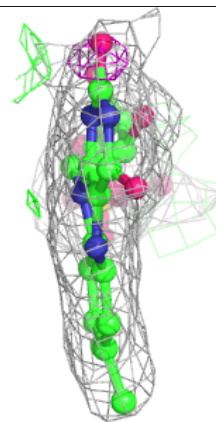
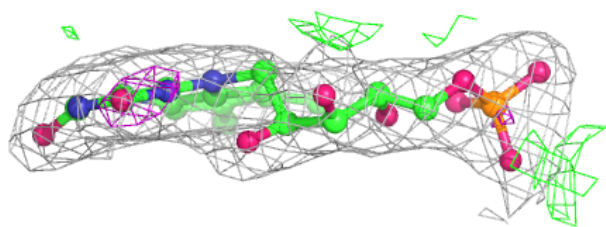
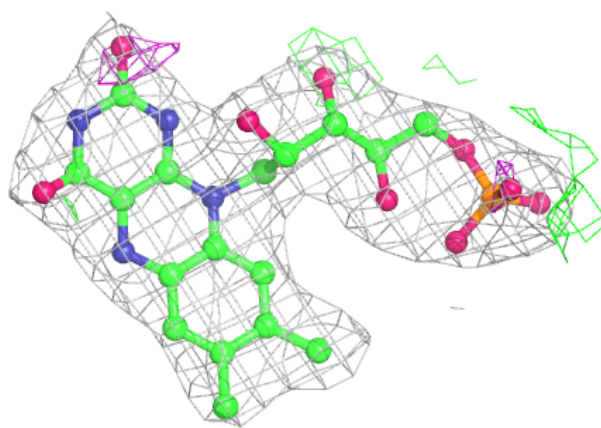
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
5	ONL	A	2511	10/10	0.90	0.16	24,32,34,36	0
5	ONL	B	2511	10/10	0.92	0.14	32,36,37,38	0
4	AKG	B	2510	10/10	0.93	0.17	36,37,40,41	0
2	FMN	B	2508	31/31	0.96	0.13	22,29,32,35	0
4	AKG	A	2510	10/10	0.96	0.11	24,27,35,37	0
3	F3S	B	2509	7/7	0.97	0.04	27,29,30,31	0
2	FMN	A	2508	31/31	0.97	0.14	18,23,28,31	0
3	F3S	A	2509	7/7	0.99	0.09	19,22,24,24	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.

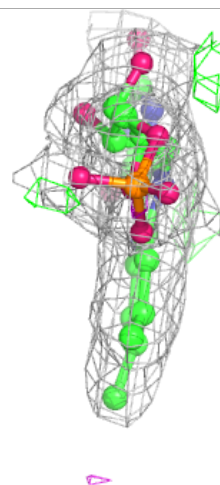
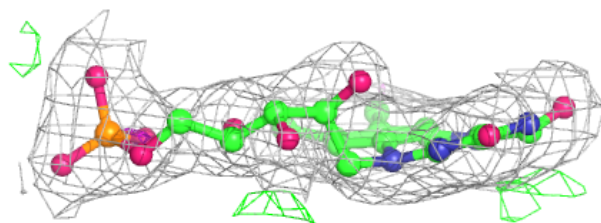
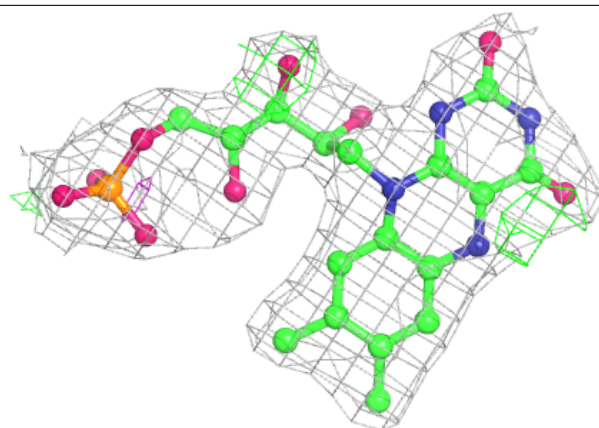
Electron density around FMN B 2508:

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 FMN A 2508:

$2mF_o-DF_c$ (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.