



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 7, 2026 – 12:44 AM UTC

PDB ID : 2HQM / pdb_00002hqm
Title : Crystal Structure of Glutathione Reductase Glr1 from the Yeast *Saccharomyces cerevisiae*
Authors : Yu, J.; Zhou, C.Z.
Deposited on : 2006-07-19
Resolution : 2.40 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.

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A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

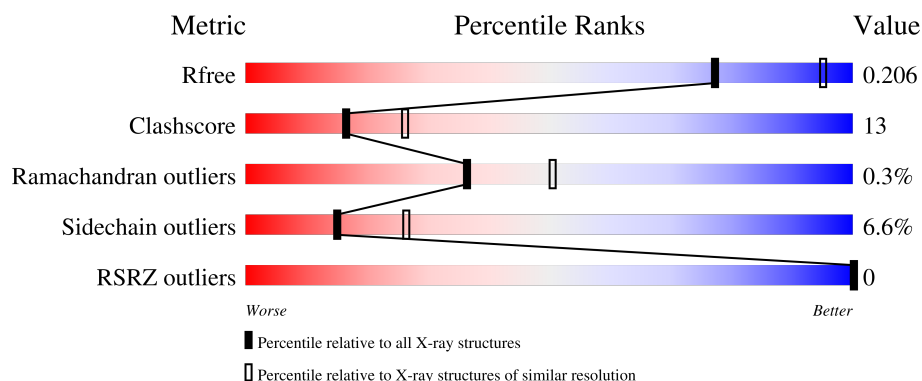
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	479	 77% 16% . .
1	B	479	 72% 21% . .

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	GSH	A	3001	-	-	X	-
5	GSH	B	4001	-	-	X	-
6	GOL	A	6001	-	X	X	-
6	GOL	A	6006	-	X	X	-
6	GOL	B	6003	-	X	-	-
6	GOL	B	6004	-	X	X	-
6	GOL	B	6007	-	X	X	-
6	GOL	B	6009	-	X	X	-
6	GOL	B	6010	-	X	-	-

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 8152 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.

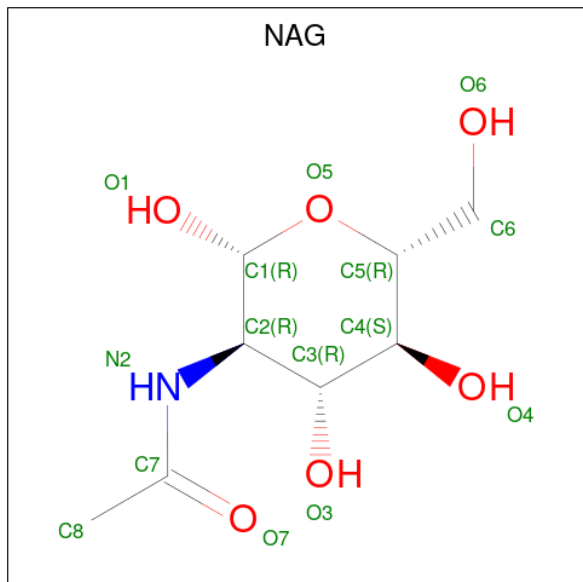
- Molecule 1 is a protein called Glutathione reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	461	Total	C	N	O	S	0	0	0
			3588	2277	620	679	12			
1	B	461	Total	C	N	O	S	0	0	0
			3588	2277	620	679	12			

There are 24 discrepancies between the modelled and reference sequences:

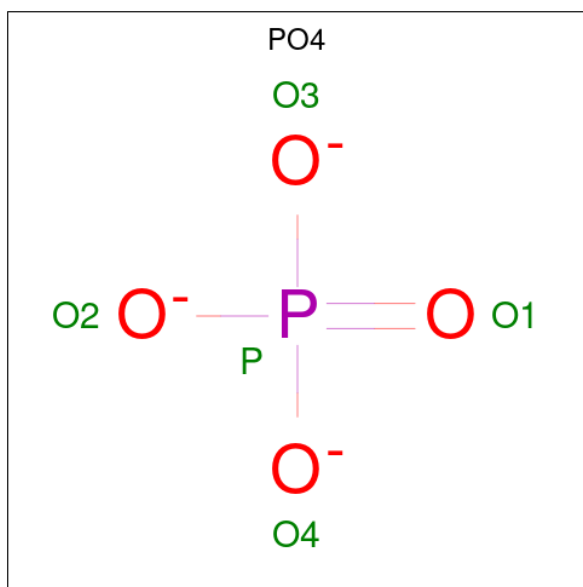
Chain	Residue	Modelled	Actual	Comment	Reference
A	13	TYR	-	cloning artifact	UNP P41921
A	14	VAL	-	cloning artifact	UNP P41921
A	15	GLU	-	cloning artifact	UNP P41921
A	16	PHE	-	cloning artifact	UNP P41921
A	484	GLY	-	cloning artifact	UNP P41921
A	485	SER	-	cloning artifact	UNP P41921
A	486	HIS	-	cloning artifact	UNP P41921
A	487	HIS	-	cloning artifact	UNP P41921
A	488	HIS	-	cloning artifact	UNP P41921
A	489	HIS	-	cloning artifact	UNP P41921
A	490	HIS	-	cloning artifact	UNP P41921
A	491	HIS	-	cloning artifact	UNP P41921
B	13	TYR	-	cloning artifact	UNP P41921
B	14	VAL	-	cloning artifact	UNP P41921
B	15	GLU	-	cloning artifact	UNP P41921
B	16	PHE	-	cloning artifact	UNP P41921
B	484	GLY	-	cloning artifact	UNP P41921
B	485	SER	-	cloning artifact	UNP P41921
B	486	HIS	-	cloning artifact	UNP P41921
B	487	HIS	-	cloning artifact	UNP P41921
B	488	HIS	-	cloning artifact	UNP P41921
B	489	HIS	-	cloning artifact	UNP P41921
B	490	HIS	-	cloning artifact	UNP P41921
B	491	HIS	-	cloning artifact	UNP P41921

- Molecule 2 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



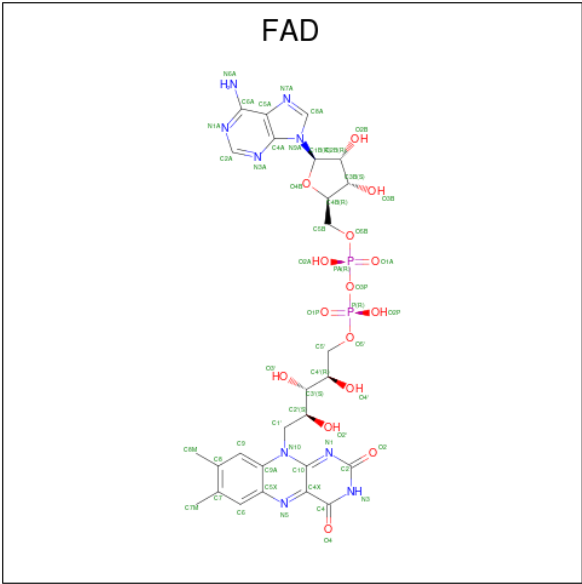
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	B	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 3 is PHOSPHATE ION (CCD ID: PO4) (formula: O_4P).



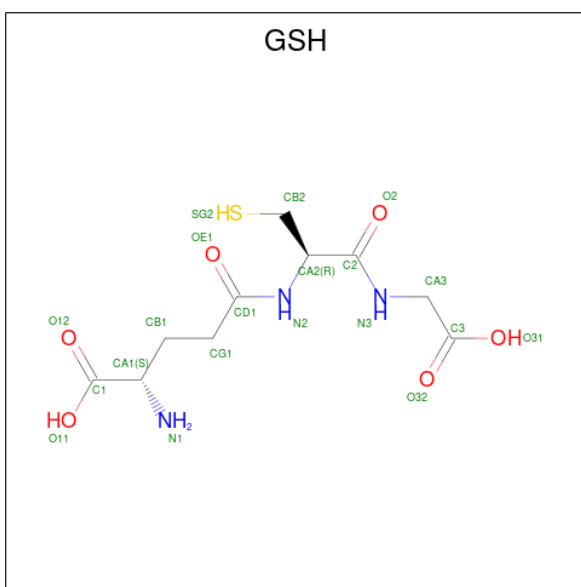
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	P	0	0
			5	4	1		
3	B	1	Total	O	P	0	0
			5	4	1		

- Molecule 4 is FLAVIN-ADENINE DINUCLEOTIDE (CCD ID: FAD) (formula: C₂₇H₃₃N₉O₁₅P₂).



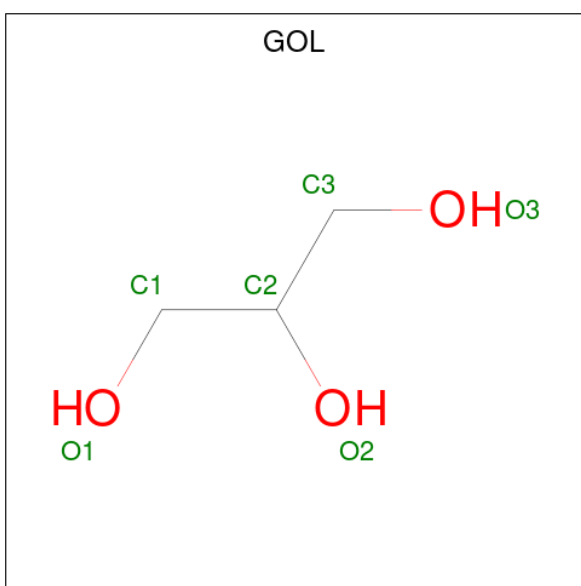
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total 53	C 27	N 9	O 15	P 2	0	0
4	B	1	Total 53	C 27	N 9	O 15	P 2	0	0

- Molecule 5 is GLUTATHIONE (CCD ID: GSH) (formula: C₁₀H₁₇N₃O₆S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	S	0	0
			20	10	3	6	1		
5	B	1	Total	C	N	O	S	0	0
			20	10	3	6	1		

- Molecule 6 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

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

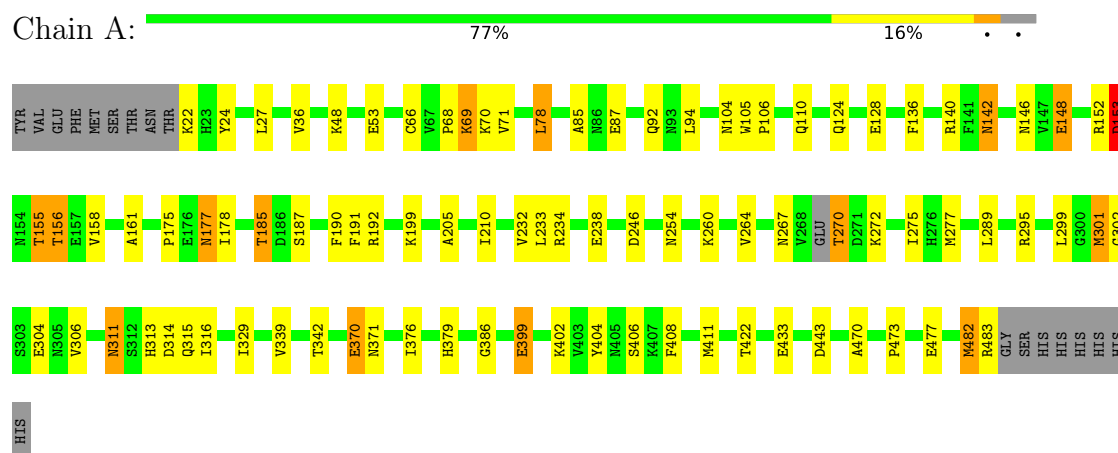
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	377	Total	O	0	0
			377	377		
7	B	349	Total	O	0	0
			349	349		

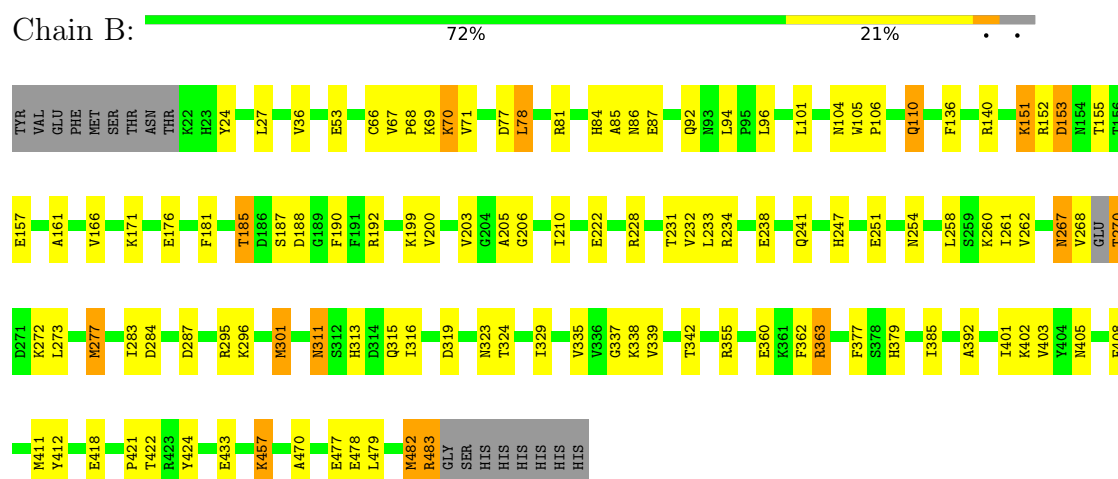
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: Glutathione reductase



• Molecule 1: Glutathione reductase



4 Data and refinement statistics

Property	Value	Source
Space group	P 42	Depositor
Cell constants a, b, c, α , β , γ	116.97Å 116.97Å 85.24Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	21.05 – 2.40 21.05 – 2.40	Depositor EDS
% Data completeness (in resolution range)	92.8 (21.05-2.40) 92.9 (21.05-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.95 (at 2.41Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.202 , 0.205 0.203 , 0.206	Depositor DCC
R_{free} test set	2118 reflections (4.70%)	wwPDB-VP
Wilson B-factor (Å ²)	22.4	Xtriage
Anisotropy	0.062	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 33.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.487 for h,-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	8152	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.58% 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: FAD, PO4, NAG, GSH, GOL

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	0.72	1/3658 (0.0%)	0.95	7/4939 (0.1%)
1	B	0.76	4/3658 (0.1%)	0.94	6/4939 (0.1%)
All	All	0.74	5/7316 (0.1%)	0.94	13/9878 (0.1%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	482	MET	C-N	9.26	1.46	1.33
1	B	482	MET	CA-C	7.89	1.64	1.53
1	B	87	GLU	CA-CB	-6.49	1.42	1.53
1	B	482	MET	N-CA	5.68	1.52	1.45
1	A	482	MET	C-N	5.46	1.41	1.33

The worst 5 of 13 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	153	ASP	N-CA-C	-14.81	95.20	114.31
1	B	87	GLU	CB-CG-CD	-7.29	100.21	112.60
1	A	482	MET	N-CA-C	6.59	120.63	110.42
1	B	482	MET	CA-C-N	6.58	133.54	121.70
1	B	482	MET	C-N-CA	6.58	133.54	121.70

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	3588	0	3579	84	0
1	B	3588	0	3579	98	0
2	A	14	0	13	0	0
2	B	14	0	13	0	0
3	A	5	0	0	1	0
3	B	5	0	0	1	0
4	A	53	0	31	2	0
4	B	53	0	31	0	0
5	A	20	0	14	8	0
5	B	20	0	14	8	0
6	A	30	0	40	15	0
6	B	36	0	44	21	0
7	A	377	0	0	16	0
7	B	349	0	0	11	0
All	All	8152	0	7358	190	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 13.

The worst 5 of 190 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)
6:B:6009:GOL:C2	6:B:6009:GOL:O2	1.63	1.46
6:A:6001:GOL:O2	6:A:6001:GOL:C2	1.64	1.43
6:A:6006:GOL:O2	6:A:6006:GOL:C2	1.67	1.40
6:B:6004:GOL:C2	6:B:6004:GOL:O2	1.73	1.37
1:B:110:GLN:HE21	1:B:110:GLN:HA	1.23	1.03

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	457/479 (95%)	438 (96%)	17 (4%)	2 (0%)	30	43
1	B	457/479 (95%)	438 (96%)	18 (4%)	1 (0%)	43	58
All	All	914/958 (95%)	876 (96%)	35 (4%)	3 (0%)	36	50

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	156	THR
1	A	339	VAL
1	B	339	VAL

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	386/403 (96%)	363 (94%)	23 (6%)	17	31
1	B	386/403 (96%)	358 (93%)	28 (7%)	13	22
All	All	772/806 (96%)	721 (93%)	51 (7%)	15	26

5 of 51 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	101	LEU
1	B	260	LYS
1	B	457	LYS
1	B	110	GLN
1	B	152	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 33 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	311	ASN

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Mol	Chain	Res	Type
1	B	322	GLN
1	B	405	ASN
1	A	298	HIS
1	A	276	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 [i](#)

19 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	$\# Z > 2$
6	GOL	B	6007	-	5,5,5	2.07	2 (40%)	5,5,5	3.50	3 (60%)
6	GOL	B	6010	-	5,5,5	2.42	1 (20%)	5,5,5	2.94	5 (100%)
6	GOL	B	6009	-	5,5,5	3.66	3 (60%)	5,5,5	3.47	5 (100%)
5	GSH	B	4001	1	18,19,19	1.33	2 (11%)	21,24,24	2.22	7 (33%)
6	GOL	A	6005	-	5,5,5	0.35	0	5,5,5	1.08	1 (20%)
6	GOL	A	6006	-	5,5,5	4.54	4 (80%)	5,5,5	4.37	5 (100%)
3	PO4	A	5001	-	4,4,4	1.45	1 (25%)	6,6,6	0.48	0
6	GOL	A	6001	-	5,5,5	3.66	3 (60%)	5,5,5	4.29	5 (100%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	GOL	B	6003	-	5,5,5	5.03	4 (80%)	5,5,5	5.96	5 (100%)
4	FAD	B	2001	-	58,58,58	1.24	7 (12%)	85,89,89	1.54	14 (16%)
4	FAD	A	1001	-	58,58,58	1.25	5 (8%)	85,89,89	1.59	14 (16%)
3	PO4	B	5002	-	4,4,4	1.52	1 (25%)	6,6,6	0.62	0
6	GOL	A	6008	-	5,5,5	1.35	1 (20%)	5,5,5	1.90	2 (40%)
6	GOL	A	6011	-	5,5,5	1.78	1 (20%)	5,5,5	2.00	1 (20%)
6	GOL	B	6004	-	5,5,5	4.93	4 (80%)	5,5,5	4.21	4 (80%)
2	NAG	B	1303	1	14,14,15	0.35	0	17,19,21	0.85	1 (5%)
2	NAG	A	1301	1	14,14,15	0.59	0	17,19,21	0.70	0
5	GSH	A	3001	1	18,19,19	1.35	3 (16%)	21,24,24	1.68	7 (33%)
6	GOL	B	6002	-	5,5,5	1.03	1 (20%)	5,5,5	1.31	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
2	NAG	B	1303	1	-	2/6/23/26	0/1/1/1
6	GOL	B	6009	-	-	4/4/4/4	-
6	GOL	B	6010	-	-	4/4/4/4	-
2	NAG	A	1301	1	-	0/6/23/26	0/1/1/1
5	GSH	B	4001	1	-	3/24/24/24	-
6	GOL	A	6008	-	-	2/4/4/4	-
6	GOL	A	6005	-	-	0/4/4/4	-
5	GSH	A	3001	1	-	7/24/24/24	-
6	GOL	A	6011	-	-	3/4/4/4	-
6	GOL	A	6006	-	-	4/4/4/4	-
6	GOL	B	6002	-	-	2/4/4/4	-
6	GOL	B	6004	-	-	2/4/4/4	-
6	GOL	B	6007	-	-	4/4/4/4	-
6	GOL	A	6001	-	-	3/4/4/4	-
6	GOL	B	6003	-	-	3/4/4/4	-
4	FAD	A	1001	-	-	4/34/50/50	0/6/6/6
4	FAD	B	2001	-	-	4/34/50/50	0/6/6/6

The worst 5 of 43 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	6004	GOL	O2-C2	10.29	1.73	1.43
6	B	6003	GOL	O2-C2	-9.09	1.17	1.43
6	A	6006	GOL	O2-C2	8.33	1.67	1.43
6	A	6001	GOL	O2-C2	7.24	1.64	1.43
6	B	6009	GOL	O2-C2	6.95	1.63	1.43

The worst 5 of 80 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	B	6003	GOL	O2-C2-C1	6.76	137.15	109.18
6	B	6007	GOL	C3-C2-C1	-6.44	88.18	111.80
6	B	6003	GOL	C3-C2-C1	-6.34	88.54	111.80
6	B	6003	GOL	O2-C2-C3	6.00	134.01	109.18
4	A	1001	FAD	C5A-C4A-N3A	-5.92	118.57	126.72

There are no chirality outliers.

5 of 51 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	1303	NAG	C8-C7-N2-C2
5	A	3001	GSH	N1-CA1-CB1-CG1
5	A	3001	GSH	C1-CA1-CB1-CG1
6	A	6001	GOL	C1-C2-C3-O3
6	A	6006	GOL	O1-C1-C2-C3

There are no ring outliers.

15 monomers are involved in 56 short contacts:

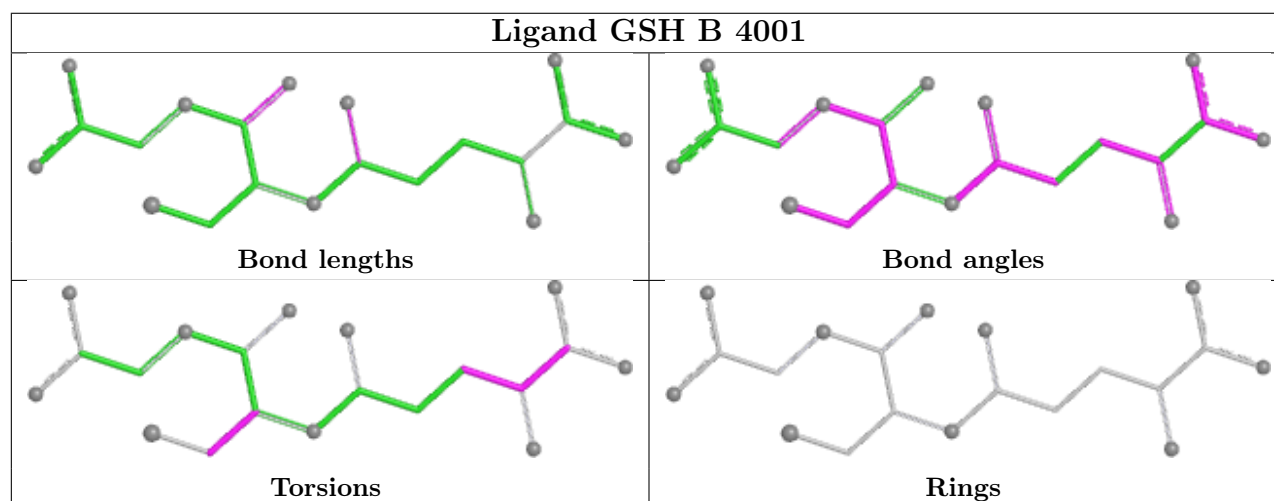
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	B	6007	GOL	4	0
6	B	6010	GOL	3	0
6	B	6009	GOL	6	0
5	B	4001	GSH	8	0
6	A	6005	GOL	2	0
6	A	6006	GOL	4	0
3	A	5001	PO4	1	0
6	A	6001	GOL	8	0
6	B	6003	GOL	3	0
4	A	1001	FAD	2	0
3	B	5002	PO4	1	0
6	A	6011	GOL	1	0
6	B	6004	GOL	4	0

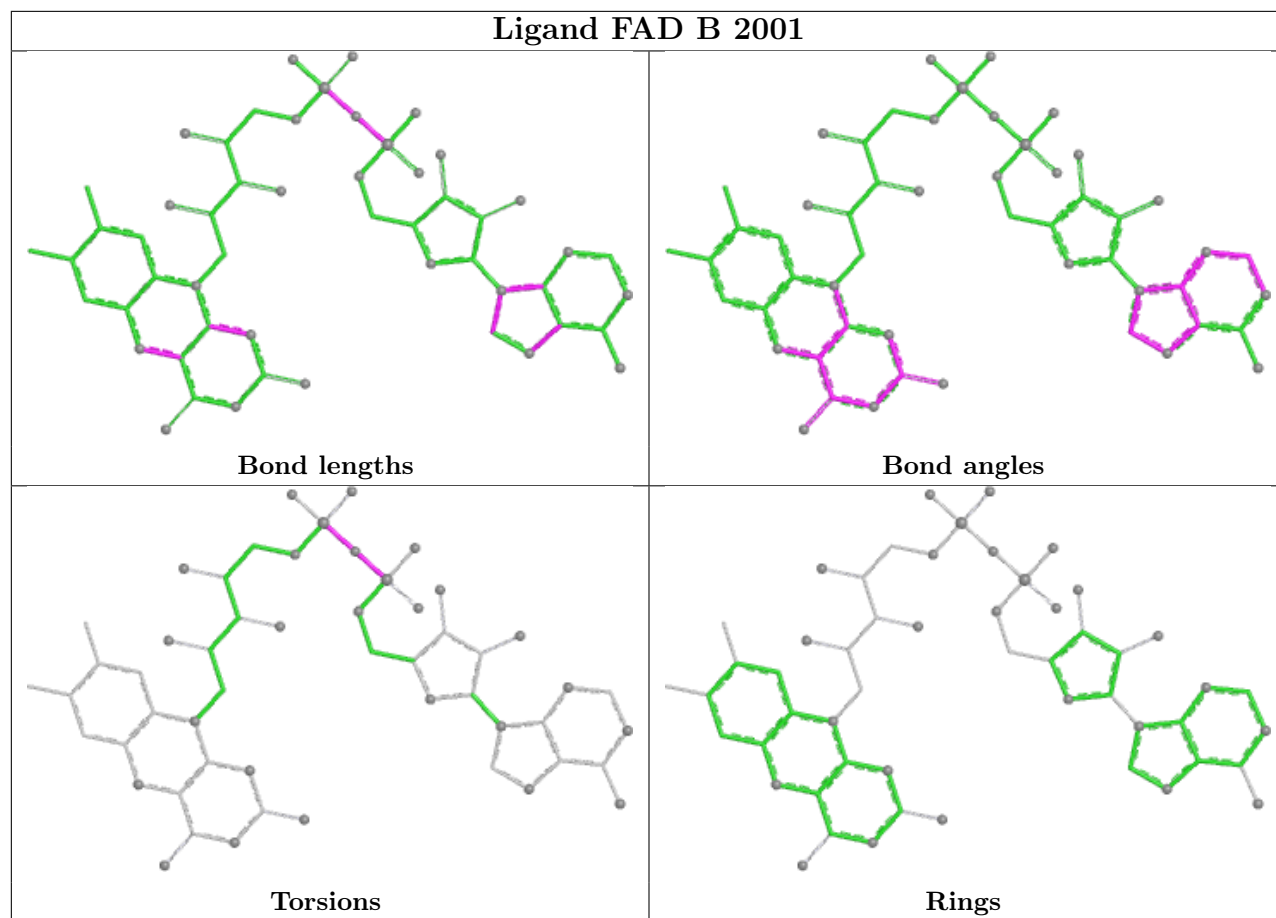
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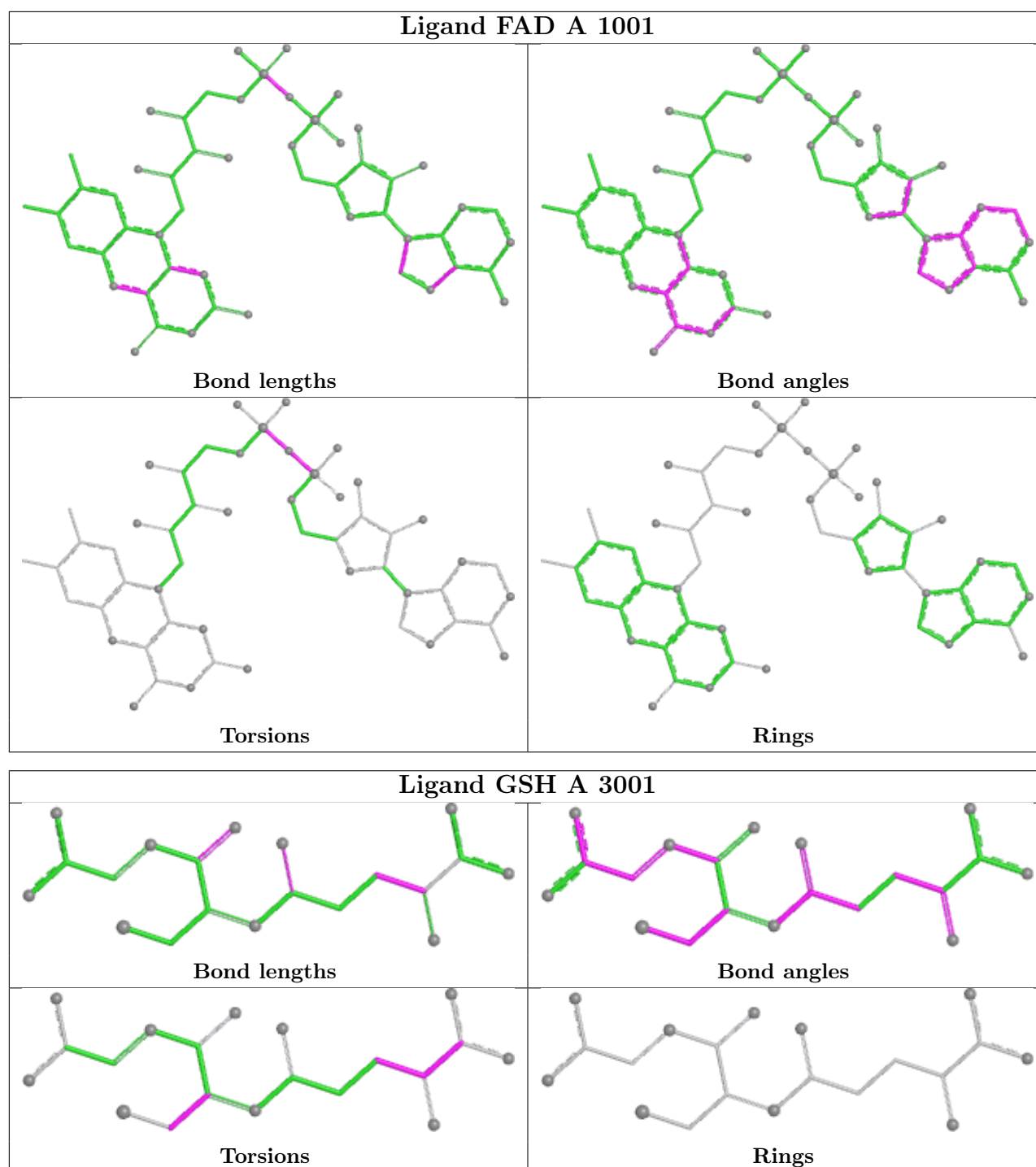
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	3001	GSH	8	0
6	B	6002	GOL	1	0

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	461/479 (96%)	-1.76	0 100 100	12, 22, 48, 73	0
1	B	461/479 (96%)	-1.76	0 100 100	13, 22, 47, 68	0
All	All	922/958 (96%)	-1.76	0 100 100	12, 22, 48, 73	0

There are no RSRZ outliers to report.

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

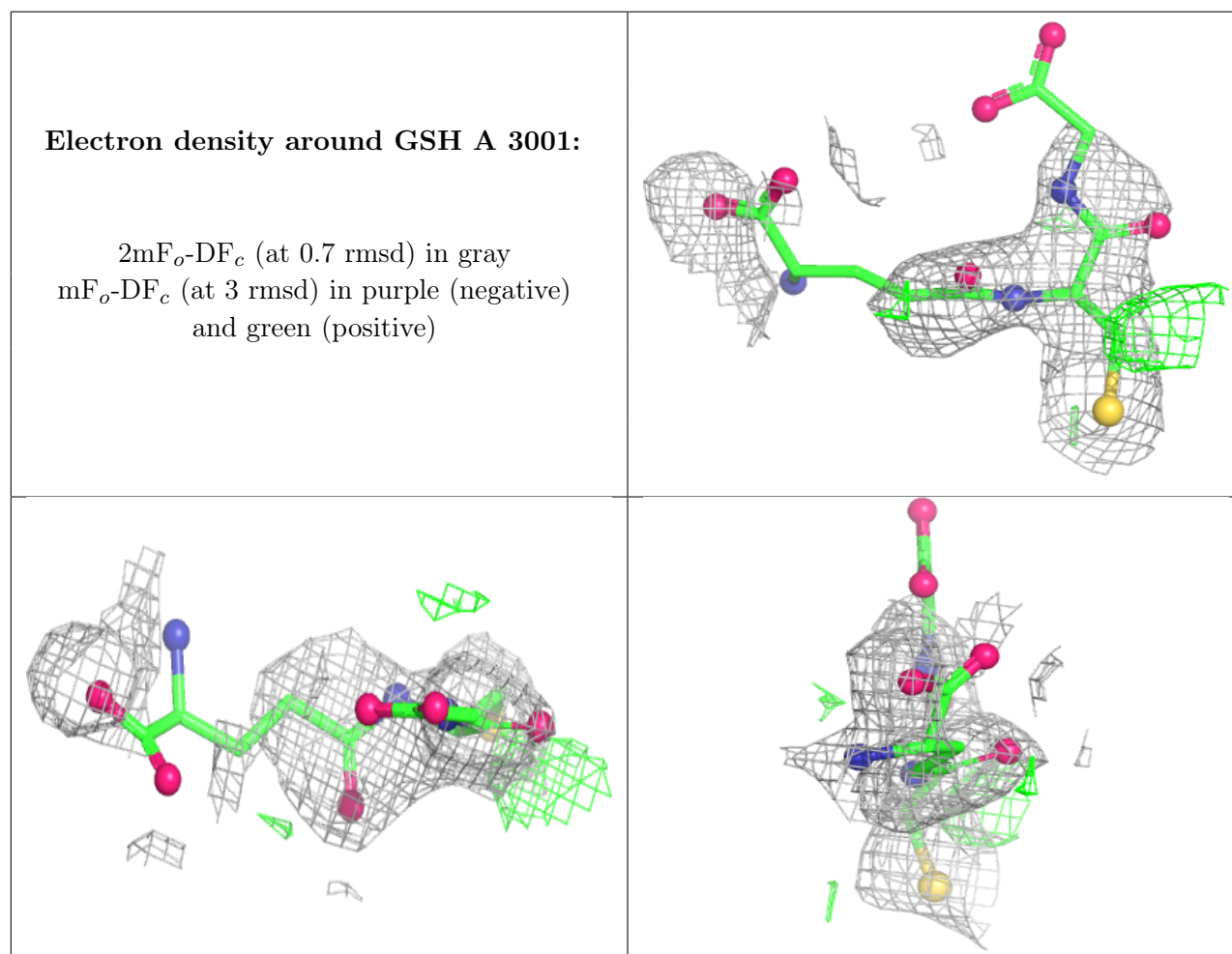
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
6	GOL	B	6009	6/6	0.98	0.06	40,42,43,43	6
3	PO4	A	5001	5/5	0.99	0.04	41,42,43,44	5
3	PO4	B	5002	5/5	0.99	0.04	40,40,41,41	5
5	GSH	A	3001	20/20	0.99	0.05	27,43,54,54	20
5	GSH	B	4001	20/20	0.99	0.05	23,40,49,49	20
6	GOL	A	6005	6/6	0.99	0.03	32,33,33,34	6
6	GOL	A	6006	6/6	0.99	0.06	38,38,39,41	6

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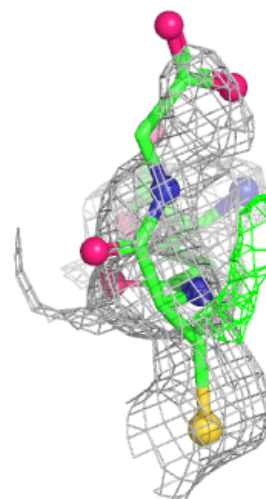
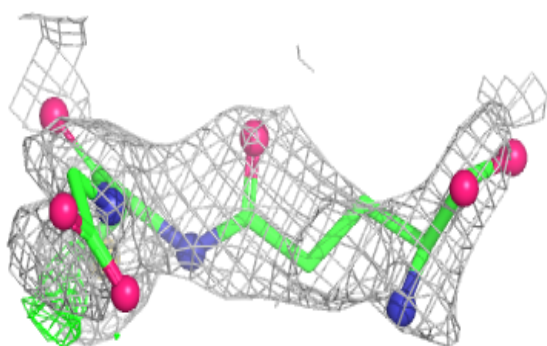
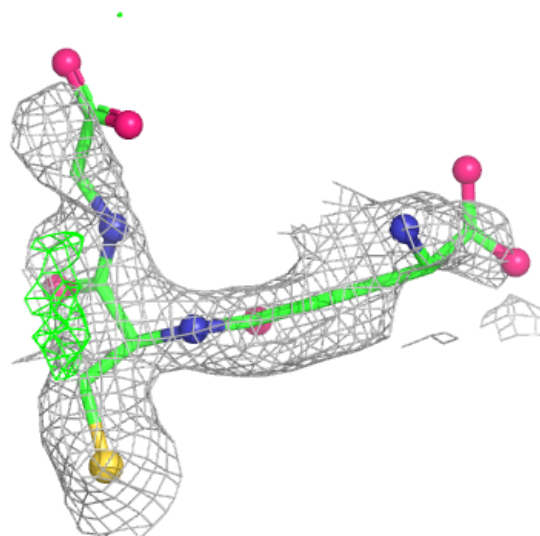
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
6	GOL	A	6008	6/6	0.99	0.06	41,46,47,49	6
6	GOL	A	6011	6/6	0.99	0.04	39,41,42,44	6
6	GOL	B	6002	6/6	0.99	0.07	39,43,44,45	6
6	GOL	B	6003	6/6	0.99	0.03	28,30,33,35	6
6	GOL	B	6004	6/6	0.99	0.03	25,30,31,35	6
6	GOL	B	6007	6/6	0.99	0.04	29,30,32,33	6
2	NAG	B	1303	14/15	0.99	0.02	31,35,38,40	0
6	GOL	B	6010	6/6	0.99	0.03	46,47,48,48	6
4	FAD	A	1001	53/53	1.00	0.02	10,17,24,26	0
6	GOL	A	6001	6/6	1.00	0.04	33,36,39,42	6
4	FAD	B	2001	53/53	1.00	0.02	12,16,22,24	0
2	NAG	A	1301	14/15	1.00	0.02	30,33,36,36	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.



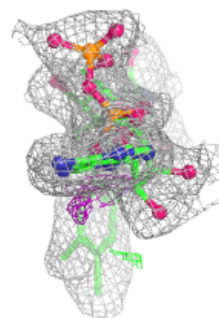
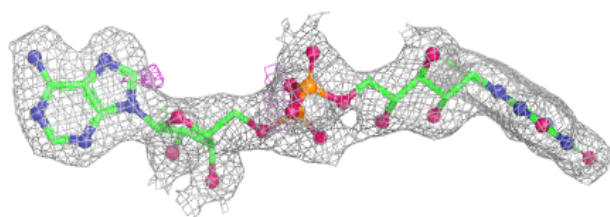
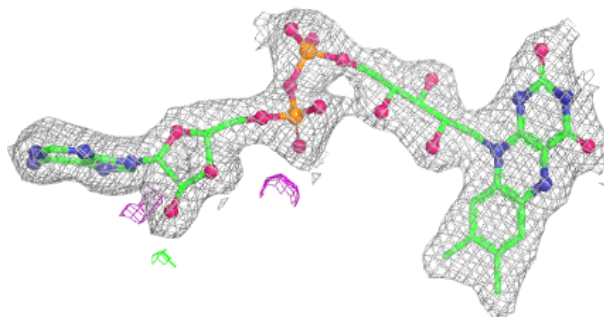
Electron density around GSH B 4001:

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

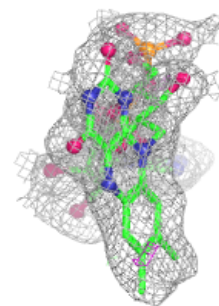
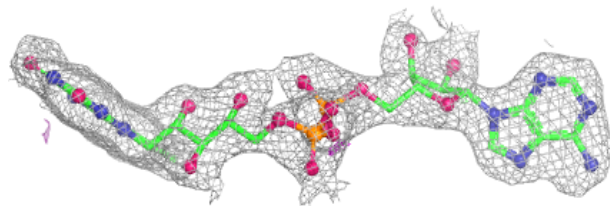
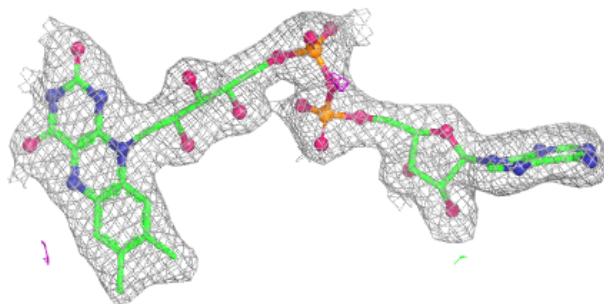


Electron density around FAD A 1001:

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

**Electron density around FAD B 2001:**

$2mF_o-DF_c$ (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.