



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 01:27 AM UTC

PDB ID : 3BHK / pdb_00003bhk
Title : Crystal structure of R49K mutant of monomeric sarcosine oxidase crystallized in phosphate as precipitant
Authors : Hassan-Abdallah, A.; Zhao, G.; Chen, Z.; Mathews, F.S.; Jorns, M.S.
Deposited on : 2007-11-28
Resolution : 1.71 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

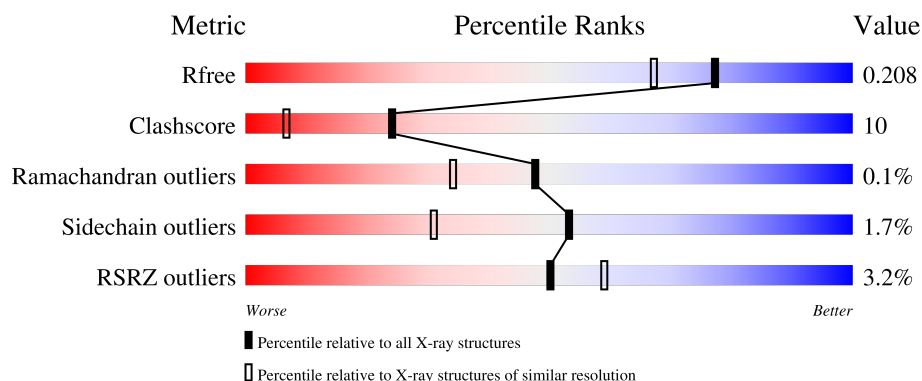
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

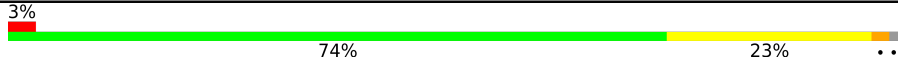
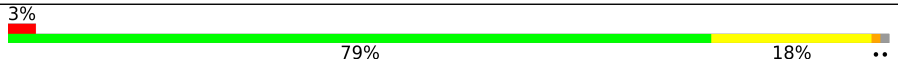
The reported resolution of this entry is 1.71 Å.

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	1039 (1.72-1.72)
Clashscore	190562	1049 (1.72-1.72)
Ramachandran outliers	187476	1041 (1.72-1.72)
Sidechain outliers	187428	1041 (1.72-1.72)
RSRZ outliers	180081	1039 (1.72-1.72)

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	390	
1	B	390	

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
4	GOL	A	901	-	-	X	-
4	GOL	B	902	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6790 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	385	Total	C	N	O	S	0	0	0
			3006	1916	503	577	10			
1	B	385	Total	C	N	O	S	0	0	0
			3006	1916	503	577	10			

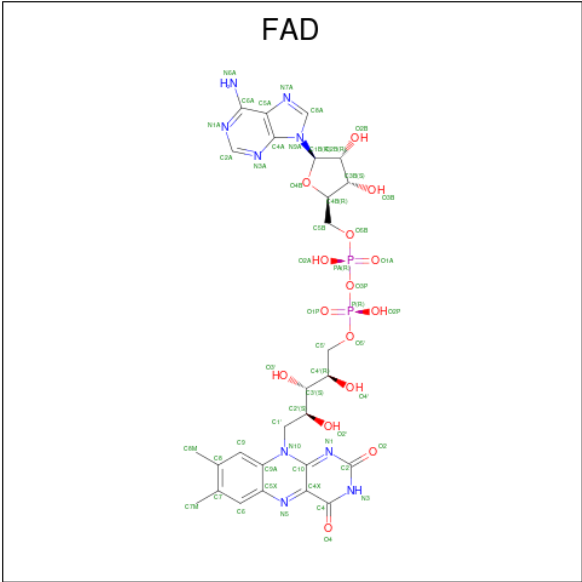
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	49	LYS	ARG	engineered mutation	UNP P40859
B	49	LYS	ARG	engineered mutation	UNP P40859

- Molecule 2 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

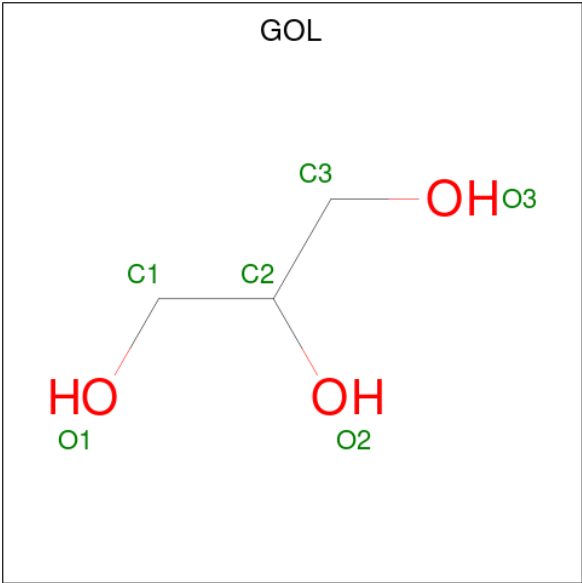
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Cl	0	0
			1	1		
2	B	1	Total	Cl	0	0
			1	1		

- Molecule 3 is FLAVIN-ADENINE DINUCLEOTIDE (CCD ID: FAD) (formula: C₂₇H₃₃N₉O₁₅P₂).



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

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



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

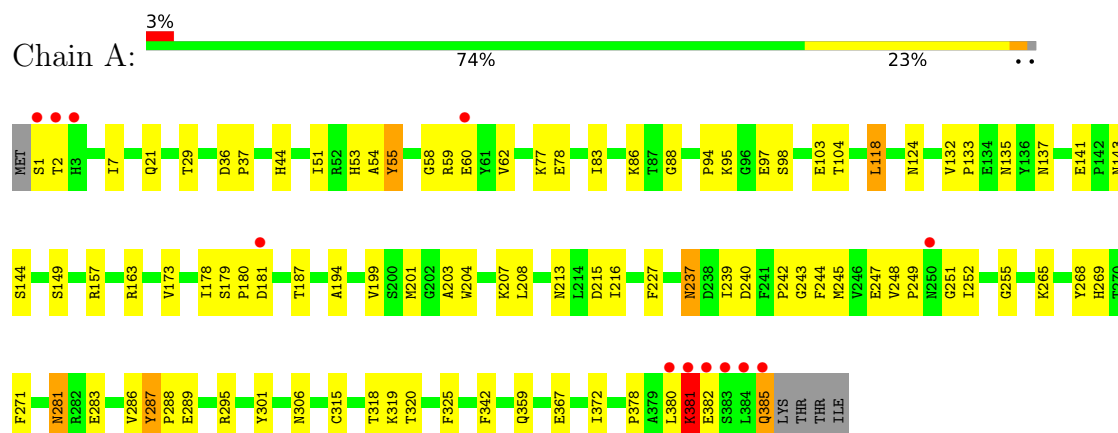
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	328	Total 328	O 328	0	0
5	B	330	Total 330	O 330	0	0

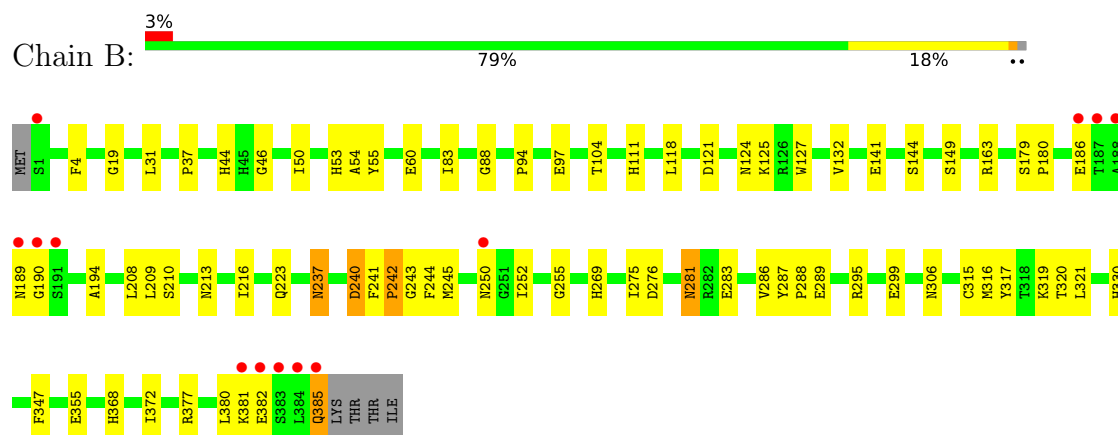
3 Residue-property plots [i](#)

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

• Molecule 1: MONOMERIC SARCOSINE OXIDASE



• 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	72.65Å 69.66Å 73.36Å 90.00° 93.74° 90.00°	Depositor
Resolution (Å)	32.40 – 1.71 32.40 – 1.71	Depositor EDS
% Data completeness (in resolution range)	92.0 (32.40-1.71) 92.6 (32.40-1.71)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.41 (at 1.71Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.172 , 0.208 0.172 , 0.208	Depositor DCC
R_{free} test set	3788 reflections (5.10%)	wwPDB-VP
Wilson B-factor (Å ²)	15.5	Xtriage
Anisotropy	0.511	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 41.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.025 for l,-k,h	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	6790	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.75% 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: GOL, FAD, CL

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	0.58	0/3081	1.05	18/4172 (0.4%)
1	B	0.58	0/3081	1.04	25/4172 (0.6%)
All	All	0.58	0/6162	1.05	43/8344 (0.5%)

There are no bond length outliers.

All (43) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	243	GLY	N-CA-C	-9.62	99.54	112.82
1	A	54	ALA	N-CA-C	-8.55	97.01	109.59
1	B	243	GLY	N-CA-C	-8.34	99.42	112.85
1	A	88	GLY	N-CA-C	-8.13	102.08	112.65
1	B	54	ALA	N-CA-C	-7.14	99.10	109.59
1	A	44	HIS	N-CA-C	6.96	121.48	112.92
1	A	149	SER	N-CA-C	6.96	119.75	111.33
1	B	149	SER	N-CA-C	6.60	118.47	111.28
1	A	269	HIS	N-CA-C	6.46	119.63	111.69
1	B	190	GLY	N-CA-C	6.38	118.55	110.96
1	B	250	ASN	N-CA-C	6.34	119.95	112.72
1	A	83	ILE	N-CA-C	6.28	116.96	111.56
1	B	88	GLY	N-CA-C	-6.26	104.51	112.65
1	B	377	ARG	CA-C-N	6.24	125.46	118.97
1	B	377	ARG	C-N-CA	6.24	125.46	118.97
1	B	55	TYR	N-CA-C	6.23	118.92	108.02
1	B	37	PRO	N-CA-C	6.15	118.20	110.70
1	B	83	ILE	N-CA-C	6.03	116.75	111.56
1	B	104	THR	N-CA-C	-5.69	105.16	111.36
1	A	320	THR	N-CA-C	-5.57	102.22	110.46
1	A	55	TYR	N-CA-C	5.55	117.81	107.99
1	A	301	TYR	N-CA-C	5.51	120.28	113.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	251	GLY	N-CA-C	5.49	118.31	110.46
1	A	29	THR	N-CA-C	5.48	118.33	109.40
1	B	330	HIS	N-CA-C	-5.45	102.72	109.64
1	B	269	HIS	N-CA-C	5.44	118.39	111.69
1	B	46	GLY	N-CA-C	-5.44	107.60	115.27
1	B	241	PHE	N-CA-C	-5.44	101.50	109.50
1	B	208	LEU	N-CA-C	5.42	120.56	113.30
1	B	320	THR	N-CA-C	-5.42	102.44	110.46
1	B	276	ASP	N-CA-C	-5.42	101.19	109.64
1	A	37	PRO	N-CA-C	5.41	117.30	110.70
1	A	237	ASN	N-CA-C	-5.36	105.55	111.71
1	B	44	HIS	N-CA-C	5.34	119.90	113.17
1	A	319	LYS	N-CA-C	5.34	118.52	109.76
1	B	240	ASP	N-CA-C	5.33	119.08	112.58
1	B	319	LYS	N-CA-C	5.31	118.47	109.76
1	B	237	ASN	N-CA-C	-5.26	105.66	111.71
1	B	50	ILE	CB-CA-C	-5.25	103.89	111.34
1	A	287	TYR	N-CA-C	-5.24	102.23	110.14
1	A	104	THR	N-CA-C	-5.10	105.63	111.14
1	B	316	MET	N-CA-C	5.04	117.66	109.24
1	A	178	ILE	N-CA-C	5.02	115.53	108.36

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	3006	0	2910	73	1
1	B	3006	0	2910	44	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	53	0	29	1	0
3	B	53	0	29	1	0
4	A	6	0	8	4	0
4	B	6	0	8	6	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	328	0	0	17	0
5	B	330	0	0	6	0
All	All	6790	0	5894	120	1

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 10.

All (120) 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:103:GLU:HG2	5:A:1225:HOH:O	1.42	1.16
1:A:213:ASN:ND2	1:A:385:GLN:HB2	1.84	0.93
1:B:237:ASN:O	4:B:902:GOL:H2	1.70	0.91
4:B:902:GOL:H12	5:B:1142:HOH:O	1.73	0.87
1:A:55:TYR:HB3	5:A:1226:HOH:O	1.76	0.85
1:B:213:ASN:ND2	1:B:385:GLN:HB2	1.93	0.84
1:A:237:ASN:O	4:A:901:GOL:O2	1.98	0.82
1:A:240:ASP:HA	4:A:901:GOL:O2	1.82	0.79
1:B:240:ASP:OD1	4:B:902:GOL:H11	1.84	0.78
1:A:382:GLU:HB3	5:A:1223:HOH:O	1.87	0.75
1:A:124:ASN:HD21	1:A:132:VAL:H	1.36	0.74
1:A:60:GLU:H	1:A:60:GLU:CD	1.96	0.73
1:A:1:SER:HA	5:A:1220:HOH:O	1.91	0.70
1:A:58:GLY:HA2	5:A:1225:HOH:O	1.90	0.70
1:A:58:GLY:N	5:A:1225:HOH:O	2.24	0.70
1:A:213:ASN:HD22	1:A:385:GLN:HB2	1.54	0.70
1:B:213:ASN:HD22	1:B:385:GLN:HB2	1.56	0.70
1:A:58:GLY:HA3	1:A:60:GLU:OE1	1.94	0.68
1:B:281:ASN:HD22	1:B:283:GLU:H	1.43	0.66
1:A:163:ARG:HG2	1:A:163:ARG:HH11	1.60	0.66
1:A:59:ARG:HA	5:A:1226:HOH:O	1.95	0.65
1:A:58:GLY:CA	5:A:1225:HOH:O	2.44	0.65
1:A:237:ASN:OD1	4:A:901:GOL:H2	1.98	0.63
1:B:295:ARG:HH11	1:B:306:ASN:ND2	1.97	0.61
1:A:380:LEU:O	1:A:381:LYS:HB2	2.01	0.61
1:A:53:HIS:HB2	1:A:141:GLU:OE2	2.00	0.61
1:A:124:ASN:HD21	1:A:132:VAL:HG22	1.66	0.61
1:A:244:PHE:CZ	1:A:255:GLY:HA3	2.38	0.59
1:B:121:ASP:OD1	1:B:125:LYS:HE2	2.03	0.59
1:B:163:ARG:NE	5:B:1185:HOH:O	2.36	0.58
4:B:902:GOL:H31	5:B:1250:HOH:O	2.03	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:213:ASN:HD21	1:A:385:GLN:HB2	1.66	0.57
1:B:124:ASN:HD21	1:B:132:VAL:H	1.52	0.57
1:B:281:ASN:HD22	1:B:281:ASN:C	2.13	0.56
1:A:281:ASN:C	1:A:281:ASN:HD22	2.14	0.56
1:B:60:GLU:HG3	1:B:372:ILE:HG21	1.87	0.56
1:B:288:PRO:HG2	1:B:289:GLU:OE2	2.05	0.56
1:B:124:ASN:HD21	1:B:132:VAL:HG22	1.71	0.56
1:B:94:PRO:HB2	1:B:97:GLU:HB2	1.87	0.55
1:A:163:ARG:NE	5:A:938:HOH:O	2.40	0.54
1:A:135:ASN:H	1:A:135:ASN:HD22	1.54	0.54
1:A:281:ASN:HD22	1:A:283:GLU:H	1.55	0.54
1:A:173:VAL:HA	1:A:187:THR:HG22	1.90	0.54
1:A:215:ASP:HB2	1:A:385:GLN:OE1	2.07	0.54
1:A:58:GLY:C	5:A:1226:HOH:O	2.51	0.54
1:A:77:LYS:HE3	5:A:1127:HOH:O	2.08	0.54
1:A:378:PRO:HD2	5:A:1005:HOH:O	2.07	0.53
1:A:359:GLN:HE22	1:A:367:GLU:HG2	1.73	0.53
1:B:281:ASN:ND2	1:B:283:GLU:H	2.07	0.53
1:B:141:GLU:OE1	1:B:144:SER:OG	2.25	0.52
1:B:295:ARG:O	1:B:299:GLU:HG2	2.09	0.52
1:A:163:ARG:HG2	1:A:163:ARG:NH1	2.24	0.52
1:A:281:ASN:ND2	1:A:283:GLU:H	2.09	0.51
1:A:95:LYS:HG2	1:A:137:ASN:HB3	1.93	0.51
1:A:59:ARG:O	1:A:62:VAL:HG23	2.11	0.50
1:A:179:SER:HB2	1:A:180:PRO:HD2	1.92	0.50
1:A:55:TYR:O	5:A:1226:HOH:O	2.20	0.50
1:B:380:LEU:O	1:B:381:LYS:HB2	2.11	0.50
1:B:210:SER:HA	1:B:385:GLN:HE22	1.78	0.49
1:B:355:GLU:OE2	1:B:368:HIS:HE1	1.95	0.49
1:A:295:ARG:HH11	1:A:306:ASN:ND2	2.10	0.49
1:B:286:VAL:HG23	1:B:287:TYR:CD2	2.47	0.49
1:B:179:SER:HB2	1:B:180:PRO:HD2	1.95	0.49
1:B:245:MET:HE3	1:B:252:ILE:HG21	1.94	0.49
1:A:94:PRO:HB2	1:A:97:GLU:HB2	1.95	0.48
1:B:237:ASN:ND2	4:B:902:GOL:O2	2.46	0.48
1:A:36:ASP:OD2	1:A:157:ARG:NH1	2.37	0.47
1:A:207:LYS:C	1:A:208:LEU:HD12	2.38	0.47
1:B:19:GLY:HA3	1:B:31:LEU:HD21	1.95	0.47
1:B:244:PHE:CZ	1:B:255:GLY:HA3	2.49	0.47
1:B:186:GLU:O	5:B:1159:HOH:O	2.20	0.47
1:B:189:ASN:HD22	1:B:189:ASN:N	2.13	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:4:PHE:O	1:B:194:ALA:HA	2.16	0.46
1:A:245:MET:HE3	1:A:252:ILE:HG21	1.98	0.46
1:B:275:ILE:C	1:B:275:ILE:HD12	2.40	0.46
1:A:372:ILE:HG12	5:A:1171:HOH:O	2.16	0.45
1:A:98:SER:OG	1:A:247:GLU:OE2	2.27	0.45
1:B:163:ARG:HG2	1:B:163:ARG:HH11	1.81	0.45
1:A:213:ASN:ND2	1:A:385:GLN:H	2.13	0.45
1:A:58:GLY:CA	1:A:60:GLU:OE1	2.62	0.45
1:B:53:HIS:HB2	1:B:141:GLU:OE2	2.17	0.45
1:B:321:LEU:CD1	1:B:381:LYS:HD3	2.47	0.45
1:B:240:ASP:HA	4:B:902:GOL:H11	1.99	0.44
1:B:163:ARG:CZ	5:B:1185:HOH:O	2.64	0.44
1:A:133:PRO:HB2	1:A:135:ASN:ND2	2.31	0.44
1:A:288:PRO:HG2	1:A:289:GLU:OE2	2.18	0.44
1:B:141:GLU:CD	1:B:144:SER:OG	2.61	0.44
1:A:199:VAL:CG1	1:A:201:MET:HE2	2.48	0.44
1:A:7:ILE:HG13	1:A:194:ALA:HB2	1.99	0.43
1:A:216:ILE:HG12	1:A:382:GLU:HG2	2.00	0.43
1:B:223:GLN:HG3	1:B:317:TYR:CG	2.53	0.43
1:B:216:ILE:HG12	1:B:382:GLU:HG2	2.00	0.43
1:A:124:ASN:ND2	1:A:132:VAL:HG22	2.30	0.43
1:A:239:ILE:HD12	1:A:239:ILE:HA	1.92	0.43
1:A:21:GLN:NE2	5:A:1013:HOH:O	2.52	0.43
1:A:51:ILE:O	1:A:144:SER:HB2	2.19	0.43
1:B:381:LYS:HB3	1:B:381:LYS:HE2	1.67	0.42
1:A:95:LYS:HG3	1:A:135:ASN:O	2.20	0.42
1:A:286:VAL:HG23	1:A:287:TYR:CD2	2.55	0.42
1:A:248:VAL:HB	1:A:249:PRO:CD	2.50	0.42
1:A:86:LYS:NZ	1:A:143:ASN:HD22	2.18	0.41
1:B:121:ASP:O	1:B:125:LYS:HG3	2.20	0.41
1:A:2:THR:HB	5:A:1163:HOH:O	2.20	0.41
1:A:135:ASN:H	1:A:135:ASN:ND2	2.18	0.41
1:B:111:HIS:HD2	5:B:1286:HOH:O	2.03	0.41
1:A:203:ALA:HA	1:A:318:THR:OG1	2.21	0.41
1:B:209:LEU:HD23	1:B:209:LEU:HA	1.94	0.41
1:A:227:PHE:CZ	1:A:265:LYS:HD2	2.56	0.41
1:B:347:PHE:HB2	3:B:400:FAD:H2'	2.03	0.41
1:A:268:TYR:CE2	1:A:271:PHE:HB3	2.56	0.41
1:A:60:GLU:CG	1:A:372:ILE:HD13	2.51	0.40
1:A:124:ASN:ND2	1:A:132:VAL:H	2.12	0.40
1:A:163:ARG:NH1	1:A:163:ARG:CG	2.84	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:181:ASP:OD1	1:A:181:ASP:N	2.54	0.40
1:A:289:GLU:H	1:A:289:GLU:CD	2.29	0.40
1:B:127:TRP:CD1	1:B:242:PRO:HD3	2.56	0.40
1:A:118:LEU:N	1:A:118:LEU:HD23	2.36	0.40
1:A:204:TRP:CE3	3:A:400:FAD:H8A	2.57	0.40
1:A:325:PHE:CD1	1:A:342:PHE:HA	2.57	0.40
4:A:901:GOL:H2	5:A:973:HOH:O	2.22	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:78:GLU:O	4:B:902:GOL:O2[1_455]	2.07	0.13

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	383/390 (98%)	370 (97%)	12 (3%)	1 (0%)	36	24
1	B	383/390 (98%)	372 (97%)	11 (3%)	0	100	100
All	All	766/780 (98%)	742 (97%)	23 (3%)	1 (0%)	48	34

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	381	LYS

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	322/327 (98%)	316 (98%)	6 (2%)	50	27
1	B	322/327 (98%)	317 (98%)	5 (2%)	55	34
All	All	644/654 (98%)	633 (98%)	11 (2%)	53	31

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	118	LEU
1	A	242	PRO
1	A	281	ASN
1	A	315	CYS
1	A	381	LYS
1	A	385	GLN
1	B	118	LEU
1	B	242	PRO
1	B	281	ASN
1	B	315	CYS
1	B	385	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	21	GLN
1	A	41	ASN
1	A	124	ASN
1	A	135	ASN
1	A	143	ASN
1	A	189	ASN
1	A	205	ASN
1	A	213	ASN
1	A	219	GLN
1	A	250	ASN
1	A	281	ASN
1	A	293	ASN
1	A	306	ASN
1	A	359	GLN
1	A	368	HIS

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Mol	Chain	Res	Type
1	B	111	HIS
1	B	124	ASN
1	B	135	ASN
1	B	189	ASN
1	B	205	ASN
1	B	213	ASN
1	B	219	GLN
1	B	281	ASN
1	B	293	ASN
1	B	306	ASN
1	B	359	GLN
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 [i](#)

Of 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 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	$\# Z > 2$
4	GOL	A	901	-	5,5,5	0.84	0	5,5,5	0.63	0
3	FAD	A	400	1	58,58,58	2.39	12 (20%)	85,89,89	1.57	13 (15%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	FAD	B	400	1	58,58,58	2.49	16 (27%)	85,89,89	1.57	17 (20%)
4	GOL	B	902	-	5,5,5	0.47	0	5,5,5	0.94	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
4	GOL	A	901	-	-	2/4/4/4	-
3	FAD	A	400	1	-	1/34/50/50	0/6/6/6
3	FAD	B	400	1	-	1/34/50/50	0/6/6/6
4	GOL	B	902	-	-	0/4/4/4	-

All (28) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	400	FAD	C4'-C3'	9.44	1.69	1.53
3	B	400	FAD	C4'-C3'	9.12	1.69	1.53
3	A	400	FAD	P-O3P	7.08	1.67	1.59
3	A	400	FAD	PA-O3P	6.98	1.67	1.59
3	B	400	FAD	P-O3P	6.94	1.67	1.59
3	B	400	FAD	PA-O3P	6.91	1.67	1.59
3	B	400	FAD	C8M-C8	-5.20	1.41	1.51
3	A	400	FAD	C8M-C8	-4.67	1.42	1.51
3	B	400	FAD	C4X-N5	4.39	1.40	1.30
3	A	400	FAD	C4X-N5	4.13	1.39	1.30
3	A	400	FAD	O4'-C4'	-3.71	1.35	1.43
3	B	400	FAD	C6-C5X	3.64	1.45	1.40
3	B	400	FAD	C4-N3	3.48	1.45	1.38
3	B	400	FAD	O4'-C4'	-3.35	1.36	1.43
3	B	400	FAD	C2A-N3A	2.97	1.39	1.33
3	A	400	FAD	C6-C5X	2.90	1.44	1.40
3	A	400	FAD	O2'-C2'	2.77	1.49	1.43
3	B	400	FAD	C5'-C4'	-2.73	1.48	1.51
3	B	400	FAD	C10-N10	2.65	1.43	1.37
3	A	400	FAD	C2A-N3A	2.51	1.38	1.33
3	A	400	FAD	C9A-N10	2.39	1.45	1.41
3	B	400	FAD	C9-C8	-2.38	1.36	1.39
3	B	400	FAD	C1'-N10	2.11	1.52	1.47
3	B	400	FAD	C5B-C4B	-2.08	1.45	1.51
3	A	400	FAD	C10-N10	2.08	1.41	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	400	FAD	P-O2P	-2.07	1.45	1.55
3	B	400	FAD	O2'-C2'	2.04	1.47	1.43
3	A	400	FAD	C4-N3	2.01	1.42	1.38

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	400	FAD	C4'-C3'-C2'	-5.30	104.75	113.57
3	A	400	FAD	O4B-C1B-N9A	5.20	118.07	108.09
3	B	400	FAD	C4'-C3'-C2'	-5.06	105.14	113.57
3	B	400	FAD	C5'-C4'-C3'	-4.90	102.98	112.22
3	A	400	FAD	C5'-C4'-C3'	-4.88	103.01	112.22
3	B	400	FAD	O4B-C1B-N9A	4.68	117.08	108.09
3	A	400	FAD	O4'-C4'-C5'	3.75	118.26	109.99
3	B	400	FAD	O4'-C4'-C5'	3.55	117.81	109.99
3	B	400	FAD	C8M-C8-C9	-3.22	113.89	119.57
3	B	400	FAD	O3P-P-O1P	-3.18	101.12	110.70
3	A	400	FAD	C8M-C8-C9	-3.13	114.06	119.57
3	A	400	FAD	O3P-P-O1P	-3.02	101.63	110.70
3	A	400	FAD	O2P-P-O1P	2.92	126.03	112.44
3	B	400	FAD	O2P-P-O1P	2.91	125.97	112.44
3	A	400	FAD	O2A-PA-O3P	-2.78	99.75	107.27
3	A	400	FAD	O4'-C4'-C3'	2.61	115.37	109.25
3	B	400	FAD	O2A-PA-O3P	-2.57	100.33	107.27
3	B	400	FAD	O4'-C4'-C3'	2.54	115.20	109.25
3	B	400	FAD	C1'-C2'-C3'	-2.34	103.30	109.66
3	A	400	FAD	O2A-PA-O1A	2.31	123.19	112.44
3	B	400	FAD	N3A-C2A-N1A	-2.28	125.13	128.58
3	B	400	FAD	O2A-PA-O1A	2.26	122.96	112.44
3	A	400	FAD	C5X-N5-C4X	-2.26	114.43	118.09
3	A	400	FAD	C1'-C2'-C3'	-2.22	103.64	109.66
3	A	400	FAD	N3A-C2A-N1A	-2.19	125.27	128.58
3	B	400	FAD	C9-C8-C7	2.14	122.83	119.69
3	B	400	FAD	O2B-C2B-C1B	2.13	117.45	110.10
3	B	400	FAD	C2B-C1B-N9A	2.11	118.54	113.30
3	B	400	FAD	O5'-C5'-C4'	-2.07	103.84	109.36
3	B	400	FAD	C4-C4X-N5	-2.01	115.44	118.21

There are no chirality outliers.

All (4) torsion outliers are listed below:

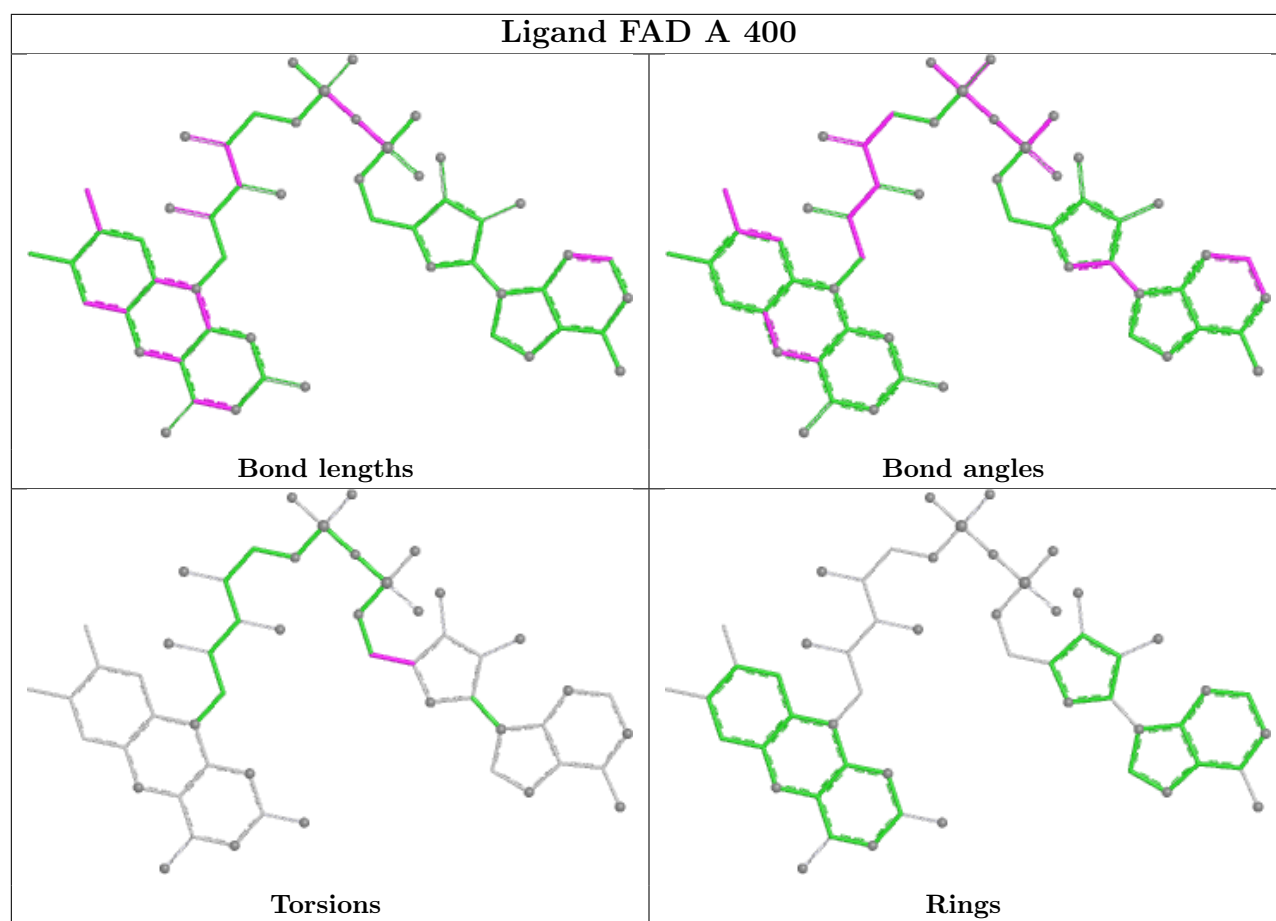
Mol	Chain	Res	Type	Atoms
4	A	901	GOL	C1-C2-C3-O3
4	A	901	GOL	O2-C2-C3-O3
3	B	400	FAD	O4B-C4B-C5B-O5B
3	A	400	FAD	O4B-C4B-C5B-O5B

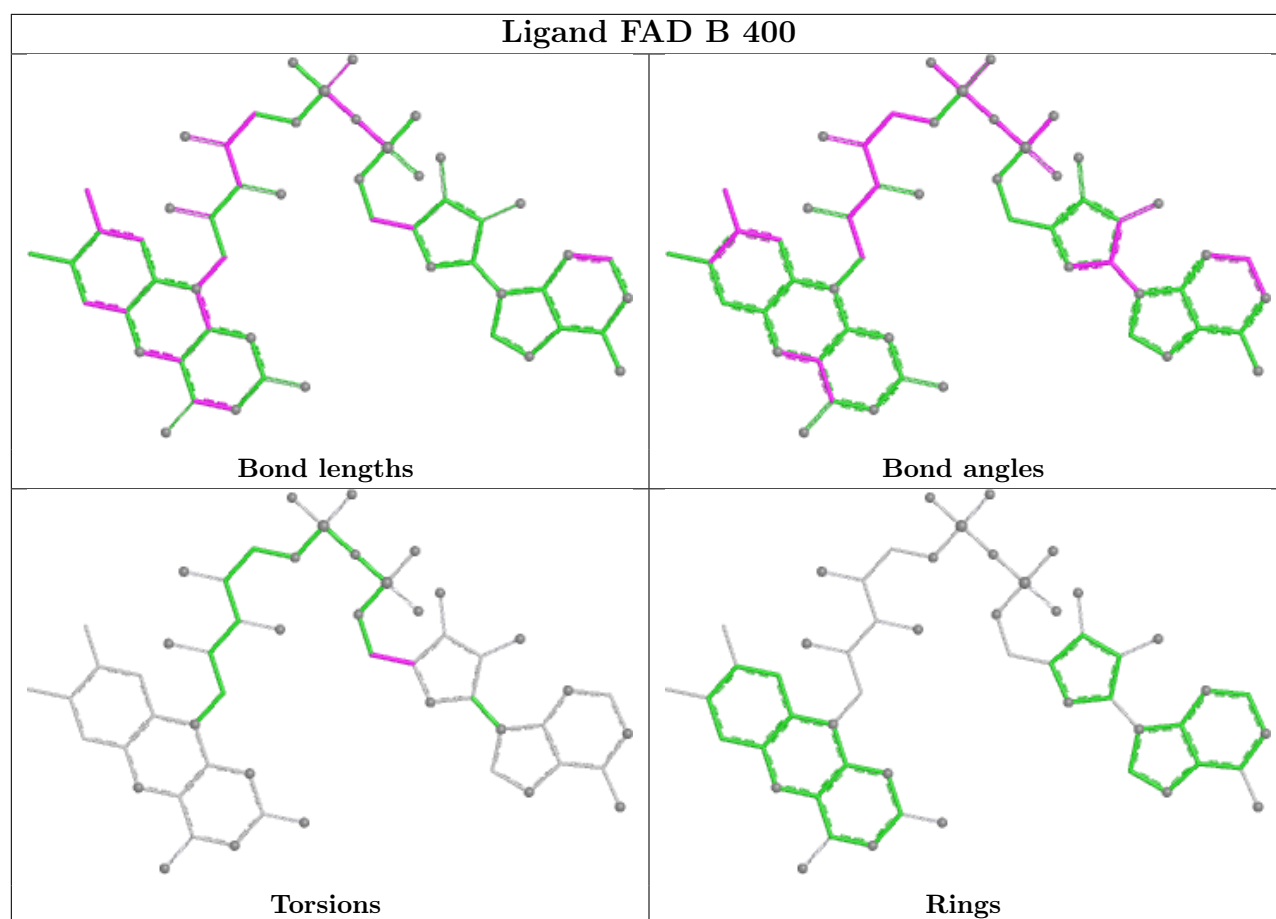
There are no ring outliers.

4 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	901	GOL	4	0
3	A	400	FAD	1	0
3	B	400	FAD	1	0
4	B	902	GOL	6	1

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	385/390 (98%)	-0.09	12 (3%)	51	61	11, 19, 33, 71	0
1	B	385/390 (98%)	-0.06	13 (3%)	48	57	10, 20, 35, 71	0
All	All	770/780 (98%)	-0.07	25 (3%)	50	59	10, 20, 34, 71	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	382	GLU	5.5
1	A	385	GLN	4.6
1	A	384	LEU	4.0
1	B	384	LEU	4.0
1	B	385	GLN	4.0
1	A	383	SER	3.8
1	B	381	LYS	3.8
1	A	381	LYS	3.5
1	A	2	THR	3.5
1	B	383	SER	3.5
1	B	382	GLU	3.2
1	B	191	SER	3.0
1	A	60	GLU	2.7
1	B	190	GLY	2.6
1	A	380	LEU	2.6
1	B	189	ASN	2.6
1	A	181	ASP	2.5
1	B	188	ALA	2.5
1	B	250	ASN	2.4
1	A	1	SER	2.4
1	B	1	SER	2.3
1	B	187	THR	2.3
1	A	250	ASN	2.3
1	A	3	HIS	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	186	GLU	2.1

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](#)

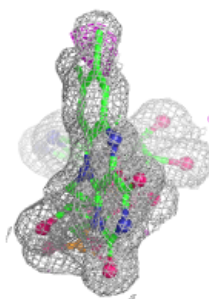
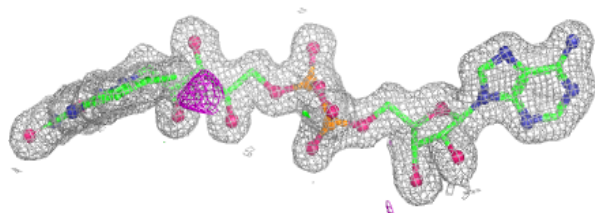
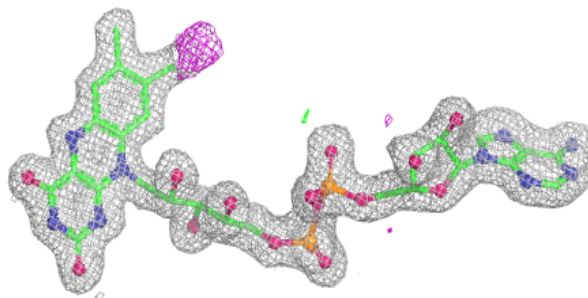
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	GOL	A	901	6/6	0.74	0.18	30,34,39,40	0
4	GOL	B	902	6/6	0.74	0.21	20,29,38,40	0
3	FAD	A	400	53/53	0.98	0.05	11,14,17,18	0
3	FAD	B	400	53/53	0.98	0.04	12,14,17,21	0
2	CL	A	401	1/1	0.99	0.02	16,16,16,16	0
2	CL	B	401	1/1	0.99	0.03	16,16,16,16	0

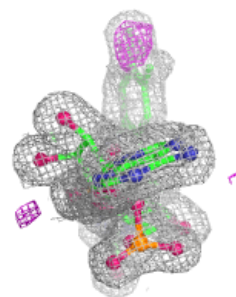
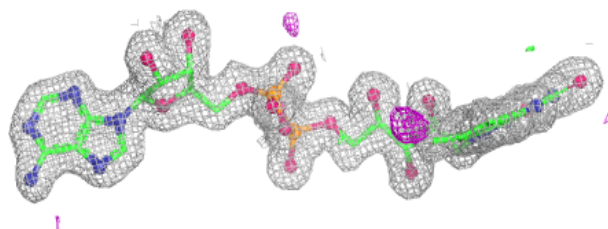
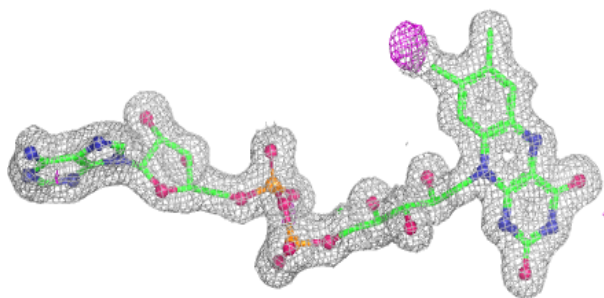
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around FAD A 400:

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

**Electron density around FAD B 400:**

$2mF_o-DF_c$ (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.