



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 15, 2026 – 12:38 PM UTC

PDB ID : 6FCX / pdb_00006fcx
Title : Structure of human 5,10-methylenetetrahydrofolate reductase (MTHFR)
Authors : Kopec, J.; Bezerra, G.A.; Oberholzer, A.E.; Rembeza, E.; Sorrell, F.J.; Chalk, R.; Borkowska, O.; Ellis, K.; Kupinska, K.; Krojer, T.; Burgess-Brown, N.; Von Delft, F.; Arrowsmith, C.; Edwards, A.; Bountra, C.; Froese, D.S.; Baumgartner, M.; Yue, W.W.; Structural Genomics Consortium (SGC)
Deposited on : 2017-12-21
Resolution : 2.50 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](#) ①) 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

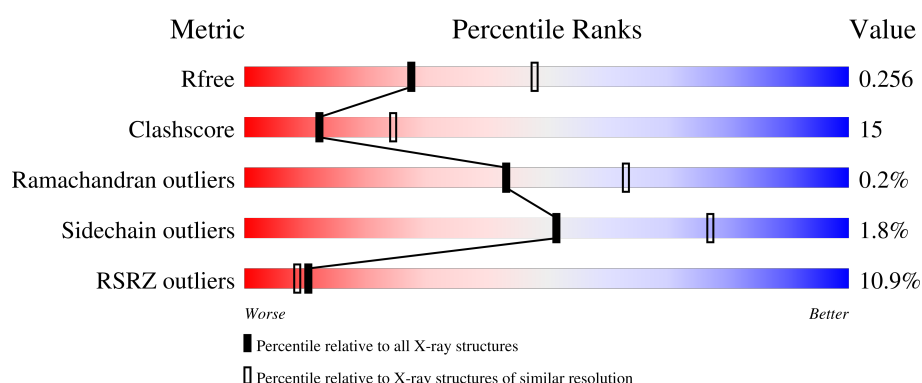
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	615	3% 78% 17% ..
1	B	615	18% 73% 22% ..

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 9180 atoms, of which 10 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.

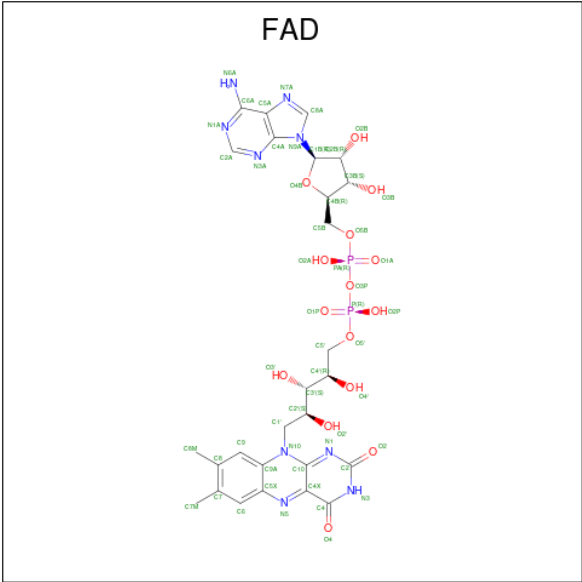
- Molecule 1 is a protein called Methylenetetrahydrofolate reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	596	Total	C	N	O	S	0	0	0
			4723	3048	786	870	19			
1	B	590	Total	C	N	O	S	0	0	0
			4191	2660	727	791	13			

There are 20 discrepancies between the modelled and reference sequences:

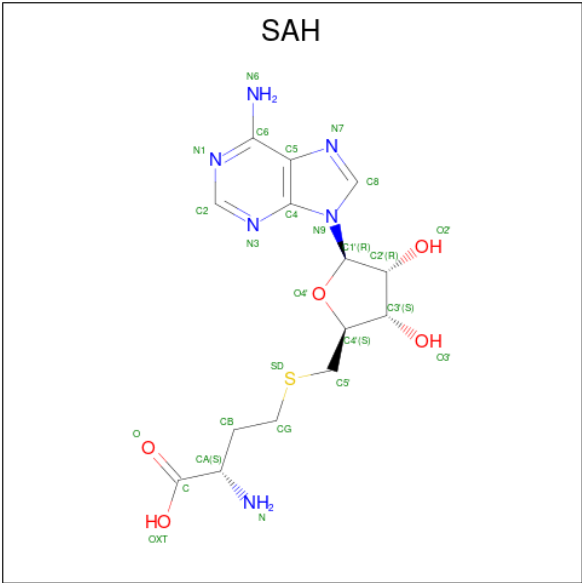
Chain	Residue	Modelled	Actual	Comment	Reference
A	37	MET	LEU	initiating methionine	UNP P42898
A	429	ALA	GLU	variant	UNP P42898
A	594	GLN	ARG	variant	UNP P42898
A	645	ALA	-	expression tag	UNP P42898
A	646	GLU	-	expression tag	UNP P42898
A	647	ASN	-	expression tag	UNP P42898
A	648	LEU	-	expression tag	UNP P42898
A	649	TYR	-	expression tag	UNP P42898
A	650	PHE	-	expression tag	UNP P42898
A	651	GLN	-	expression tag	UNP P42898
B	37	MET	LEU	initiating methionine	UNP P42898
B	429	ALA	GLU	variant	UNP P42898
B	594	GLN	ARG	variant	UNP P42898
B	645	ALA	-	expression tag	UNP P42898
B	646	GLU	-	expression tag	UNP P42898
B	647	ASN	-	expression tag	UNP P42898
B	648	LEU	-	expression tag	UNP P42898
B	649	TYR	-	expression tag	UNP P42898
B	650	PHE	-	expression tag	UNP P42898
B	651	GLN	-	expression tag	UNP P42898

- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (CCD ID: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).



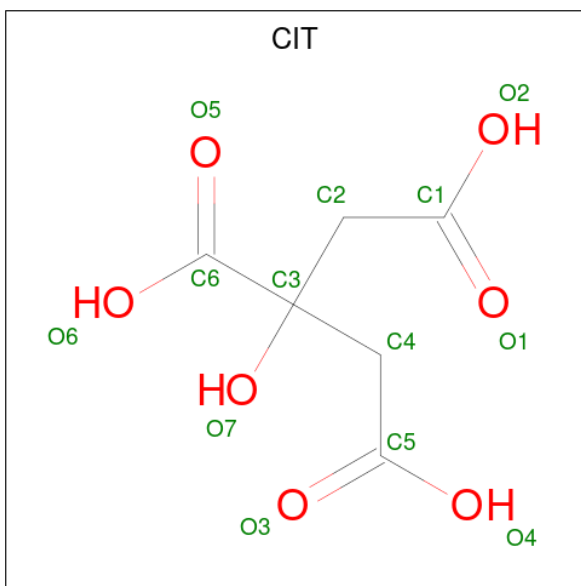
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
2	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 3 is S-ADENOSYL-L-HOMOCYSTEINE (CCD ID: SAH) (formula: C₁₄H₂₀N₆O₅S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			26	14	6	5	1		
3	B	1	Total	C	N	O	S	0	0
			26	14	6	5	1		

- Molecule 4 is CITRIC ACID (CCD ID: CIT) (formula: $C_6H_8O_7$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	H	O	0	0
			18	6	5	7		
4	B	1	Total	C	H	O	0	0
			18	6	5	7		

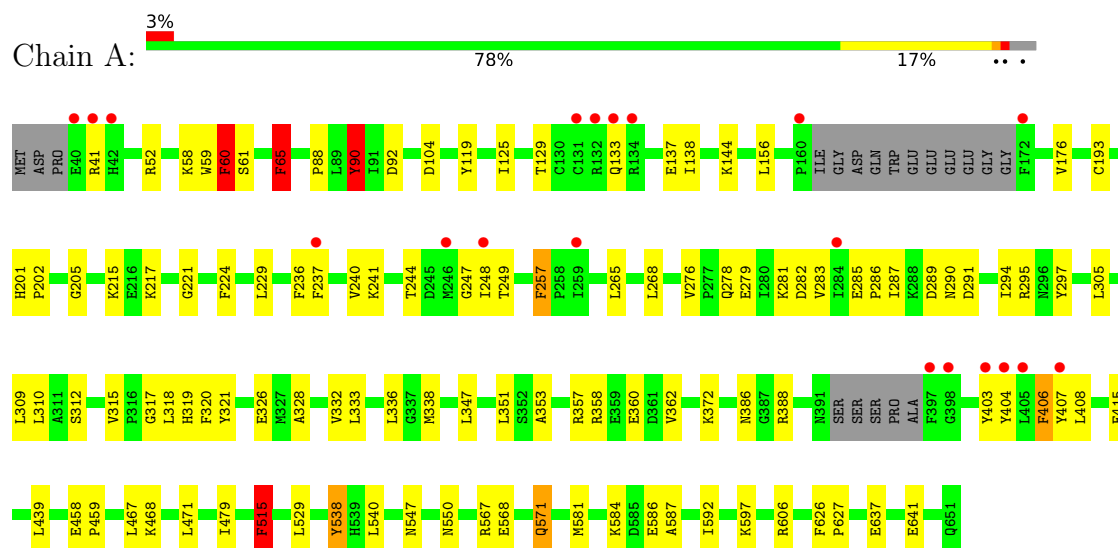
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	48	Total	O	0	0
			48	48		
5	B	24	Total	O	0	0
			24	24		

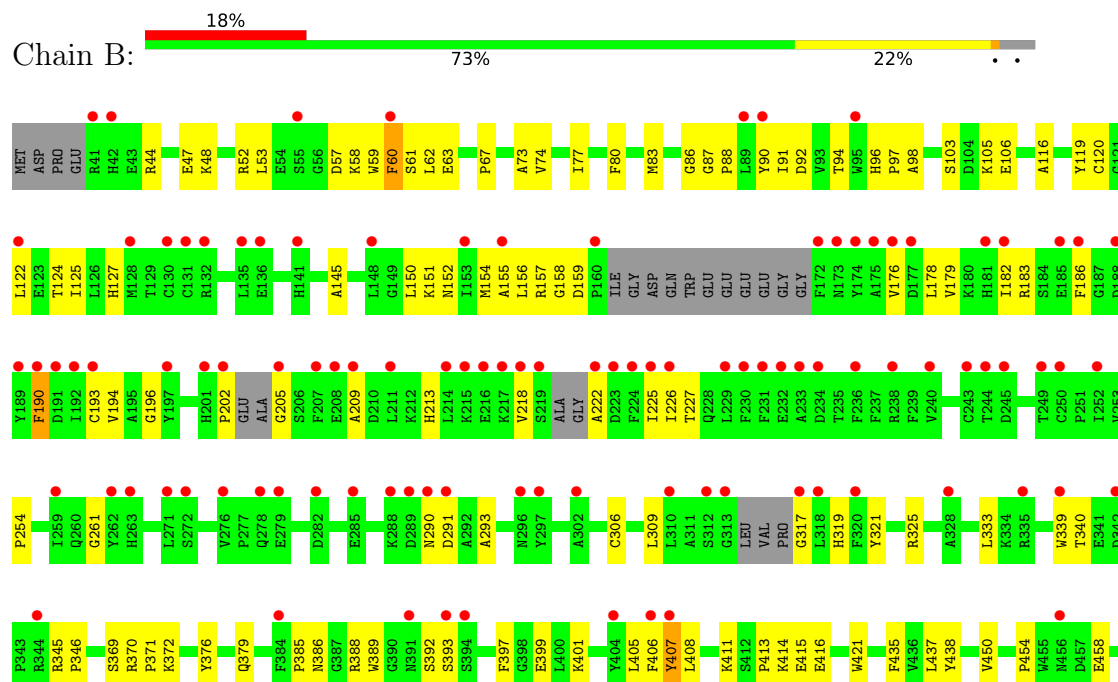
3 Residue-property plots

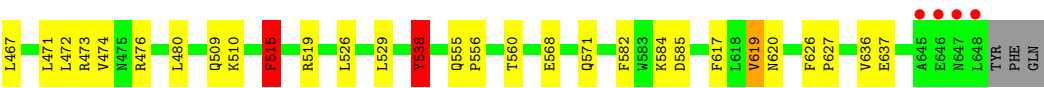
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Methylenetetrahydrofolate reductase



• Molecule 1: Methylenetetrahydrofolate reductase





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	97.34Å 127.94Å 147.01Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	63.97 – 2.50 63.97 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.9 (63.97-2.50) 100.0 (63.97-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.44 (at 2.51Å)	Xtriage
Refinement program	PHENIX (1.13_2998: ???)	Depositor
R, R_{free}	0.210 , 0.247 0.222 , 0.256	Depositor DCC
R_{free} test set	3266 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å ²)	69.2	Xtriage
Anisotropy	0.359	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 61.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	9180	wwPDB-VP
Average B, all atoms (Å ²)	92.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.73% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CIT, FAD, SAH

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	0.61	0/4853	0.81	3/6606 (0.0%)
1	B	0.49	0/4295	0.69	0/5885
All	All	0.56	0/9148	0.75	3/12491 (0.0%)

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	5
1	B	0	2
All	All	0	7

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	606	ARG	CA-CB-CG	-5.72	102.66	114.10
1	A	571	GLN	CB-CG-CD	5.41	121.80	112.60
1	A	90	TYR	CA-CB-CG	5.20	123.27	113.90

There are no chirality outliers.

All (7) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	404	TYR	Peptide
1	A	515	PHE	Sidechain
1	A	60	PHE	Sidechain

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Mol	Chain	Res	Type	Group
1	A	65	PHE	Sidechain
1	A	90	TYR	Sidechain
1	B	515	PHE	Sidechain
1	B	538	TYR	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	4723	0	4516	128	0
1	B	4191	0	3577	147	0
2	A	53	0	31	2	0
2	B	53	0	30	2	0
3	A	26	0	19	0	0
3	B	26	0	17	0	0
4	A	13	5	5	1	0
4	B	13	5	5	1	0
5	A	48	0	0	0	0
5	B	24	0	0	3	0
All	All	9170	10	8200	266	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 15.

All (266) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:62:LEU:HB2	1:B:91:ILE:HD11	1.31	1.11
1:B:62:LEU:HB2	1:B:91:ILE:CD1	1.81	1.09
1:A:529:LEU:HD13	1:A:538:TYR:CD2	1.92	1.05
1:A:439:LEU:O	1:A:468:LYS:NZ	1.93	0.99
1:B:157:ARG:HB3	2:B:701:FAD:H5'1	1.44	0.99
1:B:411:LYS:NZ	1:B:458:GLU:OE2	1.95	0.98
1:B:61:SER:HB3	1:B:319:HIS:HA	1.48	0.94
1:B:529:LEU:HD13	1:B:538:TYR:CD2	2.03	0.93
1:A:571:GLN:O	1:A:571:GLN:HG3	1.67	0.92

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:529:LEU:HD13	1:A:538:TYR:HD2	1.29	0.92
1:A:276:VAL:HG13	1:A:281:LYS:HD2	1.53	0.91
1:A:567:ARG:HG2	1:B:555:GLN:HE21	1.35	0.90
1:B:154:MET:HE2	1:B:156:LEU:HD11	1.52	0.89
1:A:92:ASP:OD1	1:A:125:ILE:HB	1.72	0.87
1:A:357:ARG:HD3	1:A:360:GLU:OE2	1.75	0.86
1:A:215:LYS:HA	1:A:248:ILE:HD11	1.58	0.84
1:A:567:ARG:HG2	1:B:555:GLN:NE2	1.93	0.82
1:A:129:THR:HG22	1:A:156:LEU:HB2	1.61	0.82
1:B:159:ASP:OD1	2:B:701:FAD:O4'	2.00	0.79
1:B:405:LEU:HD13	1:B:407:TYR:HE1	1.48	0.79
1:B:190:PHE:HD1	1:B:190:PHE:O	1.65	0.78
1:B:154:MET:HG2	1:B:156:LEU:CD1	2.14	0.77
1:B:571:GLN:O	1:B:571:GLN:HG3	1.83	0.77
1:A:332:VAL:O	1:A:336:LEU:HD13	1.84	0.77
1:A:318:LEU:HB3	1:A:320:PHE:HE1	1.50	0.76
1:B:62:LEU:CD1	1:B:83:MET:HE2	2.16	0.75
1:A:201:HIS:CD2	1:A:202:PRO:HD2	2.22	0.75
1:B:529:LEU:HD13	1:B:538:TYR:HD2	1.52	0.74
1:B:62:LEU:HD13	1:B:83:MET:HE2	1.72	0.72
1:B:53:LEU:HG	1:B:59:TRP:CZ3	2.24	0.72
1:A:283:VAL:HG21	1:A:297:TYR:CZ	2.25	0.71
1:A:215:LYS:HA	1:A:248:ILE:CD1	2.20	0.71
1:B:62:LEU:CB	1:B:91:ILE:HD11	2.15	0.71
1:B:538:TYR:O	1:B:538:TYR:HD1	1.74	0.71
1:B:218:VAL:O	1:B:222:ALA:HB3	1.91	0.70
1:A:129:THR:HG22	1:A:156:LEU:CB	2.22	0.70
1:A:626:PHE:CG	1:A:627:PRO:HD3	2.27	0.69
1:A:290:ASN:O	1:A:294:ILE:HD12	1.91	0.69
1:B:515:PHE:CZ	1:B:619:VAL:HG21	2.28	0.69
1:B:476:ARG:NH1	5:B:801:HOH:O	2.25	0.68
1:B:435:PHE:CE2	1:B:480:LEU:HB3	2.27	0.68
1:A:156:LEU:HD13	2:A:701:FAD:C4X	2.24	0.67
1:B:438:TYR:HA	1:B:450:VAL:HG23	1.76	0.67
1:A:318:LEU:HB3	1:A:320:PHE:CE1	2.29	0.67
1:A:406:PHE:HE1	1:A:407:TYR:CE2	2.15	0.65
1:B:346:PRO:HD2	5:B:806:HOH:O	1.97	0.64
1:B:405:LEU:HD13	1:B:407:TYR:CE1	2.31	0.64
1:B:154:MET:CE	1:B:156:LEU:HD11	2.26	0.64
1:B:155:ALA:C	1:B:156:LEU:HD12	2.22	0.64
1:A:58:LYS:HG3	1:A:310:LEU:HD12	1.79	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:568:GLU:OE2	1:B:556:PRO:HG2	1.98	0.64
1:B:145:ALA:O	1:B:150:LEU:HB2	1.99	0.63
1:B:125:ILE:HA	1:B:152:ASN:O	1.98	0.63
1:B:60:PHE:HZ	1:B:333:LEU:CB	2.12	0.63
1:A:41:ARG:O	1:A:41:ARG:HG3	1.99	0.62
1:A:58:LYS:HE3	1:A:315:VAL:O	2.00	0.62
1:B:413:PRO:HG2	1:B:416:GLU:HG2	1.80	0.62
1:B:90:TYR:HD2	1:B:125:ILE:CD1	2.12	0.62
1:B:472:LEU:O	1:B:476:ARG:HG3	2.00	0.62
1:B:155:ALA:O	1:B:156:LEU:HD12	2.00	0.62
1:A:41:ARG:NH1	1:A:52:ARG:HH21	1.97	0.61
1:A:236:PHE:O	1:A:240:VAL:HG23	1.99	0.61
1:B:92:ASP:OD2	1:B:127:HIS:NE2	2.33	0.61
1:A:283:VAL:HG21	1:A:297:TYR:CE2	2.35	0.61
1:A:60:PHE:CD1	1:A:60:PHE:C	2.78	0.61
1:B:626:PHE:CG	1:B:627:PRO:HD3	2.35	0.61
1:B:154:MET:HG2	1:B:156:LEU:HD11	1.83	0.61
1:B:438:TYR:N	1:B:450:VAL:HG21	2.16	0.60
1:B:209:ALA:O	1:B:213:HIS:N	2.32	0.60
1:B:190:PHE:O	1:B:190:PHE:CD1	2.52	0.60
1:A:279:GLU:OE1	1:A:279:GLU:N	2.26	0.60
1:B:87:GLY:N	1:B:88:PRO:HD3	2.17	0.60
1:A:41:ARG:HH12	1:A:52:ARG:NE	1.99	0.60
1:A:295:ARG:HD2	1:A:328:ALA:HB2	1.83	0.59
1:A:637:GLU:O	1:A:641:GLU:HG3	2.02	0.59
1:B:196:GLY:O	1:B:227:THR:HA	2.02	0.59
1:A:119:TYR:CE2	1:A:372:LYS:HE3	2.38	0.59
1:A:241:LYS:HA	1:A:244:THR:OG1	2.03	0.59
1:A:317:GLY:O	1:A:318:LEU:HD12	2.03	0.59
1:B:397:PHE:HB2	1:B:582:PHE:CD2	2.38	0.58
1:B:405:LEU:HD12	1:B:408:LEU:HD12	1.85	0.58
1:A:41:ARG:HH12	1:A:52:ARG:CZ	2.16	0.58
1:A:215:LYS:CA	1:A:248:ILE:HD11	2.30	0.58
1:A:388:ARG:NH2	1:B:386:ASN:OD1	2.36	0.58
1:B:53:LEU:HA	1:B:57:ASP:O	2.04	0.58
1:B:190:PHE:CD1	1:B:190:PHE:C	2.82	0.58
1:B:151:LYS:O	1:B:190:PHE:CD1	2.57	0.58
1:B:157:ARG:HD3	1:B:158:GLY:N	2.19	0.58
1:A:279:GLU:H	1:A:279:GLU:CD	2.12	0.57
1:A:65:PHE:CD1	1:A:65:PHE:C	2.83	0.57
1:A:406:PHE:CE1	1:A:407:TYR:CE2	2.92	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:376:TYR:O	1:B:379:GLN:HG2	2.05	0.56
1:B:154:MET:HG3	1:B:193:CYS:O	2.05	0.56
1:B:62:LEU:HD11	1:B:83:MET:HE2	1.86	0.56
1:B:119:TYR:CE2	1:B:407:TYR:HE2	2.24	0.56
1:A:247:GLY:O	1:A:249:THR:HG23	2.06	0.56
1:B:48:LYS:O	1:B:52:ARG:NE	2.39	0.56
1:A:41:ARG:NH1	1:A:52:ARG:NH2	2.54	0.56
1:B:60:PHE:CZ	1:B:333:LEU:CB	2.88	0.55
1:B:119:TYR:CE2	1:B:407:TYR:CE2	2.94	0.55
1:B:106:GLU:OE1	1:B:406:PHE:HB3	2.07	0.55
1:A:52:ARG:HH11	1:A:59:TRP:HB3	1.71	0.54
1:B:182:ILE:O	1:B:186:PHE:N	2.40	0.54
1:B:155:ALA:HB1	1:B:178:LEU:CB	2.37	0.54
1:B:399:GLU:OE1	1:B:401:LYS:HE3	2.08	0.54
1:B:510:LYS:NZ	5:B:802:HOH:O	2.41	0.54
1:B:48:LYS:O	1:B:52:ARG:CD	2.56	0.54
1:B:405:LEU:HB3	1:B:407:TYR:CE1	2.42	0.54
1:A:276:VAL:O	1:A:281:LYS:HD2	2.08	0.53
1:B:90:TYR:CD2	1:B:125:ILE:CD1	2.92	0.53
1:B:306:CYS:HA	1:B:309:LEU:CB	2.38	0.53
1:A:281:LYS:O	1:A:285:GLU:HG2	2.08	0.53
1:A:309:LEU:O	1:A:312:SER:HB3	2.08	0.53
1:A:41:ARG:HH12	1:A:52:ARG:NH2	2.05	0.53
1:B:450:VAL:HG23	1:B:450:VAL:O	2.09	0.53
1:A:265:LEU:HA	1:A:268:LEU:CD2	2.39	0.52
1:A:439:LEU:O	1:A:468:LYS:CE	2.58	0.52
1:B:515:PHE:CD1	1:B:515:PHE:C	2.87	0.52
1:B:515:PHE:CZ	1:B:619:VAL:CG2	2.93	0.52
1:B:59:TRP:CD1	1:B:317:GLY:HA3	2.44	0.52
1:B:62:LEU:HD13	1:B:83:MET:CE	2.39	0.52
1:A:229:LEU:C	1:A:229:LEU:HD12	2.35	0.52
1:B:96:HIS:CD2	1:B:98:ALA:HB3	2.45	0.51
1:B:438:TYR:HA	1:B:450:VAL:CG2	2.40	0.51
1:A:581:MET:O	1:A:584:LYS:HE3	2.09	0.51
1:A:133:GLN:HB3	1:A:138:ILE:HD11	1.93	0.51
1:B:369:SER:C	1:B:371:PRO:HD3	2.34	0.51
1:A:60:PHE:C	1:A:60:PHE:HD1	2.17	0.51
1:B:62:LEU:HB2	1:B:91:ILE:HD12	1.82	0.51
1:A:133:GLN:HA	1:A:137:GLU:OE1	2.10	0.51
1:B:196:GLY:N	1:B:226:ILE:O	2.39	0.50
1:A:92:ASP:OD2	1:A:319:HIS:HE1	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:91:ILE:HG13	1:B:92:ASP:H	1.77	0.50
1:B:538:TYR:CD1	1:B:538:TYR:C	2.90	0.50
1:B:67:PRO:HG3	1:B:73:ALA:HA	1.93	0.49
1:B:538:TYR:HD1	1:B:538:TYR:C	2.20	0.49
1:A:265:LEU:O	1:A:268:LEU:HD23	2.12	0.49
1:B:421:TRP:CZ2	1:B:454:PRO:HD3	2.48	0.49
1:A:278:GLN:O	1:A:282:ASP:N	2.36	0.49
1:B:145:ALA:C	1:B:150:LEU:HB2	2.38	0.49
1:A:65:PHE:C	1:A:65:PHE:HD1	2.20	0.48
1:A:567:ARG:NH1	1:B:555:GLN:HE22	2.11	0.48
1:A:626:PHE:CD2	1:A:627:PRO:HD3	2.47	0.48
1:B:86:GLY:HA3	1:B:339:TRP:NE1	2.28	0.48
1:A:236:PHE:CE2	1:A:315:VAL:HG12	2.48	0.48
1:A:279:GLU:O	1:A:283:VAL:HG13	2.12	0.48
1:A:538:TYR:HD1	1:A:538:TYR:O	1.97	0.48
1:A:538:TYR:HE1	1:A:550:ASN:N	2.11	0.48
1:A:406:PHE:CD1	1:A:406:PHE:C	2.90	0.48
1:A:133:GLN:CB	1:A:138:ILE:HD11	2.44	0.48
1:A:479:ILE:HG23	1:A:515:PHE:CG	2.49	0.48
1:A:61:SER:HB2	1:A:90:TYR:CD1	2.49	0.48
1:A:567:ARG:HH11	1:B:555:GLN:HE22	1.61	0.48
1:B:96:HIS:HD2	1:B:98:ALA:HB3	1.79	0.47
1:B:407:TYR:HD1	1:B:407:TYR:H	1.62	0.47
1:A:119:TYR:CD2	1:A:372:LYS:HE3	2.49	0.47
1:B:176:VAL:O	1:B:178:LEU:N	2.39	0.47
1:A:257:PHE:CD1	1:A:257:PHE:C	2.92	0.47
1:A:467:LEU:CD1	1:A:471:LEU:CD1	2.92	0.47
1:A:406:PHE:CE1	1:A:407:TYR:CD2	3.02	0.47
1:A:257:PHE:C	1:A:257:PHE:HD1	2.23	0.46
1:B:370:ARG:N	1:B:371:PRO:HD3	2.30	0.46
1:B:386:ASN:OD1	1:B:386:ASN:C	2.58	0.46
1:A:276:VAL:HG22	1:A:281:LYS:HG3	1.96	0.46
1:B:74:VAL:HG12	1:B:406:PHE:O	2.15	0.46
1:B:435:PHE:CD2	1:B:480:LEU:HB3	2.50	0.46
1:A:317:GLY:C	1:A:318:LEU:HD12	2.41	0.46
1:B:90:TYR:HD2	1:B:125:ILE:HD13	1.78	0.46
1:A:41:ARG:HH12	1:A:52:ARG:HE	1.63	0.46
1:B:94:THR:HA	1:B:127:HIS:ND1	2.31	0.46
1:A:257:PHE:HB2	1:A:321:TYR:HD2	1.80	0.46
1:A:467:LEU:HD13	1:A:471:LEU:HG	1.97	0.46
1:B:584:LYS:HG3	1:B:585:ASP:N	2.31	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:283:VAL:CG2	1:A:297:TYR:CE2	3.00	0.45
1:B:202:PRO:HA	1:B:205:GLY:N	2.32	0.45
1:A:353:ALA:HA	1:A:358:ARG:NH1	2.31	0.45
1:A:60:PHE:CE1	1:A:88:PRO:HB3	2.52	0.45
1:A:201:HIS:CG	1:A:202:PRO:HD2	2.49	0.45
1:A:347:LEU:HD12	1:A:351:LEU:HD21	1.98	0.45
1:A:406:PHE:HE1	1:A:407:TYR:CZ	2.35	0.45
1:A:586:GLU:O	1:A:587:ALA:C	2.60	0.45
1:B:421:TRP:CH2	1:B:454:PRO:HD3	2.52	0.45
1:A:406:PHE:HE1	1:A:407:TYR:CD2	2.34	0.45
1:B:59:TRP:O	1:B:317:GLY:HA2	2.17	0.45
1:B:405:LEU:CD1	1:B:407:TYR:HE1	2.24	0.45
1:A:60:PHE:HB3	1:A:338:MET:HE2	1.99	0.45
1:A:305:LEU:C	1:A:305:LEU:HD23	2.42	0.45
1:B:405:LEU:HB3	1:B:407:TYR:CD1	2.52	0.44
1:B:467:LEU:HD12	1:B:471:LEU:CD1	2.46	0.44
1:A:415:GLU:H	1:A:415:GLU:CD	2.25	0.44
1:B:473:ARG:NE	1:B:637:GLU:OE2	2.48	0.44
1:A:538:TYR:CE1	1:A:550:ASN:N	2.85	0.44
1:B:73:ALA:O	1:B:77:ILE:HG12	2.16	0.44
1:B:515:PHE:C	1:B:515:PHE:HD1	2.25	0.44
1:A:156:LEU:HD13	2:A:701:FAD:C4	2.47	0.44
1:A:592:ILE:O	1:A:597:LYS:HG2	2.18	0.44
1:B:53:LEU:HD23	1:B:57:ASP:O	2.17	0.44
1:B:59:TRP:NE1	1:B:317:GLY:HA3	2.33	0.44
1:B:120:CYS:HB3	1:B:122:LEU:HD23	2.00	0.44
1:B:254:PRO:HD2	1:B:317:GLY:O	2.18	0.44
1:B:509:GLN:HG2	1:B:560:THR:HG23	2.00	0.44
1:B:519:ARG:HG3	1:B:617:PHE:HE1	1.83	0.44
1:A:290:ASN:C	1:A:294:ILE:HD12	2.41	0.43
1:A:467:LEU:C	1:A:467:LEU:HD12	2.43	0.43
1:A:305:LEU:HD23	1:A:309:LEU:HG	2.00	0.43
1:A:626:PHE:CD1	1:A:627:PRO:HD3	2.53	0.43
1:B:155:ALA:CB	1:B:178:LEU:CB	2.96	0.43
1:B:340:THR:HG23	1:B:340:THR:O	2.18	0.43
1:A:515:PHE:CD1	1:A:515:PHE:C	2.96	0.43
1:A:236:PHE:HE2	1:A:315:VAL:HG12	1.83	0.43
1:A:278:GLN:HA	1:A:281:LYS:HB2	1.99	0.43
1:A:538:TYR:CE1	1:A:550:ASN:HB3	2.52	0.43
1:B:119:TYR:O	1:B:372:LYS:HE3	2.18	0.43
1:B:183:ARG:HA	1:B:186:PHE:CB	2.48	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:467:LEU:CD1	1:A:471:LEU:HD11	2.49	0.43
1:A:291:ASP:HA	1:A:294:ILE:HD13	2.01	0.43
1:B:261:GLY:HA2	1:B:291:ASP:CB	2.49	0.43
1:B:437:LEU:C	1:B:450:VAL:HG21	2.44	0.43
1:B:194:VAL:O	1:B:225:ILE:HA	2.19	0.43
1:A:104:ASP:OD1	1:A:144:LYS:NZ	2.49	0.43
1:B:44:ARG:CB	1:B:47:GLU:OE1	2.66	0.43
1:B:91:ILE:HG22	1:B:124:THR:HG22	2.01	0.43
1:B:345:ARG:HA	1:B:346:PRO:HD3	1.92	0.42
1:B:106:GLU:HB2	1:B:406:PHE:CD1	2.54	0.42
1:A:458:GLU:HB3	1:A:459:PRO:CD	2.50	0.42
1:B:96:HIS:ND1	1:B:97:PRO:HD2	2.34	0.42
1:B:157:ARG:HD3	1:B:158:GLY:C	2.44	0.42
1:A:276:VAL:HG13	1:A:281:LYS:CD	2.36	0.42
1:A:403:TYR:HE2	1:A:408:LEU:HD11	1.84	0.42
1:A:571:GLN:NE2	1:B:568:GLU:OE1	2.48	0.42
1:B:399:GLU:OE1	1:B:401:LYS:CE	2.66	0.42
1:A:568:GLU:OE1	1:B:571:GLN:NE2	2.51	0.42
1:B:474:VAL:HG12	1:B:636:VAL:HG11	2.02	0.42
1:B:179:VAL:O	1:B:183:ARG:CB	2.68	0.42
1:B:90:TYR:CD2	1:B:125:ILE:HD13	2.54	0.42
1:B:290:ASN:O	1:B:293:ALA:N	2.53	0.42
1:A:90:TYR:CD1	1:A:90:TYR:C	2.98	0.41
1:B:63:GLU:OE1	1:B:321:TYR:CE1	2.73	0.41
1:B:77:ILE:O	1:B:80:PHE:HB2	2.20	0.41
1:A:467:LEU:HD11	1:A:471:LEU:HD11	2.01	0.41
1:A:540:LEU:O	1:A:547:ASN:HA	2.20	0.41
1:A:529:LEU:CD1	1:A:538:TYR:HD2	2.16	0.41
1:B:103:SER:OG	1:B:105:LYS:CG	2.69	0.41
1:A:265:LEU:HA	1:A:265:LEU:HD23	1.71	0.41
1:B:91:ILE:HD12	1:B:91:ILE:HA	1.77	0.41
1:B:414:LYS:HE3	1:B:415:GLU:OE2	2.21	0.41
1:A:237:PHE:CZ	1:A:312:SER:HB2	2.56	0.41
1:B:392:SER:O	1:B:393:SER:C	2.63	0.41
1:A:65:PHE:CD1	1:A:65:PHE:O	2.74	0.41
1:A:386:ASN:OD1	1:A:386:ASN:C	2.64	0.41
1:B:60:PHE:HD1	1:B:60:PHE:HA	1.61	0.41
1:B:385:PRO:HG2	1:B:389:TRP:CZ3	2.56	0.41
1:A:282:ASP:O	1:A:286:PRO:CD	2.69	0.41
1:A:287:ILE:O	1:A:289:ASP:N	2.54	0.41
1:A:386:ASN:OD1	1:B:388:ARG:NH2	2.48	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:116:ALA:HB3	1:B:124:THR:HG21	2.02	0.41
1:A:176:VAL:HG23	1:A:217:LYS:HA	2.03	0.41
1:A:326:GLU:N	4:A:703:CIT:O3	2.45	0.41
1:B:120:CYS:O	1:B:122:LEU:HD22	2.21	0.41
1:B:154:MET:CG	1:B:156:LEU:HD11	2.50	0.41
1:A:193:CYS:HB3	1:A:224:PHE:CE1	2.55	0.40
1:A:41:ARG:O	1:A:41:ARG:CG	2.69	0.40
1:B:53:LEU:CD2	1:B:58:LYS:HA	2.51	0.40
1:B:325:ARG:HB3	4:B:703:CIT:O6	2.21	0.40
1:A:60:PHE:CE2	1:A:333:LEU:HD13	2.56	0.40
1:B:526:LEU:HD23	1:B:526:LEU:HA	1.87	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	590/615 (96%)	570 (97%)	18 (3%)	2 (0%)	36	55
1	B	580/615 (94%)	548 (94%)	32 (6%)	0	100	100
All	All	1170/1230 (95%)	1118 (96%)	50 (4%)	2 (0%)	43	63

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	205	GLY
1	A	221	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	491/537 (91%)	483 (98%)	8 (2%)	55	79
1	B	358/537 (67%)	351 (98%)	7 (2%)	48	75
All	All	849/1074 (79%)	834 (98%)	15 (2%)	51	77

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	60	PHE
1	A	65	PHE
1	A	90	TYR
1	A	257	PHE
1	A	362	VAL
1	A	406	PHE
1	A	515	PHE
1	A	538	TYR
1	B	60	PHE
1	B	190	PHE
1	B	407	TYR
1	B	515	PHE
1	B	538	TYR
1	B	619	VAL
1	B	620	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	391	ASN
1	A	509	GLN
1	A	557	ASN
1	B	96	HIS
1	B	147	GLN
1	B	448	HIS
1	B	509	GLN

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Mol	Chain	Res	Type
1	B	555	GLN
1	B	557	ASN
1	B	594	GLN
1	B	615	ASN

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](#)

6 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	$\# Z > 2$
4	CIT	A	703	-	12,12,12	1.09	0	17,17,17	1.65	3 (17%)
2	FAD	B	701	-	58,58,58	0.47	1 (1%)	85,89,89	0.52	0
4	CIT	B	703	-	12,12,12	1.03	0	17,17,17	1.41	1 (5%)
3	SAH	A	702	-	27,28,28	1.27	5 (18%)	36,40,40	1.99	11 (30%)
2	FAD	A	701	-	58,58,58	0.37	0	85,89,89	0.55	0
3	SAH	B	702	-	27,28,28	1.16	3 (11%)	36,40,40	2.01	12 (33%)

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	CIT	A	703	-	-	8/16/16/16	-
2	FAD	B	701	-	-	17/34/50/50	0/6/6/6
4	CIT	B	703	-	-	7/16/16/16	-
3	SAH	A	702	-	-	1/15/31/31	0/3/3/3
2	FAD	A	701	-	-	4/34/50/50	0/6/6/6
3	SAH	B	702	-	-	8/15/31/31	0/3/3/3

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	702	SAH	C2-N1	3.13	1.39	1.33
2	B	701	FAD	P-O3P	2.82	1.62	1.59
3	A	702	SAH	C2-N1	2.72	1.38	1.33
3	A	702	SAH	C2-N3	2.67	1.38	1.33
3	B	702	SAH	C5-N7	-2.64	1.34	1.39
3	A	702	SAH	C8-N7	2.57	1.36	1.31
3	B	702	SAH	C2-N3	2.53	1.38	1.33
3	A	702	SAH	OXT-C	-2.53	1.22	1.30
3	A	702	SAH	C5-N7	-2.11	1.35	1.39

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	702	SAH	N3-C2-N1	-6.35	118.97	128.58
3	B	702	SAH	N3-C2-N1	-4.84	121.26	128.58
3	B	702	SAH	N9-C8-N7	-4.28	107.86	113.94
3	A	702	SAH	C5-C4-N3	-3.72	121.60	126.72
3	B	702	SAH	C5-N7-C8	3.71	109.28	103.45
4	A	703	CIT	O6-C6-C3	3.58	120.01	113.14
3	A	702	SAH	C2-N3-C4	3.57	120.54	111.83
3	B	702	SAH	C5-C4-N3	-3.52	121.87	126.72
4	B	703	CIT	O6-C6-C3	3.28	119.44	113.14
3	A	702	SAH	C5'-SD-CG	-3.03	93.27	102.26
3	B	702	SAH	C6-C5-C4	2.87	121.09	117.18
3	A	702	SAH	OXT-C-O	-2.69	117.97	124.08
3	B	702	SAH	C2-N3-C4	2.62	118.22	111.83
3	A	702	SAH	N9-C8-N7	-2.60	110.25	113.94
3	A	702	SAH	N3-C4-N9	2.59	131.58	127.17

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	702	SAH	O4'-C1'-N9	2.59	113.06	108.09
4	A	703	CIT	O4-C5-C4	2.52	122.32	114.35
3	A	702	SAH	C5-C6-N6	-2.49	117.11	123.29
3	B	702	SAH	O4'-C4'-C5'	2.41	115.05	108.83
3	A	702	SAH	O3'-C3'-C4'	-2.35	104.33	111.08
3	B	702	SAH	C4'-C5'-SD	-2.34	105.43	113.78
3	A	702	SAH	N6-C6-N1	2.29	123.49	118.38
3	B	702	SAH	C5'-SD-CG	-2.25	95.57	102.26
3	B	702	SAH	N3-C4-N9	2.22	130.94	127.17
4	A	703	CIT	O7-C3-C2	-2.20	104.36	109.38
3	B	702	SAH	C4-C5-N7	-2.13	108.14	110.58
3	B	702	SAH	OXT-C-O	-2.08	119.36	124.08

There are no chirality outliers.

All (45) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	701	FAD	C1'-C2'-C3'-O3'
2	A	701	FAD	C1'-C2'-C3'-C4'
2	B	701	FAD	C5B-O5B-PA-O1A
2	B	701	FAD	C5B-O5B-PA-O3P
2	B	701	FAD	C1'-C2'-C3'-C4'
2	B	701	FAD	C3'-C4'-C5'-O5'
2	B	701	FAD	O4'-C4'-C5'-O5'
2	B	701	FAD	C5'-O5'-P-O2P
2	B	701	FAD	C5'-O5'-P-O3P
4	A	703	CIT	C1-C2-C3-O7
4	A	703	CIT	O7-C3-C6-O5
4	B	703	CIT	O7-C3-C6-O5
4	B	703	CIT	O7-C3-C6-O6
4	B	703	CIT	C4-C3-C6-O5
4	B	703	CIT	C4-C3-C6-O6
2	A	701	FAD	O2'-C2'-C3'-C4'
4	A	703	CIT	C1-C2-C3-C4
4	A	703	CIT	C1-C2-C3-C6
3	B	702	SAH	CA-CB-CG-SD
4	A	703	CIT	C4-C3-C6-O5
2	B	701	FAD	O3'-C3'-C4'-O4'
2	B	701	FAD	O2'-C2'-C3'-C4'
2	B	701	FAD	C2'-C3'-C4'-O4'
2	B	701	FAD	C2'-C3'-C4'-C5'
4	B	703	CIT	C6-C3-C4-C5

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Mol	Chain	Res	Type	Atoms
4	A	703	CIT	O7-C3-C6-O6
4	A	703	CIT	C4-C3-C6-O6
2	A	701	FAD	O2'-C2'-C3'-O3'
4	B	703	CIT	C1-C2-C3-O7
2	B	701	FAD	C1'-C2'-C3'-O3'
3	B	702	SAH	C2'-C1'-N9-C8
2	B	701	FAD	C5B-O5B-PA-O2A
2	B	701	FAD	C5'-O5'-P-O1P
4	B	703	CIT	C2-C3-C4-C5
3	B	702	SAH	C2'-C1'-N9-C4
4	A	703	CIT	C2-C3-C6-O6
3	B	702	SAH	OXT-C-CA-CB
3	A	702	SAH	C2'-C1'-N9-C8
2	B	701	FAD	O2'-C2'-C3'-O3'
2	B	701	FAD	O3'-C3'-C4'-C5'
3	B	702	SAH	O-C-CA-CB
3	B	702	SAH	O4'-C1'-N9-C8
2	B	701	FAD	C2B-C1B-N9A-C8A
3	B	702	SAH	OXT-C-CA-N
3	B	702	SAH	O-C-CA-N

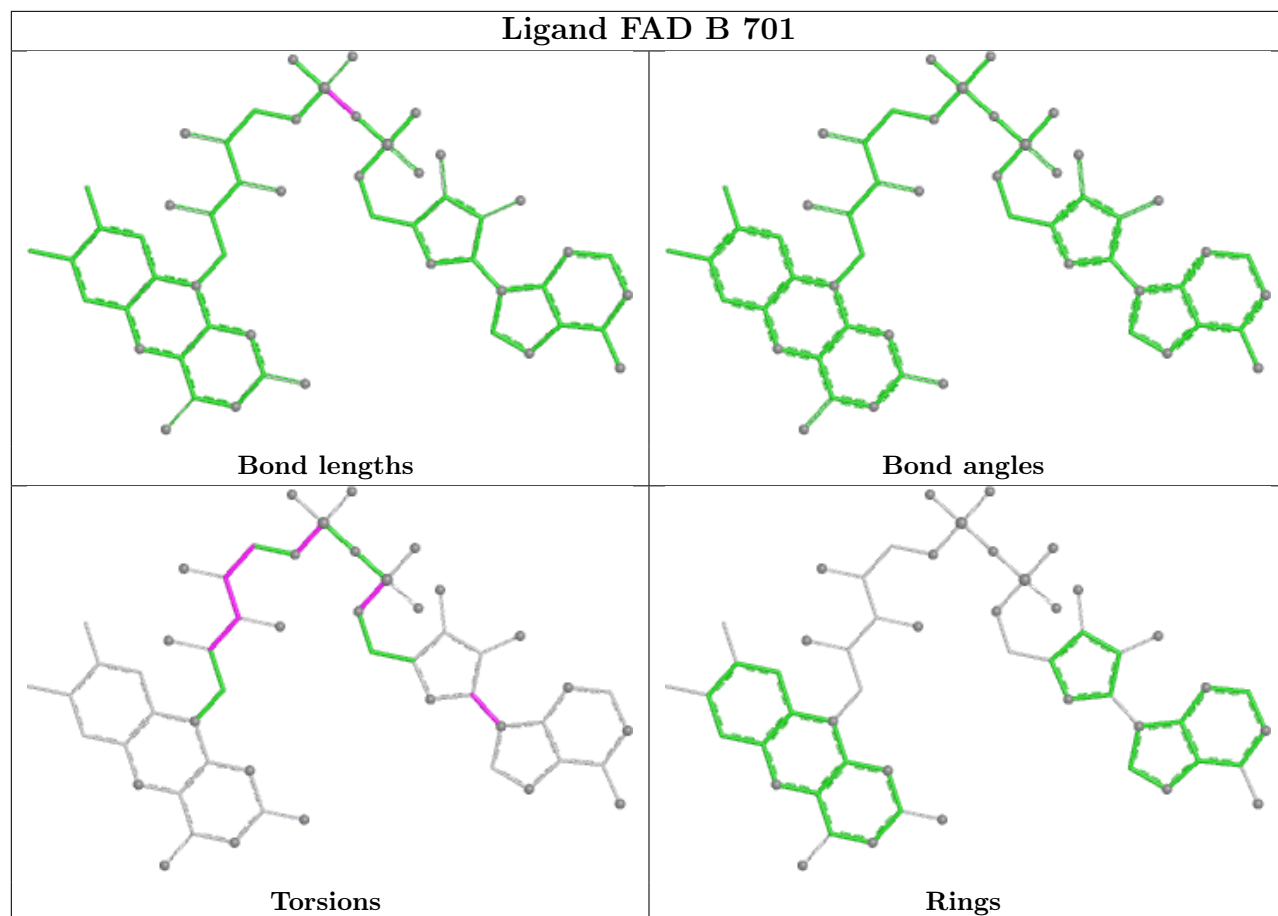
There are no ring outliers.

4 monomers are involved in 6 short contacts:

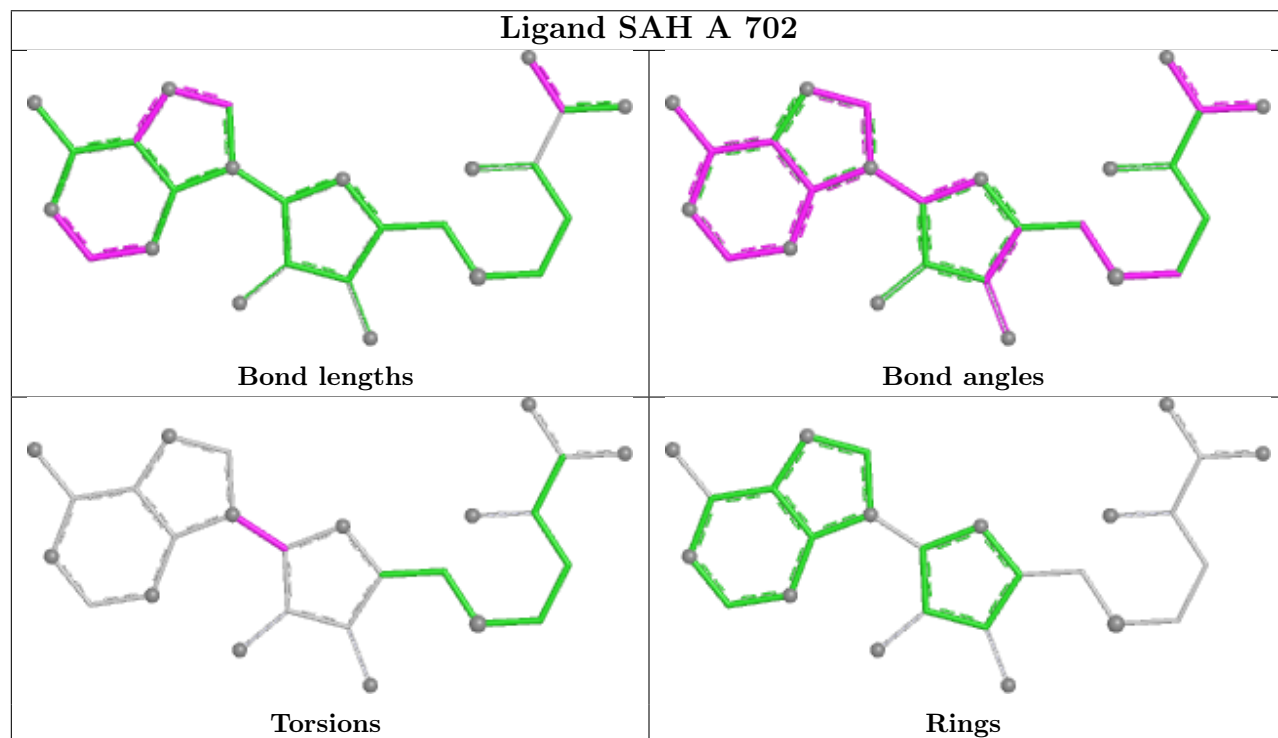
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	703	CIT	1	0
2	B	701	FAD	2	0
4	B	703	CIT	1	0
2	A	701	FAD	2	0

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.

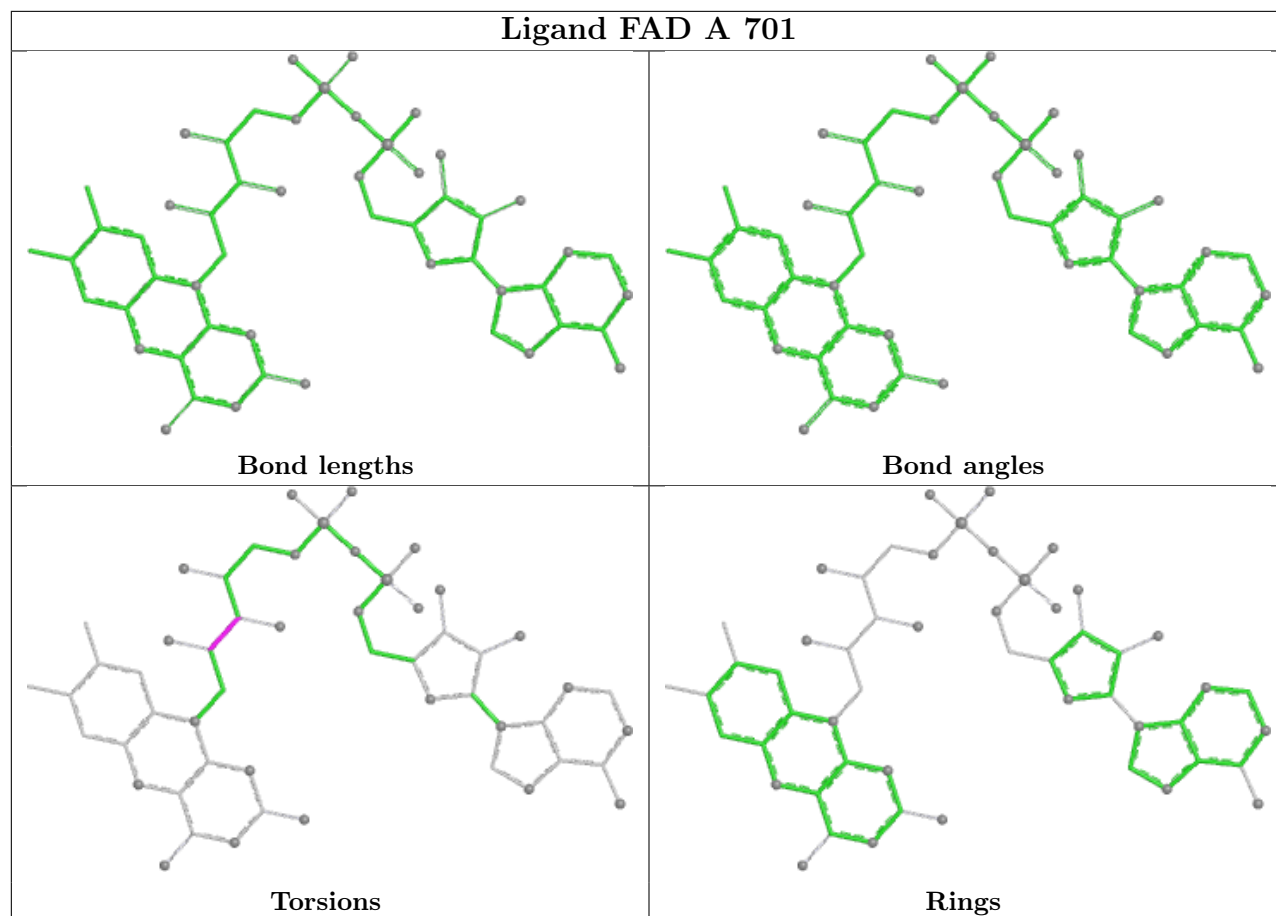
Ligand FAD B 701



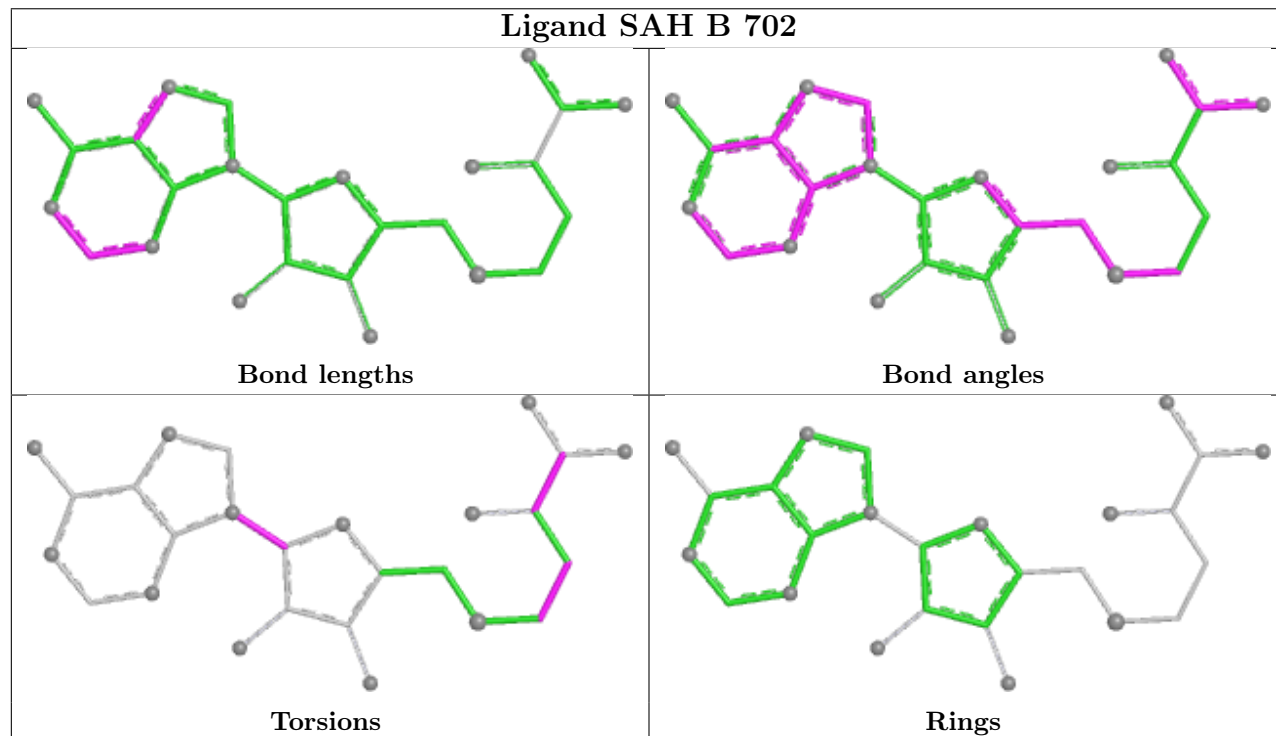
Ligand SAH A 702



Ligand FAD A 701



Ligand SAH B 702



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	596/615 (96%)	0.11	20 (3%) 48 43	48, 74, 131, 160	0
1	B	590/615 (95%)	0.86	109 (18%) 3 2	52, 97, 194, 241	0
All	All	1186/1230 (96%)	0.48	129 (10%) 10 9	48, 83, 181, 241	0

All (129) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	397	PHE	6.2
1	B	646	GLU	5.3
1	A	131	CYS	5.3
1	A	160	PRO	5.1
1	B	172	PHE	5.0
1	B	130	CYS	4.8
1	B	202	PRO	4.8
1	B	648	LEU	4.8
1	A	40	GLU	4.8
1	B	645	ALA	4.7
1	B	647	ASN	4.7
1	B	214	LEU	4.7
1	A	403	TYR	4.5
1	B	174	TYR	4.5
1	B	290	ASN	4.5
1	B	288	LYS	4.3
1	A	404	TYR	4.2
1	B	160	PRO	4.1
1	A	398	GLY	4.0
1	B	250	CYS	4.0
1	B	191	ASP	3.9
1	B	192	ILE	3.9
1	A	405	LEU	3.9
1	B	262	TYR	3.8

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Mol	Chain	Res	Type	RSRZ
1	B	205	GLY	3.8
1	B	42	HIS	3.7
1	B	155	ALA	3.7
1	B	217	LYS	3.6
1	B	289	ASP	3.6
1	B	176	VAL	3.5
1	A	42	HIS	3.5
1	B	240	VAL	3.4
1	B	189	TYR	3.4
1	B	320	PHE	3.3
1	B	131	CYS	3.3
1	B	197	TYR	3.3
1	B	278	GLN	3.3
1	B	128	MET	3.3
1	B	234	ASP	3.2
1	B	208	GLU	3.2
1	B	229	LEU	3.2
1	A	246	MET	3.2
1	B	407	TYR	3.2
1	A	132	ARG	3.1
1	B	209	ALA	3.1
1	B	219	SER	3.1
1	B	177	ASP	3.1
1	B	41	ARG	3.0
1	B	89	LEU	3.0
1	B	224	PHE	2.9
1	B	252	ILE	2.9
1	B	233	ALA	2.9
1	A	237	PHE	2.9
1	B	182	ILE	2.8
1	B	218	VAL	2.8
1	B	271	LEU	2.8
1	B	272	SER	2.8
1	B	296	ASN	2.8
1	B	185	GLU	2.8
1	B	456	ASN	2.7
1	B	249	THR	2.7
1	B	230	PHE	2.7
1	B	231	PHE	2.7
1	B	95	TRP	2.7
1	B	245	ASP	2.6
1	B	317	GLY	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	285	GLU	2.6
1	B	313	GLY	2.6
1	A	248	ILE	2.6
1	B	259	ILE	2.6
1	B	238	ARG	2.6
1	A	133	GLN	2.6
1	B	342	ASP	2.6
1	B	312	SER	2.6
1	B	215	LYS	2.6
1	B	181	HIS	2.6
1	B	393	SER	2.6
1	B	282	ASP	2.5
1	B	216	GLU	2.5
1	B	135	LEU	2.5
1	B	190	PHE	2.5
1	B	328	ALA	2.5
1	B	207	PHE	2.5
1	B	236	PHE	2.5
1	B	406	PHE	2.5
1	A	134	ARG	2.5
1	B	193	CYS	2.4
1	B	226	ILE	2.4
1	B	223	ASP	2.4
1	B	279	GLU	2.4
1	B	188	ASP	2.4
1	B	344	ARG	2.4
1	B	404	TYR	2.4
1	B	391	ASN	2.4
1	B	175	ALA	2.4
1	B	148	LEU	2.3
1	B	297	TYR	2.3
1	B	384	PHE	2.3
1	A	284	ILE	2.3
1	B	141	HIS	2.3
1	A	407	TYR	2.3
1	B	132	ARG	2.3
1	B	211	LEU	2.3
1	B	244	THR	2.2
1	B	232	GLU	2.2
1	B	243	CYS	2.2
1	B	201	HIS	2.2
1	B	339	TRP	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	41	ARG	2.2
1	B	310	LEU	2.2
1	B	153	ILE	2.2
1	B	222	ALA	2.2
1	A	172	PHE	2.2
1	B	335	ARG	2.1
1	B	60	PHE	2.1
1	B	186	PHE	2.1
1	B	90	TYR	2.1
1	B	394	SER	2.1
1	B	173	ASN	2.1
1	B	318	LEU	2.1
1	B	136	GLU	2.1
1	B	291	ASP	2.1
1	B	55	SER	2.1
1	B	276	VAL	2.0
1	B	225	ILE	2.0
1	B	122	LEU	2.0
1	B	302	ALA	2.0
1	A	259	ILE	2.0
1	B	263	HIS	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 [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($\text{\AA}^2$)	Q<0.9
4	CIT	B	703	13/13	0.54	0.11	144,145,175,175	0
2	FAD	B	701	53/53	0.78	0.13	102,125,139,142	0

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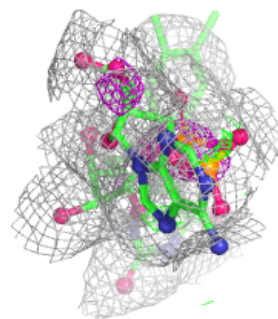
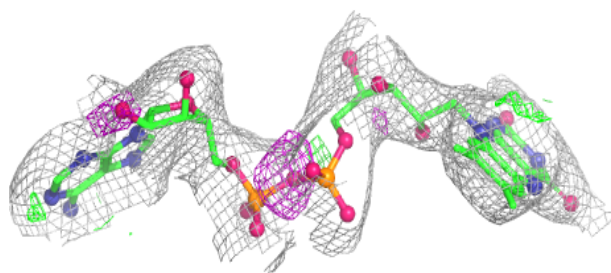
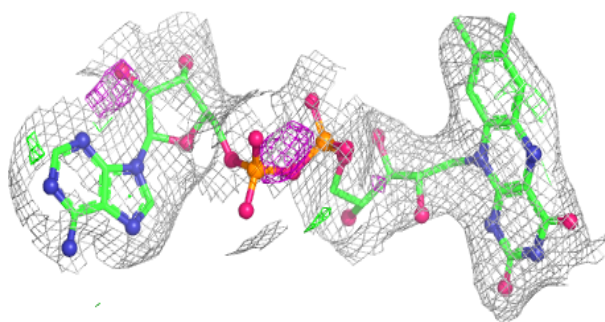
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	CIT	A	703	13/13	0.82	0.09	121,124,150,150	0
2	FAD	A	701	53/53	0.93	0.09	80,88,100,101	0
3	SAH	B	702	26/26	0.96	0.08	60,64,69,70	0
3	SAH	A	702	26/26	0.97	0.07	47,53,59,61	0

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.

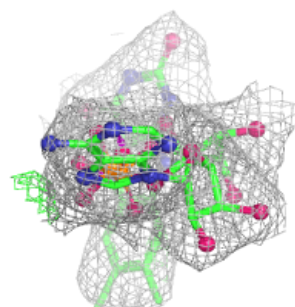
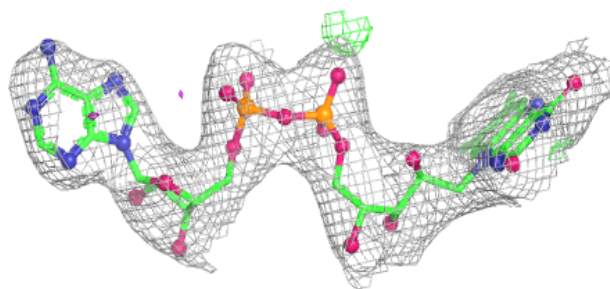
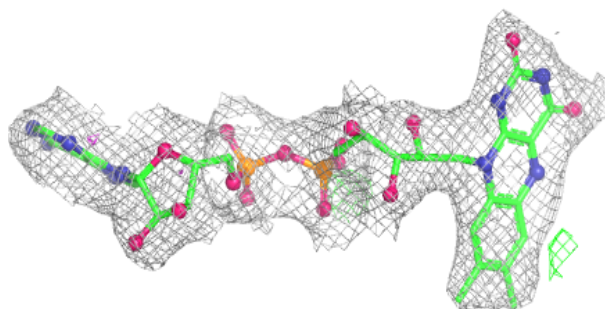
Electron density around FAD B 701:

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

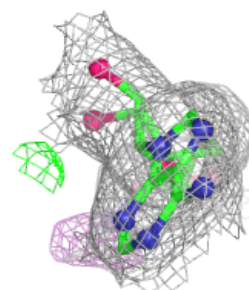
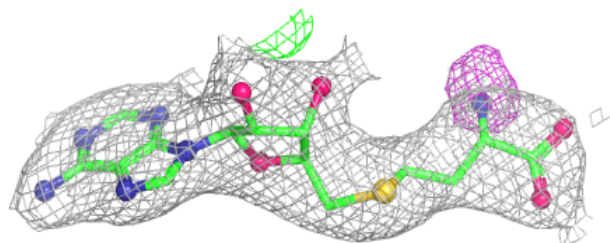
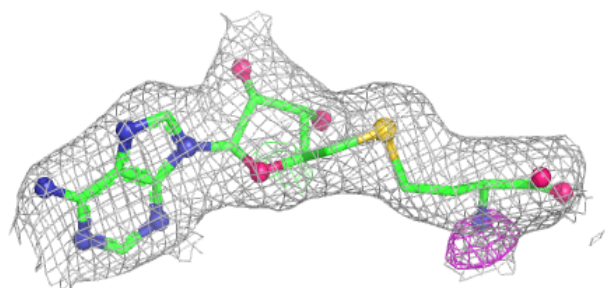


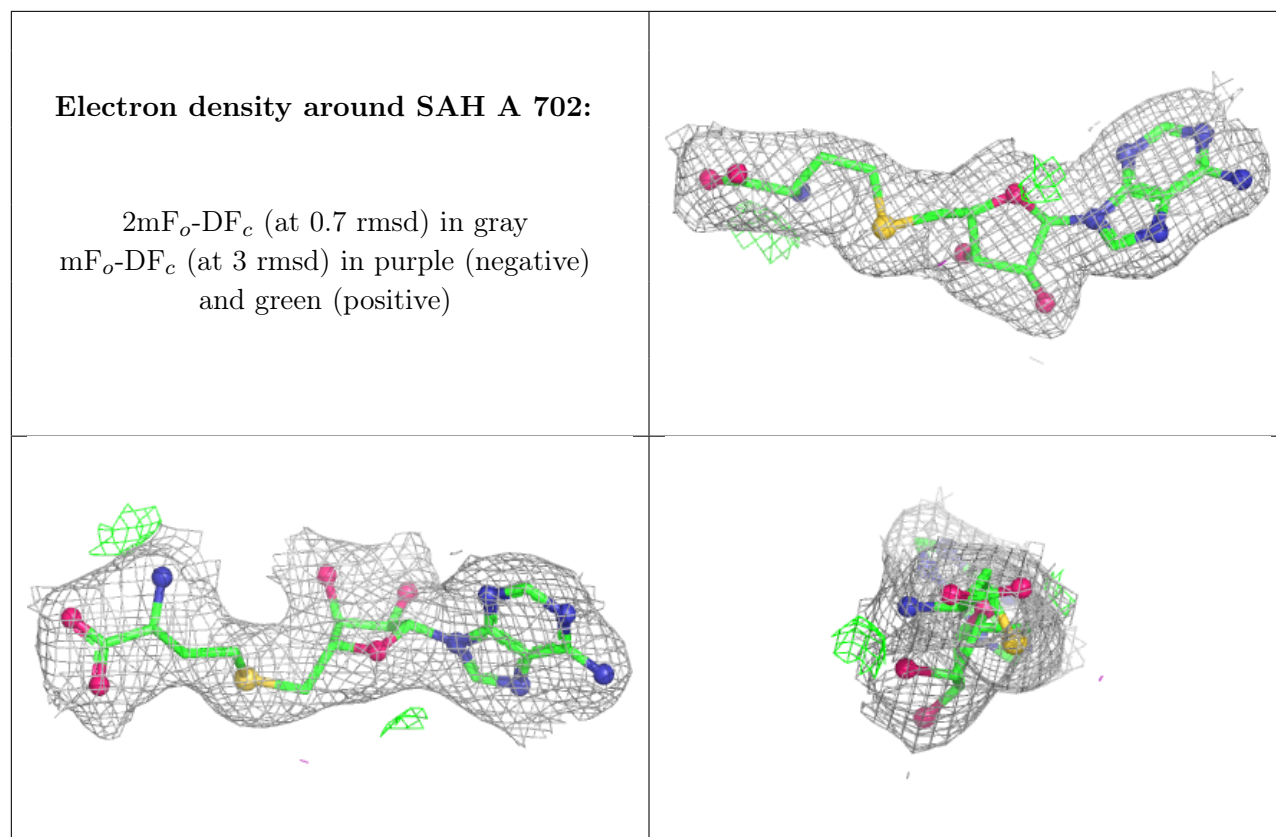
Electron density around FAD A 701:

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

**Electron density around SAH B 702:**

$2mF_o-DF_c$ (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.