

Full wwPDB X-ray Structure Validation Report (i)

May 26, 2020 – 01:31 pm BST

PDB ID Title	:	6FCX Structure of human 5.10-methylenetetrahydrofolate reductase (MTHER)
11010	•	Structure of numan 9,10-methyleneteeranydroiolate reductase (WTIII It)
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.; Baumgart-
		ner, 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

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

1 Overall quality at a glance (i)

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	$egin{array}{c} { m Whole \ archive} \ (\#{ m Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
R_{free}	130704	$4661 \ (2.50-2.50)$
Clashscore	141614	$5346\ (2.50-2.50)$
Ramachandran outliers	138981	$5231 \ (2.50-2.50)$
Sidechain outliers	138945	$5233 \ (2.50-2.50)$
RSRZ outliers	127900	4559 (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 on the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	А	615	% 7 9%	17%	
1	В	615	73%	21%	••



6 FCX

2 Entry composition (i)

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.

Mol	Chain	Residues		At	oms			ZeroOcc	AltConf	Trace
1	Λ	506	Total	С	Ν	Ο	S	0	0	0
L	л	090	4723	3048	786	870	19	0	0	
1	В	500	Total	С	Ν	Ο	S	0	0	0
	D	590	4191	2660	727	791	13	0	0	0

• Molecule 1 is a protein called Methylenetetrahydrofolate reductase.

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

There are 20 discrepancies between the modelled and reference sequences:

• Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).





Mol	Chain	Residues		At	oms			ZeroOcc	AltConf
9	Δ	1	Total	С	Ν	Ο	Р	0	0
	Z A	T	53	27	9	15	2	0	0
0	р	1	Total	С	Ν	Ο	Р	0	0
			53	27	9	15	2	0	

• Molecule 3 is S-ADENOSYL-L-HOMOCYSTEINE (three-letter code: SAH) (formula: $C_{14}H_{20}N_6O_5S$).



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



Mol	Chain	Residues		Ato	\mathbf{ms}			ZeroOcc	AltConf
2	р	1	Total	С	Ν	0	S	0	0
່ <u>ບ</u>	D	L	26	14	6	5	1	0	0

• Molecule 4 is CITRIC ACID (three-letter code: CIT) (formula: $C_6H_8O_7$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	Λ	1	Total	С	Η	Ο	0	0
4	Л	I	18	6	5	7	0	0
4	р	1	Total	С	Η	Ο	0	0
4	D	T	18	6	5	7	0	0

• Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	А	48	Total O 48 48	0	0
5	В	24	TotalO2424	0	0



3 Residue-property plots (i)

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



• Molecule 1: Methylenetetrahydrofolate reductase







4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants	97.34Å 127.94 Å 147.01 Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
$\mathbf{Bosolution} \left(\overset{\circ}{\mathbf{A}} \right)$	63.97 - 2.50	Depositor
Resolution (A)	63.97 - 2.50	EDS
% Data completeness	99.9(63.97 - 2.50)	Depositor
(in resolution range)	$100.0\ (63.97‐2.50)$	EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$< I/\sigma(I) > 1$	$1.44 (at 2.51 \text{\AA})$	Xtriage
Refinement program	PHENIX (1.13_2998: ???)	Depositor
D D.	0.210 , 0.247	Depositor
Π, Π_{free}	0.215 , 0.250	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 $k_{sol}(e/Å^3), B_{sol}(Å^2)$	0.28 , 64.3	EDS
L-test for $twinning^2$	$ < L >=0.49, < L^2>=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.

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.



¹Intensities estimated from amplitudes.

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: SAH, FAD, CIT

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).

Mal	Chain	Bond lengths		Bond angles	
		RMSZ	# Z > 5	RMSZ	# Z > 5
1	А	0.65	0/4853	0.83	13/6606~(0.2%)
1	В	0.55	0/4295	0.70	5/5885~(0.1%)
All	All	0.61	0/9148	0.77	18/12491~(0.1%)

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	А	0	5
1	В	0	2
All	All	0	7

There are no bond length outliers.

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
1	А	90	TYR	CB-CG-CD2	-9.15	115.51	121.00
1	В	515	PHE	CB-CG-CD2	-8.94	114.55	120.80
1	А	65	PHE	CB-CG-CD2	-8.56	114.81	120.80
1	А	60	PHE	CB-CG-CD2	-7.69	115.42	120.80
1	А	515	PHE	CB-CG-CD2	-7.60	115.48	120.80
1	А	361	ASP	CB-CG-OD1	7.41	124.97	118.30
1	В	515	PHE	CB-CG-CD1	6.62	125.43	120.80
1	А	515	PHE	CB-CG-CD1	6.52	125.36	120.80
1	А	65	PHE	CB-CG-CD1	6.36	125.25	120.80
1	А	90	TYR	CB-CG-CD1	6.33	124.80	121.00
1	А	400	LEU	CB-CG-CD2	5.64	120.60	111.00
1	В	190	PHE	CB-CG-CD1	5.57	124.70	120.80



Mol	Chain	Res	Type	Atoms	Z	$\mathbf{Observed}(^{o})$	$Ideal(^{o})$
1	А	60	PHE	CB-CG-CD1	5.50	124.65	120.80
1	В	538	TYR	CB-CG-CD2	-5.50	117.70	121.00
1	В	363	ARG	NE-CZ-NH2	-5.44	117.58	120.30
1	А	257	PHE	CB-CG-CD2	-5.35	117.06	120.80
1	А	90	TYR	CA-CB-CG	5.19	123.27	113.40
1	А	361	ASP	CB-CG-OD2	-5.05	113.75	118.30

There are no chirality outliers.

All (7) planarit	y outliers	are listed	below:
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Mol	Chain	\mathbf{Res}	Type	Group
1	А	404	TYR	Peptide
1	А	515	PHE	Sidechain
1	А	60	PHE	Sidechain
1	А	65	PHE	Sidechain
1	А	90	TYR	Sidechain
1	В	515	PHE	Sidechain
1	В	538	TYR	Sidechain

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	А	4723	0	4516	120	0
1	В	4191	0	3577	140	0
2	А	53	0	31	3	0
2	В	53	0	30	2	0
3	А	26	0	19	0	0
3	В	26	0	17	0	0
4	А	13	5	5	1	0
4	В	13	5	5	1	0
5	А	48	0	0	0	0
5	В	24	0	0	2	0
All	All	9170	10	8200	251	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.



• • •		Interatomic	Clash	
Atom-1	Atom-2	distance (Å)	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:HG3	1:A:571:GLN:O	1.67	0.92	
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: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:HG3	1:B:571:GLN:O	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: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: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	

All (251) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.



		Interatomic	Clash
Atom-1	Atom-2	distance $(Å)$	overlan (Å)
1:A:318:LEU:HB3	1:A:320:PHE:CE1	2.29	0.67
1:B:154:MET:CE	1:B:156:LEU:HD11	2.26	0.66
1:A:406:PHE:HE1	1:A:407:TYB: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:A:58:LYS:HG3	1:A:310:LEU:HD12	1.79	0.64
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:58:LYS:HE3	1:A:315:VAL:O	2.00	0.62
1:A:41:ARG:HG3	1:A:41:ARG:O	1.99	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: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:155:ALA:C	1:B:156:LEU:HD12	2.22	0.60
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:CD1	1:B:190:PHE:O	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:B:62:LEU:HD13	1:B:83:MET:HE1	1.85	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



		Interatomic	Clash
Atom-1	Atom-2	distance (Å)	overlap (Å)
1:B:53:LEU:HA	1:B:57:ASP:O	2.04	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:406:PHE:CE1	1:A:407:TYR:CE2	2.92	0.57
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: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:A:60:PHE:C	1:A:60:PHE:CD1	2.78	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:405:LEU:HB3	1:B:407:TYR:CE1	2.42	0.54
1:B:48:LYS:O	1:B:52:ARG:CD	2.56	0.54
1:A:276:VAL:O	1:A:281:LYS:HD2	2.08	0.53
1:B:306:CYS:HA	1:B:309:LEU:CB	2.38	0.53
1:B:90:TYR:CD2	1:B:125:ILE:CD1	2.92	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:B:190:PHE:CD1	1:B:190:PHE:C	2.82	0.53
1:A:279:GLU:CD	1:A:279:GLU:H	2.12	0.53
1:A:41:ARG:HH12	1:A:52:ARG:NH2	2.05	0.53
1:A:65:PHE:CD1	1:A:65:PHE:C	2.83	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: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:B:438:TYR:HA	1:B:450:VAL:CG2	2.40	0.51
1:B:96:HIS:CD2	1:B:98:ALA:HB3	2.45	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: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



		Interatomic	Clash
Atom-1	Atom-2	distance (\AA)	overlap (Å)
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
1:B:91:ILE:HG13	1:B:92:ASP:H	1.77	0.50
1:B:67:PRO:HG3	1:B:73:ALA:HA	1.93	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: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: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:515:PHE:CD1	1:B:515:PHE:C	2.87	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:467:LEU:CD1	1:A:471:LEU:CD1	2.92	0.47
1:A:229:LEU:C	1:A:229:LEU:HD12	2.35	0.47
1:A:406:PHE:CE1	1:A:407:TYR:CD2	3.02	0.47
1:B:369:SER:C	1:B:371:PRO:HD3	2.34	0.47
1:A:60:PHE:C	1:A:60:PHE:HD1	2.17	0.46
1:B:370:ARG:N	1:B:371:PRO:HD3	2.30	0.46
1:A:276:VAL:HG22	1:A:281:LYS:HG3	1.96	0.46
1:B:435:PHE:CD2	1:B:480:LEU:HB3	2.50	0.46
1:B:74:VAL:HG12	1:B:406:PHE:O	2.15	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
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



		Interatomic	Clash
Atom-1	Atom-2	distance $(Å)$	overlap (Å)
1:B:538:TYR:C	1:B:538:TYR:CD1	2.90	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: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: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:B:538:TYR:HD1	1:B:538:TYR:C	2.20	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: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:145:ALA:C	1:B:150:LEU:HB2	2.38	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:A:406:PHE:CD1	1:A:406:PHE:C	2.90	0.44
1:A:65:PHE:C	1:A:65:PHE:HD1	2.20	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:305:LEU:HD23	1:A:309:LEU:HG	2.00	0.43
1:A:195:ALA:HB1	2:A:701:FAD:HM83	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: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:183:ARG:HA	1:B:186:PHE:CB	2.48	0.43
1:B:119:TYR:O	1:B:372:LYS:HE3	2.18	0.43
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:154:MET:HE2	1:B:156:LEU:HD11	1.99	0.43
1:B:194:VAL:O	1:B:225:ILE:HA	2.19	0.43



		Interatomic	Clash		
Atom-1	Atom-2	distance $(Å)$	overlap (Å)		
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:A:257:PHE:C	1:A:257:PHE:CD1	2.92	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: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:B:399:GLU:OE1	1:B:401:LYS:CE	2.66	0.42		
1:A:571:GLN:NE2	1:B:568:GLU:OE1	2.48	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:A:257:PHE:HD1	1:A:257:PHE:C	2.23	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:B:63:GLU:OE1	1:B:321:TYR:CE1	2.73	0.41		
1:B:386:ASN:C	1:B:386:ASN:OD1	2.58	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:HD23	1:A:265:LEU:HA	1.71	0.41		
1:B:91:ILE:HA	1:B:91:ILE:HD12	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:A:317:GLY:C	1:A:318:LEU:HD12	2.41	0.41		
1:A:65:PHE:O	1:A:65:PHE:CD1	2.74	0.41		
1:B:385:PRO:HG2	1:B:389:TRP:CZ3	2.56	0.41		
1:B:60:PHE:HA	1:B:60:PHE:HD1	1.61	0.41		
1:A:282:ASP:O	1:A:286:PRO:CD	2.69	0.41		
1:B:116:ALA:HB3	1:B:124:THR:HG21	2.02	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		
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	$2.\overline{50}$	0.41		
1:A:193:CYS:HB3	1:A:224:PHE:CE1	2.55	0.40		



Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:41:ARG:CG	1:A:41:ARG:O	2.69	0.40
1:B:325:ARG:HB3	4:B:703:CIT:O6	2.21	0.40
1:B:53:LEU:CD2	1:B:58:LYS:HA	2.51	0.40
1:A:60:PHE:CE2	1:A:333:LEU:HD13	2.56	0.40
1:B:526:LEU:HA	1:B:526:LEU:HD23	1.87	0.40
1:A:586:GLU:O	1:A:587:ALA:C	2.60	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

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

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

Mol	Chain	Analysed Favoured Allowed		Outliers	Perce	entiles	;	
1	А	590/615~(96%)	$570 \ (97\%)$	18 (3%)	2 (0%)	41	61	
1	В	580/615~(94%)	548 (94%)	32~(6%)	0	100	100	
All	All	1170/1230~(95%)	1118 (96%)	50(4%)	2(0%)	47	68	

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	А	205	GLY
1	А	221	GLY

5.3.2 Protein sidechains (i)

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

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



Mol	Chain	Analysed Rotameric Outliers		Percentiles		
1	А	491/537~(91%)	481 (98%)	10 (2%)	55 7	79
1	В	358/537~(67%)	348~(97%)	10 (3%)	43 7	70
All	All	849/1074~(79%)	829 (98%)	20 (2%)	49 7	74

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	А	60	PHE
1	А	65	PHE
1	А	90	TYR
1	А	257	PHE
1	А	320	PHE
1	А	406	PHE
1	А	435	PHE
1	А	515	PHE
1	А	538	TYR
1	А	584	LYS
1	В	60	PHE
1	В	157	ARG
1	В	190	PHE
1	В	407	TYR
1	В	414	LYS
1	В	473	ARG
1	В	515	PHE
1	В	538	TYR
1	В	619	VAL
1	В	620	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	В	96	HIS
1	В	477	GLN
1	В	555	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.



5.4 Non-standard residues in protein, DNA, RNA chains (i)

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

5.5 Carbohydrates (i)

There are no carbohydrates in this entry.

5.6 Ligand geometry (i)

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).

Mal	Tune	Chain	Dec	Tink	Bo	ond leng	\mathbf{ths}	В	ond ang	les
	Type	Chain	nes		Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	CIT	А	703	-	3,12,12	1.24	0	$3,\!17,\!17$	2.23	1 (33%)
2	FAD	А	701	-	51, 58, 58	1.30	5 (9%)	60,89,89	2.30	7 (11%)
2	FAD	В	701	1	51, 58, 58	1.22	5 (9%)	60,89,89	2.22	7 (11%)
4	CIT	В	703	-	$3,\!12,\!12$	0.99	0	$3,\!17,\!17$	1.74	1 (33%)
3	SAH	А	702	-	21,28,28	1.37	3 (14%)	$20,\!40,\!40$	1.97	6 (30%)
3	SAH	В	702	1	21,28,28	1.34	2(9%)	$20,\!40,\!40$	1.84	<mark>5 (25%)</mark>

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	\mathbf{Res}	Link	Chirals	Torsions	Rings
4	CIT	А	703	-	-	4/6/16/16	-
2	FAD	А	701	-	-	4/30/50/50	0/6/6/6
2	FAD	В	701	1	-	16/30/50/50	0/6/6/6
4	CIT	В	703	-	-	2/6/16/16	-
3	SAH	А	702	-	-	0/7/31/31	0/3/3/3
3	SAH	В	702	1	-	1/7/31/31	0/3/3/3



6FCX

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	А	701	FAD	C4X-C10	5.82	1.44	1.38
2	В	701	FAD	C4X-C10	5.61	1.44	1.38
3	А	702	SAH	C2-N3	4.08	1.38	1.32
3	В	702	SAH	C2-N3	3.93	1.38	1.32
2	А	701	FAD	C4X-N5	-3.34	1.28	1.33
2	В	701	FAD	C4-N3	3.09	1.38	1.33
3	В	702	SAH	C2-N1	3.00	1.39	1.33
2	А	701	FAD	C4-N3	2.91	1.38	1.33
2	В	701	FAD	C4-C4X	2.66	1.45	1.41
2	В	701	FAD	C5X-N5	2.66	1.39	1.35
3	А	702	SAH	C2-N1	2.61	1.38	1.33
2	А	701	FAD	C4-C4X	2.57	1.45	1.41
2	В	701	FAD	C9A-N10	2.47	1.41	1.38
3	А	702	SAH	O4'-C1'	2.26	1.44	1.41
2	А	701	FAD	C9A-N10	2.17	1.41	1.38

All (15) bond length outliers are listed below:

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
2	А	701	FAD	C4-N3-C2	13.05	126.16	115.14
2	В	701	FAD	C4-N3-C2	12.98	126.10	115.14
2	А	701	FAD	C4X-C4-N3	-7.31	113.43	123.43
2	В	701	FAD	C4X-C4-N3	-7.00	113.86	123.43
3	А	702	SAH	N3-C2-N1	-6.21	118.97	128.68
3	В	702	SAH	N3-C2-N1	-4.74	121.26	128.68
2	А	701	FAD	C10-C4X-N5	4.61	124.44	121.26
2	В	701	FAD	C10-C4X-N5	4.46	124.34	121.26
2	А	701	FAD	C1'-N10-C9A	3.91	121.37	118.29
2	В	701	FAD	C4-C4X-C10	-3.73	117.48	119.95
4	А	703	CIT	C3-C4-C5	-3.55	109.30	114.98
2	А	701	FAD	C4-C4X-C10	-3.43	117.68	119.95
2	В	701	FAD	C4X-C10-N10	-3.41	116.79	120.30
2	А	701	FAD	C4X-C10-N10	-3.33	116.88	120.30
3	А	702	SAH	C5'-SD-CG	-3.00	93.27	102.27
3	В	702	SAH	O4'-C1'-C2'	-2.96	102.60	106.93
2	В	701	FAD	C1'-N10-C9A	2.80	120.50	118.29
3	В	702	SAH	O4'-C4'-C5'	2.41	115.05	108.83
3	А	702	SAH	N6-C6-N1	2.37	123.49	118.57
3	В	702	SAH	C4'-C5'-SD	-2.33	105.43	113.78
3	А	702	SAH	O3'-C3'-C4'	-2.32	104.33	111.05
2	А	701	FAD	C5A-C6A-N6A	2.31	123.86	120.35
3	А	702	SAH	O4'-C1'-C2'	-2.25	103.64	106.93



Mol	Chain	Res	Type	Atoms	Z	$\mathbf{Observed}(^{o})$	$Ideal(^{o})$
2	В	701	FAD	C5A-C6A-N6A	2.24	123.75	120.35
3	В	702	SAH	C5'-SD-CG	-2.23	95.57	102.27
4	В	703	CIT	C3-C4-C5	-2.14	111.56	114.98
3	А	702	SAH	C5-C6-N6	-2.13	117.11	120.35

There are no chirality outliers.

All (27) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	А	703	CIT	C1-C2-C3-O7
4	А	703	CIT	C1-C2-C3-C4
4	А	703	CIT	C1-C2-C3-C6
2	А	701	FAD	C1'-C2'-C3'-O3'
2	А	701	FAD	C1'-C2'-C3'-C4'
2	В	701	FAD	C5B-O5B-PA-O1A
2	В	701	FAD	C1'-C2'-C3'-C4'
2	В	701	FAD	C3'-C4'-C5'-O5'
2	В	701	FAD	O4'-C4'-C5'-O5'
2	В	701	FAD	C5'-O5'-P-O2P
2	В	701	FAD	C5'-O5'-P-O3P
3	В	702	SAH	CA-CB-CG-SD
2	А	701	FAD	O2'-C2'-C3'-C4'
2	В	701	FAD	O3'-C3'-C4'-O4'
2	В	701	FAD	C2'-C3'-C4'-O4'
2	В	701	FAD	O2'-C2'-C3'-C4'
2	В	701	FAD	C2'-C3'-C4'-C5'
2	А	701	FAD	O2'-C2'-C3'-O3'
4	В	703	CIT	C1-C2-C3-O7
2	В	701	FAD	C5B-O5B-PA-O3P
2	В	701	FAD	C5B-O5B-PA-O2A
2	В	701	FAD	C5'-O5'-P-O1P
2	В	701	FAD	C1'-C2'-C3'-O3'
4	A	703	CIT	C6-C3-C4-C5
4	В	703	CIT	C2-C3-C4-C5
2	В	701	FAD	O3'-C3'-C4'-C5'
2	В	701	FAD	O2'-C2'-C3'-O3'

There are no ring outliers.

4 monomers are involved in 7 short contacts:



Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	А	703	CIT	1	0
2	А	701	FAD	3	0
2	В	701	FAD	2	0
4	В	703	CIT	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 then 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 similar 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, 95^{th} 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		$\mathbf{OWAB}(\mathbf{\AA}^2)$	$Q{<}0.9$
1	А	596/615~(96%)	-0.08	8 (1%) 77 7	'9	48, 74, 131, 160	0
1	В	590/615~(95%)	0.34	48 (8%) 12	12	52, 97, 194, 241	0
All	All	1186/1230~(96%)	0.13	56 (4%) 31 3	33	48, 83, 181, 241	0

All (56) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	В	214	LEU	9.5
1	В	250	CYS	8.8
1	В	240	VAL	6.5
1	В	248	ILE	6.3
1	В	236	PHE	6.2
1	В	247	GLY	6.0
1	В	217	LYS	5.4
1	В	276	VAL	5.4
1	В	224	PHE	4.7
1	В	249	THR	4.5
1	В	194	VAL	4.4
1	А	247	GLY	4.3
1	В	243	CYS	4.2
1	В	215	LYS	4.2
1	В	302	ALA	4.2
1	В	252	ILE	3.8
1	В	310	LEU	3.8
1	В	208	GLU	3.8
1	В	317	GLY	3.6
1	В	239	PHE	3.5
1	A	246	MET	3.4
1	A	248	ILE	3.4
1	A	310	LEU	3.4
1	В	128	MET	3.4



Mol	Chain	Res	Type	RSRZ
1	В	95	TRP	3.2
1	В	277	PRO	3.1
1	В	191	ASP	3.0
1	А	403	TYR	3.0
1	В	216	GLU	2.9
1	В	339	TRP	2.8
1	В	212	LYS	2.8
1	В	406	PHE	2.7
1	В	241	LYS	2.7
1	В	318	LEU	2.7
1	В	272	SER	2.7
1	В	206	SER	2.7
1	В	223	ASP	2.5
1	В	130	CYS	2.5
1	В	218	VAL	2.4
1	В	646	GLU	2.4
1	В	238	ARG	2.4
1	В	182	ILE	2.4
1	В	246	MET	2.4
1	А	262	TYR	2.3
1	В	155	ALA	2.3
1	В	50	ARG	2.3
1	В	192	ILE	2.3
1	В	404	TYR	2.2
1	В	56	GLY	2.2
1	В	328	ALA	2.1
1	А	40	GLU	2.1
1	В	148	LEU	2.1
1	В	141	HIS	2.1
1	В	210	ASP	2.1
1	В	144	LYS	2.0
1	А	236	PHE	2.0

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

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

6.3 Carbohydrates (i)

There are no carbohydrates in this entry.



6.4 Ligands (i)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} 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
2	FAD	В	701	53/53	0.82	0.18	102,125,139,142	0
4	CIT	В	703	13/13	0.83	0.18	144,145,175,175	0
4	CIT	А	703	13/13	0.89	0.11	$121,\!124,\!150,\!150$	0
2	FAD	А	701	53/53	0.95	0.12	80,88,100,101	0
3	SAH	В	702	26/26	0.95	0.15	60,64,69,70	0
3	SAH	А	702	26/26	0.98	0.15	$47,\!53,\!59,\!61$	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.











6.5 Other polymers (i)

There are no such residues in this entry.

