



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 10:58 PM GMT

PDB ID : 1W1L  
Title : STRUCTURE OF THE OCTAMERIC FLAVOENZYME VANILLYL-ALCOHOL OXIDASE: PHE454TYR MUTANT  
Authors : Van Den Heuvel, R.H.  
Deposited on : 2004-06-22  
Resolution : 2.70 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](mailto:validation@mail.wwpdb.org)  
A user guide is available at  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

---

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

MolProbity : 4.02b-467  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

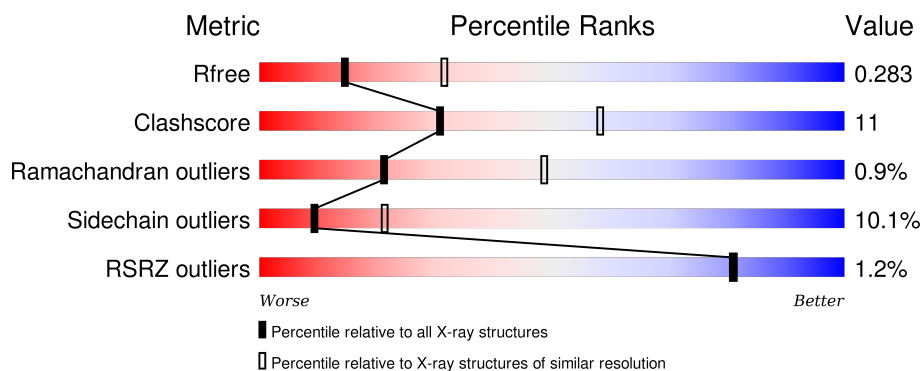
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.70 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))
$R_{free}$	91344	2103 (2.70-2.70)
Clashscore	102246	2422 (2.70-2.70)
Ramachandran outliers	100387	2382 (2.70-2.70)
Sidechain outliers	100360	2382 (2.70-2.70)
RSRZ outliers	91569	2107 (2.70-2.70)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	560	<div> <div>64%</div> <div>29%</div> <div>5%</div> <div>.</div> </div>
1	B	560	<div> <div>72%</div> <div>23%</div> <div>.</div> <div>.</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	EUG	A	1562	-	-	-	X
3	EUG	B	1562	-	-	-	X

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 8832 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 VANILLYL-ALCOHOL OXIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	550	Total	C	N	O	S	0	0	0
			4352	2793	744	791	24			
1	B	550	Total	C	N	O	S	0	0	0
			4352	2793	744	791	24			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	454	TYR	PHE	ENGINEERED MUTATION	UNP P56216
B	454	TYR	PHE	ENGINEERED MUTATION	UNP P56216

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



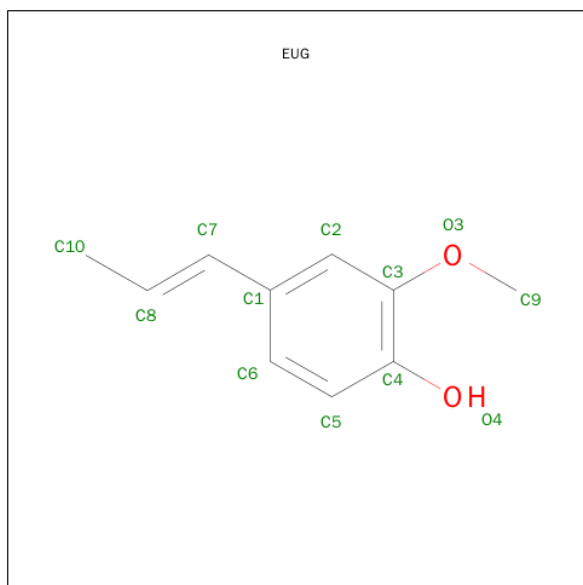
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

*Continued on next page...*

*Continued from previous page...*

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

- Molecule 3 is 2-METHOXY-4-VINYL-PHENOL (three-letter code: EUG) (formula:  $C_{10}H_{12}O_2$ ).

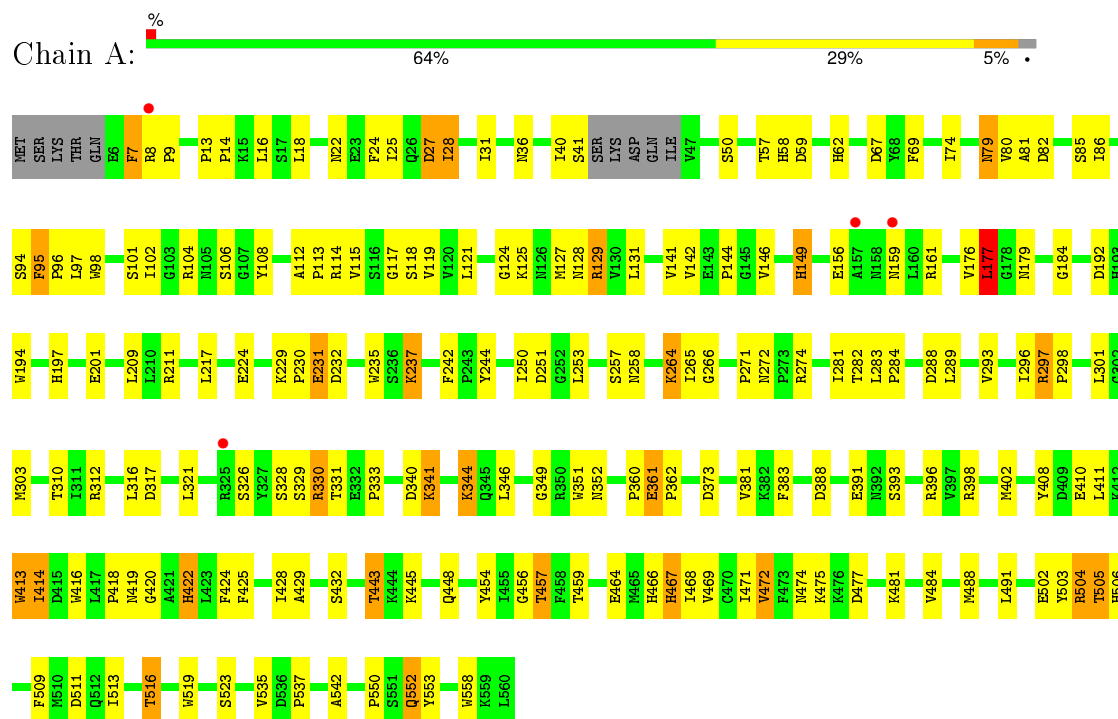


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			11	9	2		
3	B	1	Total	C	O	0	0
			11	9	2		

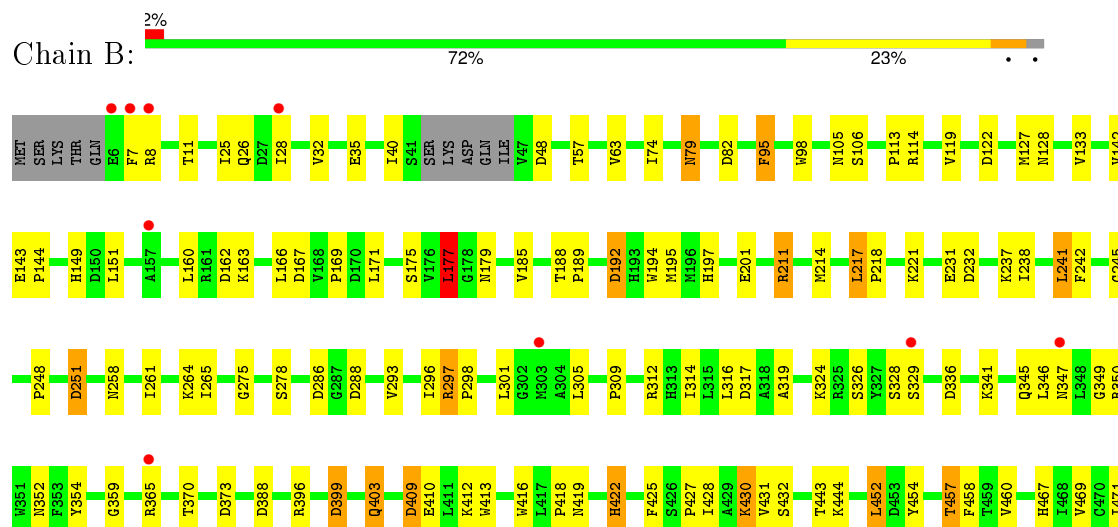
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: VANILLYL-ALCOHOL OXIDASE



#### • Molecule 1: VANILLYL-ALCOHOL OXIDASE





## 4 Data and refinement statistics

Property	Value	Source
Space group	I 4	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	129.90Å 129.90Å 133.61Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.00 – 2.70 42.13 – 2.70	Depositor EDS
% Data completeness (in resolution range)	97.6 (40.00-2.70) 97.6 (42.13-2.70)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.74 (at 2.69Å)	Xtriage
Refinement program	REFMAC 5.1.27	Depositor
R, $R_{free}$	0.203 , 0.293 0.203 , 0.283	Depositor DCC
$R_{free}$ test set	1508 reflections (5.35%)	DCC
Wilson B-factor (Å <sup>2</sup> )	46.7	Xtriage
Anisotropy	0.555	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 32.3	EDS
Estimated twinning fraction	0.007 for l,-k,h 0.020 for -l,-k,-h 0.021 for -h,-l,-k 0.008 for -h,l,k 0.037 for -h,k,-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	1 of 29696 reflections (0.003%)	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	8832	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.85% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup> Intensities estimated from amplitudes.

<sup>2</sup> Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: EUG, FAD

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	0.74	0/4471	0.91	11/6077 (0.2%)
1	B	0.54	0/4471	0.83	16/6077 (0.3%)
All	All	0.65	0/8942	0.87	27/12154 (0.2%)

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	2
1	B	0	1
All	All	0	3

There are no bond length outliers.

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	A	317	ASP	CB-CG-OD2	6.51	124.16	118.30
1	B	388	ASP	CB-CG-OD2	6.42	124.08	118.30
1	A	192	ASP	CB-CG-OD2	6.37	124.03	118.30
1	B	192	ASP	CB-CG-OD2	6.26	123.94	118.30
1	B	232	ASP	CB-CG-OD2	6.14	123.83	118.30
1	B	177	LEU	CA-CB-CG	6.04	129.20	115.30
1	B	251	ASP	CB-CG-OD2	5.96	123.66	118.30
1	B	288	ASP	CB-CG-OD2	5.86	123.58	118.30
1	A	388	ASP	CB-CG-OD2	5.86	123.57	118.30
1	A	251	ASP	CB-CG-OD2	5.73	123.46	118.30
1	A	177	LEU	CA-CB-CG	5.67	128.33	115.30
1	B	167	ASP	CB-CG-OD2	5.66	123.39	118.30

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	373	ASP	CB-CG-OD2	5.65	123.39	118.30
1	A	27	ASP	CB-CG-OD2	5.56	123.31	118.30
1	B	82	ASP	CB-CG-OD2	5.52	123.27	118.30
1	A	340	ASP	CB-CG-OD2	5.52	123.26	118.30
1	B	286	ASP	CB-CG-OD2	5.46	123.21	118.30
1	B	317	ASP	CB-CG-OD2	5.45	123.21	118.30
1	A	288	ASP	CB-CG-OD2	5.44	123.19	118.30
1	B	511	ASP	CB-CG-OD2	5.33	123.10	118.30
1	B	477	ASP	CB-CG-OD2	5.29	123.06	118.30
1	A	67	ASP	CB-CG-OD2	5.25	123.02	118.30
1	B	399	ASP	CB-CG-OD2	5.22	123.00	118.30
1	B	48	ASP	CB-CG-OD2	5.21	122.99	118.30
1	B	122	ASP	CB-CG-OD2	5.10	122.89	118.30
1	A	232	ASP	CB-CG-OD2	5.08	122.88	118.30
1	B	373	ASP	CB-CG-OD2	5.03	122.83	118.30

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	127	MET	Peptide
1	A	141	VAL	Peptide
1	B	127	MET	Peptide

## 5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4352	0	4288	118	0
1	B	4352	0	4288	81	0
2	A	53	0	29	2	0
2	B	53	0	29	2	0
3	A	11	0	7	0	0
3	B	11	0	7	1	0
All	All	8832	0	8648	191	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 11.

All (191) 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)
2:A:600:FAD:O5B	2:A:600:FAD:C5B	1.66	1.42
1:A:443:THR:HG21	1:A:469:VAL:HG21	1.50	0.93
1:B:513:ILE:O	1:B:516:THR:HB	1.76	0.83
1:A:443:THR:HA	1:A:491:LEU:HD21	1.63	0.81
1:A:552:GLN:H	1:A:552:GLN:NE2	1.79	0.81
1:A:177:LEU:HB2	1:A:265:ILE:HG21	1.64	0.80
1:A:310:THR:HG22	1:A:459:THR:HG22	1.63	0.79
1:A:211:ARG:HG3	1:B:519:TRP:CZ3	2.18	0.78
1:A:177:LEU:HB2	1:A:265:ILE:CG2	2.13	0.77
1:A:257:SER:HB3	1:B:248:PRO:HB3	1.66	0.77
1:A:349:GLY:H	1:A:352:ASN:HD21	1.35	0.74
1:A:454:TYR:OH	1:A:467:HIS:HE1	1.69	0.74
1:A:250:ILE:O	1:A:253:LEU:HB2	1.86	0.74
1:B:297:ARG:HB3	1:B:298:PRO:CD	2.20	0.70
1:B:297:ARG:HG3	1:B:431:VAL:O	1.92	0.70
1:A:454:TYR:OH	1:A:467:HIS:CE1	2.46	0.69
1:B:297:ARG:HB3	1:B:298:PRO:HD3	1.74	0.68
1:A:550:PRO:HB2	1:A:552:GLN:NE2	2.09	0.68
1:B:312:ARG:NH2	1:B:410:GLU:OE1	2.27	0.68
1:A:14:PRO:HG3	1:A:558:TRP:HZ2	1.59	0.67
1:A:513:ILE:O	1:A:516:THR:HB	1.94	0.66
1:A:328:SER:OG	1:A:330:ARG:HB2	1.96	0.66
1:A:24:PHE:HD2	1:A:25:ILE:HD12	1.60	0.66
1:A:312:ARG:HG3	1:A:457:THR:HG23	1.78	0.66
1:B:504:ARG:O	1:B:505:THR:HG23	1.95	0.66
1:A:80:VAL:HG11	1:A:209:LEU:HD11	1.77	0.66
1:A:416:TRP:HB3	1:A:472:VAL:HG21	1.78	0.65
1:B:550:PRO:HB2	1:B:552:GLN:HE21	1.61	0.65
1:B:349:GLY:H	1:B:352:ASN:HD21	1.44	0.64
1:B:275:GLY:HA3	1:B:359:GLY:O	1.98	0.64
1:A:237:LYS:NZ	1:B:498:ASN:O	2.31	0.63
1:B:95:PHE:CE1	1:B:119:VAL:HG23	2.33	0.63
1:B:79:ASN:HD22	1:B:79:ASN:C	2.03	0.62
1:A:14:PRO:HG3	1:A:558:TRP:CZ2	2.35	0.62
1:A:201:GLU:OE2	1:A:211:ARG:HD3	1.99	0.61
1:A:24:PHE:CD2	1:A:25:ILE:HD12	2.36	0.61
1:A:36:ASN:OD1	1:A:125:LYS:NZ	2.35	0.60
1:A:108:TYR:CZ	1:A:504:ARG:HG2	2.37	0.60

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:106:SER:HB2	1:A:422:HIS:HE1	1.67	0.60
1:A:7:PHE:HA	1:A:22:ASN:OD1	2.02	0.59
1:A:398:ARG:NH1	1:A:410:GLU:OE2	2.35	0.59
1:B:552:GLN:NE2	1:B:552:GLN:H	2.01	0.59
1:A:57:THR:HG22	1:A:74:ILE:HD11	1.85	0.58
1:B:550:PRO:HB2	1:B:552:GLN:NE2	2.19	0.58
1:A:297:ARG:HB3	1:A:298:PRO:CD	2.34	0.58
1:A:505:THR:OG1	1:A:513:ILE:HD12	2.05	0.57
1:A:95:PHE:CE1	1:A:119:VAL:HG23	2.40	0.57
1:B:177:LEU:HB2	1:B:265:ILE:CG2	2.34	0.56
1:B:197:HIS:HE1	1:B:251:ASP:OD2	1.89	0.56
1:A:519:TRP:CZ3	1:B:211:ARG:HG2	2.40	0.56
1:A:101:SER:O	1:A:124:GLY:HA3	2.06	0.56
1:A:398:ARG:O	1:A:402:MET:HG3	2.06	0.56
1:A:129:ARG:HH21	1:A:131:LEU:HD21	1.70	0.56
1:A:115:VAL:HG13	1:A:558:TRP:CH2	2.41	0.55
1:A:31:ILE:HD13	1:A:85:SER:HB3	1.88	0.55
1:B:28:ILE:O	1:B:32:VAL:HG22	2.06	0.55
1:A:244:TYR:OH	1:B:195:MET:HG2	2.06	0.55
1:A:464:GLU:HG2	1:B:242:PHE:HB2	1.88	0.55
1:B:177:LEU:HB2	1:B:265:ILE:HG21	1.89	0.55
1:A:82:ASP:O	1:A:86:ILE:HD12	2.06	0.55
1:B:428:ILE:HD11	1:B:503:TYR:HB3	1.88	0.54
1:A:411:LEU:O	1:A:414:ILE:HB	2.07	0.54
1:A:177:LEU:HB2	1:A:265:ILE:HG22	1.89	0.54
1:A:552:GLN:HE21	1:A:552:GLN:H	1.51	0.54
1:A:211:ARG:HG3	1:B:519:TRP:CH2	2.43	0.54
1:B:194:TRP:O	1:B:197:HIS:HD2	1.91	0.53
1:B:454:TYR:C	1:B:454:TYR:CD2	2.81	0.53
1:B:403:GLN:O	1:B:403:GLN:HG3	2.08	0.53
1:B:201:GLU:CB	1:B:264:LYS:HB2	2.38	0.53
1:A:96:PRO:HG3	1:A:553:TYR:OH	2.08	0.53
1:A:312:ARG:HG2	1:A:316:LEU:HD23	1.90	0.53
1:A:443:THR:HG21	1:A:469:VAL:CG2	2.33	0.52
1:B:261:ILE:HD11	1:B:541:ILE:HD12	1.91	0.52
1:A:297:ARG:HB3	1:A:298:PRO:HD3	1.91	0.52
1:A:506:HIS:HB3	1:A:509:PHE:HD2	1.74	0.52
1:A:502:GLU:HB2	1:A:513:ILE:HD13	1.91	0.52
1:B:114:ARG:NH2	1:B:511:ASP:OD2	2.35	0.52
1:A:69:PHE:HB2	1:A:112:ALA:HB1	1.92	0.52
1:B:171:LEU:HD21	2:B:1561:FAD:HM73	1.91	0.51

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:289:LEU:HD22	1:A:351:TRP:CZ2	2.45	0.51
1:A:194:TRP:O	1:A:197:HIS:HD2	1.93	0.51
1:A:349:GLY:H	1:A:352:ASN:ND2	2.08	0.51
1:B:177:LEU:HD12	1:B:177:LEU:C	2.30	0.51
1:A:454:TYR:CD2	1:A:454:TYR:C	2.84	0.51
1:A:101:SER:HB3	1:A:144:PRO:O	2.11	0.51
1:A:419:ASN:O	1:A:474:ASN:HA	2.11	0.51
1:A:477:ASP:O	1:A:481:LYS:HG3	2.11	0.51
1:A:114:ARG:NH2	1:A:511:ASP:OD2	2.42	0.50
1:A:361:GLU:HB3	1:A:362:PRO:HD3	1.94	0.50
1:A:24:PHE:O	1:A:27:ASP:HB2	2.12	0.50
2:B:1561:FAD:O3'	2:B:1561:FAD:N1	2.36	0.50
1:A:535:VAL:O	1:A:537:PRO:HD3	2.11	0.50
1:B:57:THR:HG22	1:B:74:ILE:HD11	1.93	0.50
1:B:409:ASP:OD1	1:B:409:ASP:N	2.44	0.50
1:A:237:LYS:HB3	1:B:430:LYS:HD2	1.94	0.50
1:B:133:VAL:HG11	1:B:160:LEU:HD13	1.92	0.50
1:B:258:ASN:OD1	1:B:532:LYS:HD3	2.12	0.50
1:A:211:ARG:HD2	1:A:235:TRP:CH2	2.48	0.49
1:B:201:GLU:HB3	1:B:264:LYS:HB2	1.95	0.49
1:A:149:HIS:HD1	1:A:408:TYR:HH	1.59	0.49
1:A:62:HIS:H	1:A:62:HIS:CD2	2.29	0.49
1:A:281:ILE:HG12	1:A:383:PHE:HD1	1.76	0.49
2:A:600:FAD:C4B	2:A:600:FAD:O5B	2.55	0.49
1:B:316:LEU:HD23	1:B:457:THR:HG23	1.94	0.49
1:B:98:TRP:CD2	1:B:113:PRO:HA	2.47	0.49
1:B:106:SER:HB3	1:B:422:HIS:HE1	1.77	0.49
1:A:9:PRO:HA	1:A:40:ILE:O	2.12	0.48
1:A:18:LEU:O	1:A:22:ASN:ND2	2.46	0.48
1:A:146:VAL:O	1:A:176:VAL:HG23	2.14	0.48
1:A:471:ILE:HG21	1:A:484:VAL:HG13	1.95	0.48
1:A:330:ARG:NH2	1:A:333:PRO:O	2.46	0.48
1:A:197:HIS:HA	1:A:266:GLY:O	2.14	0.48
1:A:293:VAL:HA	1:A:296:ILE:HD12	1.95	0.47
1:A:349:GLY:N	1:A:352:ASN:HD21	2.08	0.47
1:A:360:PRO:HB2	1:A:362:PRO:HD2	1.97	0.47
1:A:59:ASP:OD2	1:A:62:HIS:HA	2.15	0.47
1:A:445:LYS:O	1:A:448:GLN:HB3	2.14	0.47
1:B:238:ILE:HA	1:B:241:LEU:HD12	1.98	0.46
1:B:296:ILE:HG23	1:B:305:LEU:CD1	2.45	0.46
1:B:425:PHE:HZ	1:B:495:CYS:HG	1.64	0.46

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:414:ILE:O	1:A:420:GLY:HA3	2.15	0.46
1:B:346:LEU:O	1:B:347:ASN:HB2	2.15	0.46
1:A:97:LEU:HD13	1:A:121:LEU:HD22	1.97	0.46
1:A:242:PHE:CZ	1:A:244:TYR:HB2	2.51	0.45
1:A:58:HIS:C	1:A:112:ALA:HB2	2.36	0.45
1:A:79:ASN:HD21	1:A:81:ALA:HB3	1.81	0.45
1:A:422:HIS:CD2	1:A:424:PHE:CZ	3.04	0.45
1:A:258:ASN:HB2	1:A:542:ALA:HB3	1.99	0.45
1:B:506:HIS:HB3	1:B:509:PHE:HD2	1.81	0.45
1:B:319:ALA:CB	1:B:413:TRP:HB2	2.47	0.45
1:B:201:GLU:HB2	1:B:264:LYS:HB2	1.98	0.45
1:B:425:PHE:HB3	1:B:469:VAL:HB	1.98	0.45
1:B:458:PHE:CE1	1:B:467:HIS:CD2	3.06	0.45
1:B:427:PRO:HG2	1:B:467:HIS:HB3	1.98	0.45
1:B:314:ILE:HG13	1:B:314:ILE:O	2.17	0.44
1:B:143:GLU:HB3	1:B:144:PRO:HD2	1.98	0.44
1:A:281:ILE:HG12	1:A:383:PHE:CD1	2.51	0.44
1:B:443:THR:HA	1:B:491:LEU:HD21	1.98	0.44
1:A:98:TRP:CD2	1:A:113:PRO:HA	2.52	0.44
1:A:428:ILE:HG22	1:A:429:ALA:N	2.33	0.44
1:B:217:LEU:HD13	1:B:218:PRO:HD2	1.99	0.44
1:B:297:ARG:NH1	1:B:298:PRO:HD3	2.33	0.43
1:A:341:LYS:HA	1:A:344:LYS:HE2	2.00	0.43
1:A:271:PRO:O	1:A:272:ASN:C	2.56	0.43
1:A:443:THR:HA	1:A:491:LEU:CD2	2.40	0.43
1:B:293:VAL:HA	1:B:296:ILE:HD12	2.01	0.43
1:B:188:THR:HB	1:B:189:PRO:HD2	2.00	0.43
1:B:550:PRO:HG2	1:B:553:TYR:CD1	2.54	0.42
1:A:28:ILE:HD13	1:A:86:ILE:HG12	2.01	0.42
1:B:175:SER:O	1:B:179:ASN:HB2	2.19	0.42
1:A:211:ARG:HG3	1:B:519:TRP:CE3	2.52	0.42
1:A:102:ILE:HG21	1:A:104:ARG:HD2	2.02	0.42
1:B:194:TRP:O	1:B:197:HIS:CD2	2.71	0.42
1:B:309:PRO:O	1:B:460:VAL:HG23	2.20	0.42
1:A:550:PRO:HB2	1:A:552:GLN:HE21	1.80	0.42
1:B:552:GLN:HE21	1:B:552:GLN:H	1.66	0.42
1:B:185:VAL:HB	3:B:1562:EUG:C9	2.50	0.42
1:A:79:ASN:HD22	1:A:79:ASN:C	2.23	0.42
1:A:321:LEU:HD12	1:A:346:LEU:HD22	2.02	0.42
1:A:411:LEU:HD23	1:A:411:LEU:HA	1.82	0.42
1:B:151:LEU:HD23	1:B:166:LEU:HD22	2.01	0.42

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:425:PHE:HB2	1:A:488:MET:SD	2.60	0.42
1:A:179:ASN:OD1	1:A:184:GLY:HA3	2.20	0.42
1:A:456:GLY:HA2	1:A:468:ILE:O	2.20	0.41
1:A:459:THR:OG1	1:A:466:HIS:HB2	2.20	0.41
1:A:310:THR:HB	1:A:457:THR:CG2	2.49	0.41
1:A:129:ARG:NH2	1:A:131:LEU:HD21	2.35	0.41
1:A:201:GLU:OE1	1:A:264:LYS:NZ	2.40	0.41
1:A:282:THR:HG22	1:A:352:ASN:HD22	1.86	0.41
1:B:40:ILE:HD11	1:B:74:ILE:HD12	2.02	0.41
1:A:224:GLU:CD	1:A:224:GLU:H	2.23	0.41
1:B:297:ARG:CB	1:B:298:PRO:CD	2.92	0.41
1:B:319:ALA:HB3	1:B:413:TRP:HB2	2.02	0.41
1:A:13:PRO:HD3	1:A:117:GLY:O	2.21	0.41
1:B:350:ARG:HD2	1:B:350:ARG:HA	1.97	0.41
1:A:505:THR:CG2	1:A:513:ILE:HD12	2.51	0.41
1:B:177:LEU:HB2	1:B:265:ILE:HG22	2.03	0.41
1:B:278:SER:OG	1:B:399:ASP:OD1	2.33	0.41
1:A:211:ARG:CG	1:B:519:TRP:CZ3	2.98	0.41
1:A:79:ASN:ND2	1:A:79:ASN:C	2.74	0.41
1:B:452:LEU:HD11	1:B:471:ILE:HG23	2.03	0.41
1:A:244:TYR:OH	1:B:192:ASP:OD1	2.28	0.41
1:A:310:THR:CG2	1:A:459:THR:HG22	2.42	0.40
1:B:312:ARG:HD3	1:B:354:TYR:CE1	2.56	0.40
1:A:413:TRP:CH2	1:A:422:HIS:HD2	2.39	0.40
1:A:297:ARG:O	1:A:301:LEU:HB2	2.20	0.40
1:B:324:LYS:HG3	1:B:416:TRP:CZ2	2.56	0.40
1:B:214:MET:CE	1:B:245:GLY:HA2	2.52	0.40
1:B:419:ASN:HB2	1:B:474:ASN:OD1	2.21	0.40
1:B:238:ILE:HG22	1:B:238:ILE:O	2.21	0.40
1:A:283:LEU:HA	1:A:284:PRO:HD3	1.86	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	546/560 (98%)	492 (90%)	48 (9%)	6 (1%)	17	42
1	B	546/560 (98%)	514 (94%)	28 (5%)	4 (1%)	26	55
All	All	1092/1120 (98%)	1006 (92%)	76 (7%)	10 (1%)	21	49

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	231	GLU
1	B	418	PRO
1	A	159	ASN
1	B	105	ASN
1	A	274	ARG
1	A	418	PRO
1	A	361	GLU
1	B	345	GLN
1	B	169	PRO
1	A	230	PRO

### 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	470/482 (98%)	421 (90%)	49 (10%)	9	20
1	B	470/482 (98%)	424 (90%)	46 (10%)	10	23
All	All	940/964 (98%)	845 (90%)	95 (10%)	9	21

All (95) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	PHE
1	A	8	ARG
1	A	16	LEU

*Continued on next page...*



*Continued from previous page...*

Mol	Chain	Res	Type
1	A	28	ILE
1	A	41	SER
1	A	50	SER
1	A	79	ASN
1	A	94	SER
1	A	95	PHE
1	A	118	SER
1	A	128	ASN
1	A	129	ARG
1	A	142	VAL
1	A	149	HIS
1	A	156	GLU
1	A	161	ARG
1	A	177	LEU
1	A	217	LEU
1	A	229	LYS
1	A	231	GLU
1	A	237	LYS
1	A	264	LYS
1	A	297	ARG
1	A	303	MET
1	A	326	SER
1	A	329	SER
1	A	330	ARG
1	A	331	THR
1	A	341	LYS
1	A	344	LYS
1	A	381	VAL
1	A	391	GLU
1	A	393	SER
1	A	396	ARG
1	A	413	TRP
1	A	414	ILE
1	A	422	HIS
1	A	432	SER
1	A	443	THR
1	A	457	THR
1	A	467	HIS
1	A	472	VAL
1	A	475	LYS
1	A	503	TYR
1	A	504	ARG

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	A	505	THR
1	A	516	THR
1	A	523	SER
1	A	552	GLN
1	B	7	PHE
1	B	8	ARG
1	B	11	THR
1	B	25	ILE
1	B	26	GLN
1	B	35	GLU
1	B	63	VAL
1	B	79	ASN
1	B	95	PHE
1	B	128	ASN
1	B	142	VAL
1	B	149	HIS
1	B	162	ASP
1	B	163	LYS
1	B	177	LEU
1	B	211	ARG
1	B	217	LEU
1	B	221	LYS
1	B	231	GLU
1	B	237	LYS
1	B	241	LEU
1	B	297	ARG
1	B	301	LEU
1	B	326	SER
1	B	328	SER
1	B	329	SER
1	B	336	ASP
1	B	341	LYS
1	B	365	ARG
1	B	370	THR
1	B	396	ARG
1	B	403	GLN
1	B	409	ASP
1	B	412	LYS
1	B	422	HIS
1	B	430	LYS
1	B	432	SER
1	B	444	LYS

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	B	452	LEU
1	B	457	THR
1	B	503	TYR
1	B	504	ARG
1	B	513	ILE
1	B	516	THR
1	B	552	GLN
1	B	560	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	79	ASN
1	A	128	ASN
1	A	153	ASN
1	A	197	HIS
1	A	352	ASN
1	A	467	HIS
1	A	520	ASN
1	A	552	GLN
1	A	555	HIS
1	B	79	ASN
1	B	128	ASN
1	B	153	ASN
1	B	158	ASN
1	B	197	HIS
1	B	240	HIS
1	B	352	ASN
1	B	467	HIS
1	B	485	GLN
1	B	520	ASN
1	B	552	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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](#)

4 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	EUG	A	1562	-	11,11,12	1.84	1 (9%)	14,14,15	1.42	2 (14%)
2	FAD	A	600	1	48,58,58	1.48	7 (14%)	54,89,89	2.40	11 (20%)
2	FAD	B	1561	1	48,58,58	1.30	5 (10%)	54,89,89	2.34	8 (14%)
3	EUG	B	1562	-	11,11,12	1.69	1 (9%)	14,14,15	1.42	2 (14%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	EUG	A	1562	-	-	0/4/4/5	0/1/1/1
2	FAD	A	600	1	-	0/30/50/50	0/6/6/6
2	FAD	B	1561	1	-	0/30/50/50	0/6/6/6
3	EUG	B	1562	-	-	0/4/4/5	0/1/1/1

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	600	FAD	C1'-N10	2.15	1.50	1.48
2	A	600	FAD	C10-N1	2.29	1.39	1.35
2	A	600	FAD	C2A-N1A	2.39	1.38	1.33
2	B	1561	FAD	C2A-N1A	2.72	1.39	1.33
2	A	600	FAD	C4-N3	2.85	1.38	1.33
2	A	600	FAD	C4X-N5	3.08	1.38	1.33

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1561	FAD	C4X-N5	3.10	1.38	1.33
2	B	1561	FAD	C4-N3	3.13	1.38	1.33
2	B	1561	FAD	C1'-N10	3.21	1.51	1.48
2	A	600	FAD	C2A-N3A	3.80	1.38	1.32
2	B	1561	FAD	C2A-N3A	3.85	1.39	1.32
3	B	1562	EUG	C4-C3	5.19	1.48	1.40
2	A	600	FAD	O5B-C5B	5.35	1.66	1.44
3	A	1562	EUG	C4-C3	5.37	1.48	1.40

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1561	FAD	N3A-C2A-N1A	-12.62	119.23	128.89
2	A	600	FAD	N3A-C2A-N1A	-12.03	119.69	128.89
2	A	600	FAD	P-O3P-PA	-3.55	122.77	132.73
2	B	1561	FAD	C4X-C4-N3	-3.34	119.02	123.59
2	A	600	FAD	C4X-C4-N3	-3.08	119.38	123.59
2	B	1561	FAD	P-O3P-PA	-2.84	124.76	132.73
3	A	1562	EUG	O3-C3-C2	-2.32	120.25	124.21
2	A	600	FAD	O4'-C4'-C5'	-2.15	105.52	110.19
2	A	600	FAD	O3P-PA-O5B	-2.09	97.39	102.94
2	A	600	FAD	C8M-C8-C7	2.12	125.38	120.73
2	A	600	FAD	C1'-N10-C9A	2.17	121.29	118.86
2	A	600	FAD	O2A-PA-O5B	2.21	119.59	108.46
3	B	1562	EUG	C9-O3-C3	2.55	121.42	117.54
2	B	1561	FAD	C5X-C9A-N10	2.81	119.75	117.62
2	B	1561	FAD	C1'-N10-C9A	2.81	122.02	118.86
2	A	600	FAD	C4X-N5-C5X	2.94	120.14	116.76
2	B	1561	FAD	C4B-O4B-C1B	3.03	113.05	109.72
2	B	1561	FAD	C4X-N5-C5X	3.18	120.42	116.76
3	B	1562	EUG	O3-C3-C4	3.72	119.47	114.57
3	A	1562	EUG	O3-C3-C4	3.78	119.55	114.57
2	A	600	FAD	C5X-C9A-N10	4.65	121.15	117.62
2	B	1561	FAD	C4-N3-C2	6.86	121.18	115.25
2	A	600	FAD	C4-N3-C2	7.08	121.37	115.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	600	FAD	2	0
2	B	1561	FAD	2	0
3	B	1562	EUG	1	0

## 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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	550/560 (98%)	-0.38	4 (0%) 89 90	33, 42, 55, 66	0
1	B	550/560 (98%)	0.07	9 (1%) 74 75	44, 60, 90, 102	0
All	All	1100/1120 (98%)	-0.16	13 (1%) 81 81	33, 51, 81, 102	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	8	ARG	3.4
1	A	325	ARG	3.2
1	B	7	PHE	3.2
1	B	6	GLU	2.9
1	A	157	ALA	2.8
1	B	28	ILE	2.6
1	B	157	ALA	2.5
1	B	347	ASN	2.4
1	B	329	SER	2.4
1	B	303	MET	2.2
1	B	365	ARG	2.2
1	A	8	ARG	2.1
1	A	159	ASN	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 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
3	EUG	A	1562	11/12	0.87	0.29	4.87	64,65,67,67	0
3	EUG	B	1562	11/12	0.92	0.30	2.70	71,72,74,74	0
2	FAD	A	600	53/53	0.95	0.17	1.18	48,55,64,65	0
2	FAD	B	1561	53/53	0.97	0.16	-0.50	54,59,70,71	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.