



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 09:27 AM GMT

PDB ID : 3II4
Title : Structure of mycobacterial lipoamide dehydrogenase bound to a triazaspirodi
methoxybenzoyl inhibitor
Authors : Lima, C.D.
Deposited on : 2009-07-31
Resolution : 2.42 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
<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

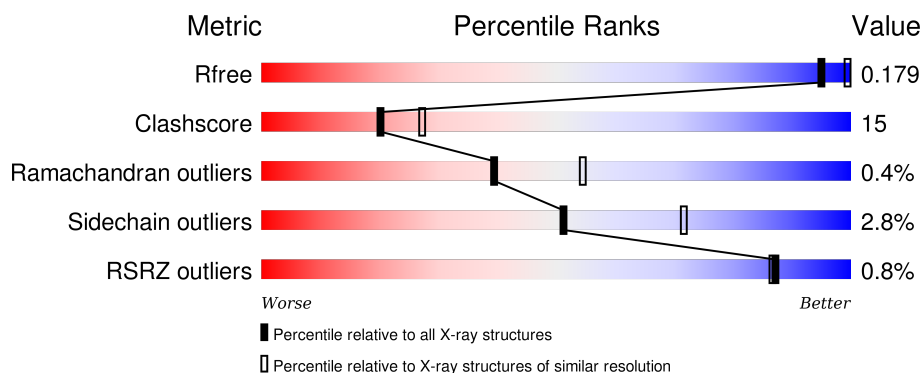
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.42 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	3386 (2.44-2.40)
Clashscore	102246	3897 (2.44-2.40)
Ramachandran outliers	100387	3837 (2.44-2.40)
Sidechain outliers	100360	3838 (2.44-2.40)
RSRZ outliers	91569	3396 (2.44-2.40)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 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	466	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 72%, green 26%, grey 3%);"></div> <div style="display: flex; justify-content: space-between; width: 90%; margin: 0 auto;"> % 72% 26% .. </div> </div>
1	B	466	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 1%, orange 1%, yellow 69%, green 29%, grey 2%);"></div> <div style="display: flex; justify-content: space-between; width: 90%; margin: 0 auto;"> % 69% 29% . </div> </div>

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 7873 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.

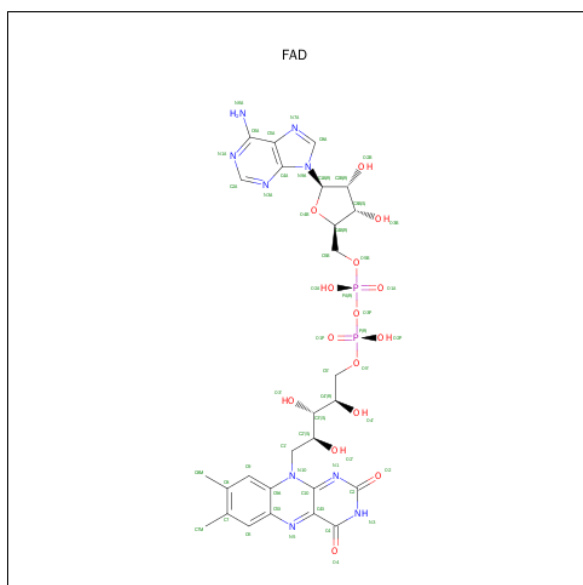
- Molecule 1 is a protein called Dihydrolipoyl dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	463	Total	C	N	O	S	0	0	0
			3461	2194	594	663	10			
1	B	464	Total	C	N	O	S	0	0	0
			3469	2199	595	664	11			

There are 4 discrepancies between the modelled and reference sequences:

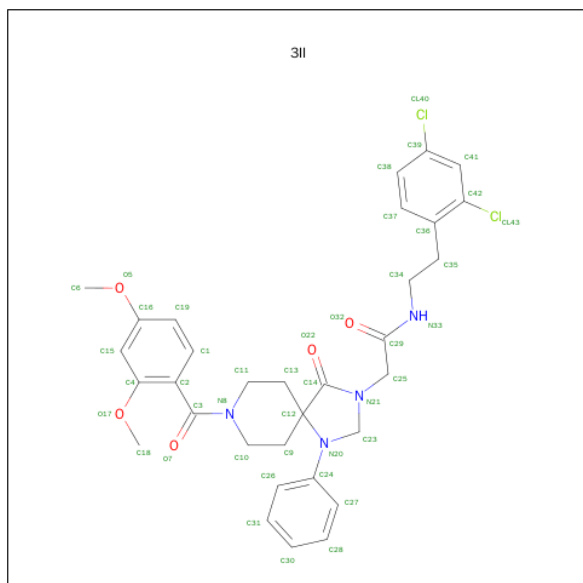
Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	EXPRESSION TAG	UNP P66004
A	0	SER	-	EXPRESSION TAG	UNP P66004
B	-1	GLY	-	EXPRESSION TAG	UNP P66004
B	0	SER	-	EXPRESSION TAG	UNP P66004

- 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		
2	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 3 is N-[2-(2,4-DICHLOROPHENYL)ETHYL]-2-{8-[(2,4-DIMETHOXYPHENYL)CARBONYL]-4-OXO-1-PHENYL-1,3,8-TRIAZASPIRO[4.5]DEC-3-YL}ACETAMIDE (three-letter code: 3II) (formula: C₃₂H₃₄Cl₂N₄O₅).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	Cl	N	O	0	0
			43	32	2	4	5		
3	B	1	Total	C	Cl	N	O	0	0
			43	32	2	4	5		

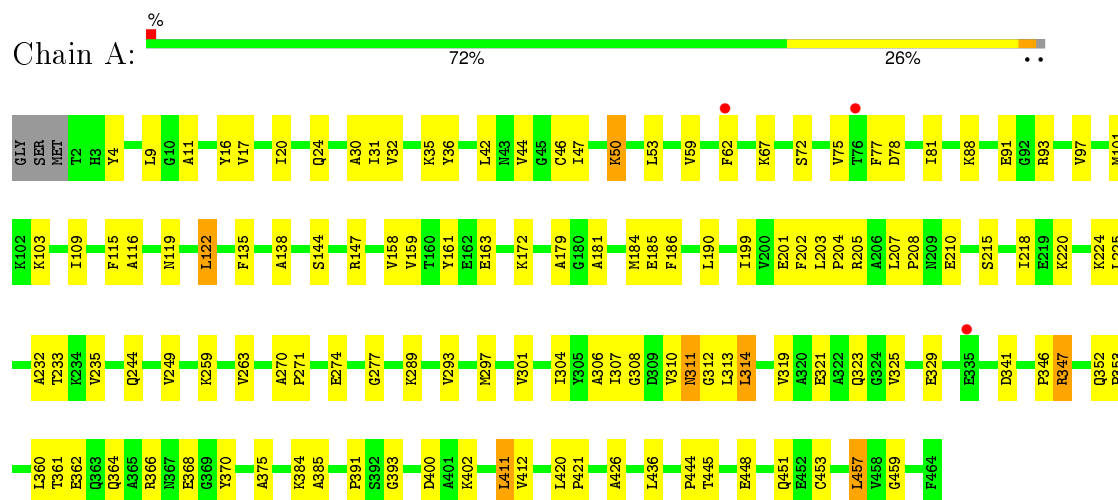
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	389	Total	O	0	0
			389	389		
4	B	362	Total	O	0	0
			362	362		

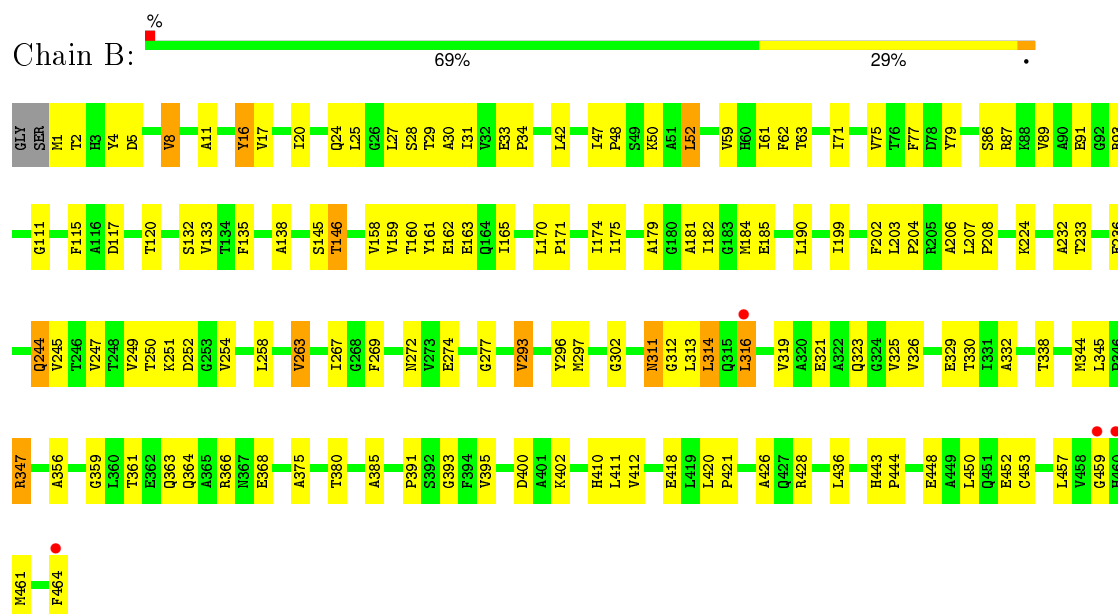
3 Residue-property plots

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: Dihydrolipoyl dehydrogenase



• Molecule 1: Dihydrolipoyl dehydrogenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	83.90 Å 98.40 Å 123.80 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.66 – 2.42 34.66 – 2.42	Depositor EDS
% Data completeness (in resolution range)	98.8 (29.66-2.42) 98.9 (34.66-2.42)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.43 (at 2.42 Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.185 , 0.242 0.189 , 0.179	Depositor DCC
R_{free} test set	1965 reflections (5.02%)	DCC
Wilson B-factor (Å ²)	28.7	Xtriage
Anisotropy	0.326	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 51.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Outliers	0 of 39116 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7873	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.59% 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, 0.375 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: 3II, 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	0.31	0/3523	0.59	0/4783
1	B	0.30	0/3531	0.58	0/4793
All	All	0.30	0/7054	0.59	0/9576

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3461	0	3461	98	0
1	B	3469	0	3473	116	0
2	A	53	0	31	2	0
2	B	53	0	31	1	0
3	A	43	0	34	2	0
3	B	43	0	34	2	0
4	A	389	0	0	6	0
4	B	362	0	0	10	0
All	All	7873	0	7064	209	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 (209) 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:314:LEU:HD22	1:B:316:LEU:HD22	1.53	0.91
1:A:364:GLN:O	1:A:368:GLU:HG2	1.80	0.81
1:A:400:ASP:OD2	1:A:402:LYS:HB2	1.81	0.80
1:B:224:LYS:NZ	1:B:224:LYS:HB2	1.98	0.77
1:B:311:ASN:C	1:B:311:ASN:HD22	1.89	0.75
1:A:50:LYS:HE3	1:A:53:LEU:HD12	1.69	0.73
1:A:244:GLN:HE22	1:A:259:LYS:NZ	1.88	0.72
1:A:311:ASN:HD22	1:A:311:ASN:C	1.93	0.70
1:B:311:ASN:ND2	1:B:313:LEU:H	1.89	0.70
1:B:91:GLU:HG2	4:B:692:HOH:O	1.90	0.69
1:A:311:ASN:HD22	1:A:312:GLY:N	1.91	0.69
1:A:72:SER:O	1:B:75:VAL:HG23	1.96	0.65
1:A:184:MET:CE	1:A:199:ILE:HG21	2.24	0.65
1:B:319:VAL:HG22	1:B:345:LEU:HD13	1.78	0.65
1:A:11:ALA:HA	1:A:31:ILE:HD11	1.77	0.65
1:A:20:ILE:O	1:A:24:GLN:HG3	1.97	0.64
1:A:361:THR:OG1	1:A:364:GLN:HG3	1.98	0.64
1:A:233:THR:HG22	1:A:249:VAL:CG2	2.28	0.64
1:B:50:LYS:N	1:B:50:LYS:HD2	2.13	0.63
1:B:20:ILE:O	1:B:24:GLN:HG3	1.99	0.63
1:B:11:ALA:HA	1:B:31:ILE:HD11	1.79	0.63
1:B:181:ALA:O	1:B:185:GLU:HG3	1.98	0.62
1:B:8:VAL:HG13	1:B:31:ILE:HA	1.82	0.62
1:B:361:THR:OG1	1:B:364:GLN:HG3	2.00	0.61
1:A:375:ALA:HB2	1:A:457:LEU:HD13	1.83	0.61
1:B:247:VAL:HG13	1:B:258:LEU:HB2	1.82	0.61
1:B:52:LEU:HD13	1:B:86:SER:HB3	1.83	0.60
1:A:426:ALA:HB2	1:A:436:LEU:HD21	1.82	0.60
1:B:338:THR:HG22	4:B:646:HOH:O	2.01	0.60
1:B:17:VAL:HG21	1:B:321:GLU:HG2	1.84	0.60
1:A:181:ALA:O	1:A:185:GLU:HG3	2.00	0.60
1:A:411:LEU:HD23	1:A:411:LEU:N	2.17	0.59
1:B:444:PRO:HA	1:B:448:GLU:OE2	2.03	0.59
1:B:393:GLY:HA3	1:B:412:VAL:O	2.03	0.58
1:B:224:LYS:HB2	1:B:224:LYS:HZ3	1.65	0.58
1:B:145:SER:OG	1:B:272:ASN:ND2	2.34	0.58
1:B:311:ASN:HD22	1:B:312:GLY:N	2.01	0.58
1:B:204:PRO:HD2	4:B:590:HOH:O	2.03	0.58
1:A:352:GLN:OE1	1:A:353:PRO:HA	2.04	0.58
1:B:146:THR:HG21	1:B:160:THR:CG2	2.34	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:158:VAL:HA	1:A:263:VAL:HG13	1.85	0.57
1:A:451:GLN:HG3	4:A:740:HOH:O	2.03	0.57
1:A:233:THR:HG22	1:A:249:VAL:HG21	1.86	0.57
1:B:146:THR:HG22	4:B:497:HOH:O	2.03	0.57
1:B:174:ILE:HG21	1:B:190:LEU:HD23	1.87	0.57
1:A:159:VAL:HB	1:A:163:GLU:HB2	1.87	0.57
1:A:44:VAL:HG23	1:A:161:TYR:HE1	1.69	0.57
1:B:5:ASP:H	1:B:28:SER:HB3	1.71	0.56
1:B:42:LEU:HA	1:B:47:ILE:CG1	2.35	0.56
1:B:366:ARG:HG3	1:B:366:ARG:HH11	1.69	0.56
1:B:311:ASN:HD22	1:B:313:LEU:H	1.52	0.55
1:A:88:LYS:HA	1:A:91:GLU:HG2	1.88	0.55
1:A:184:MET:CE	1:A:199:ILE:HD13	2.37	0.55
1:B:1:MET:HE3	1:B:132:SER:HB3	1.90	0.54
1:A:78:ASP:OD2	1:A:81:ILE:HG12	2.07	0.54
1:A:385:ALA:HB1	1:A:391:PRO:HA	1.89	0.54
1:B:61:ILE:C	1:B:63:THR:H	2.11	0.54
1:A:59:VAL:HG22	1:A:77:PHE:CD2	2.42	0.54
1:A:444:PRO:HA	1:A:448:GLU:OE2	2.07	0.54
1:B:267:ILE:HG12	4:B:504:HOH:O	2.08	0.53
1:A:274:GLU:HG2	4:A:764:HOH:O	2.08	0.53
1:A:360:LEU:HD13	1:A:368:GLU:HG3	1.89	0.53
1:B:175:ILE:HD13	1:B:247:VAL:HG11	1.91	0.53
1:B:344:MET:CE	1:B:359:GLY:HA2	2.39	0.52
1:B:175:ILE:HD13	1:B:247:VAL:CG1	2.39	0.52
1:A:453:CYS:O	1:A:457:LEU:HD22	2.10	0.52
1:A:319:VAL:O	1:A:323:GLN:HG3	2.10	0.52
1:A:17:VAL:HG21	1:A:321:GLU:HG2	1.92	0.51
1:A:421:PRO:HD3	4:B:752:HOH:O	2.09	0.51
1:B:236:GLU:HG3	1:B:250:THR:HG23	1.92	0.51
1:B:47:ILE:HB	1:B:48:PRO:CD	2.40	0.51
1:B:400:ASP:OD2	1:B:402:LYS:HE2	2.11	0.51
1:A:67:LYS:HD3	4:A:552:HOH:O	2.09	0.51
1:A:445:THR:HB	1:B:418:GLU:OE1	2.10	0.51
1:B:375:ALA:CB	1:B:457:LEU:HD13	2.40	0.51
1:B:274:GLU:HG2	4:B:760:HOH:O	2.10	0.51
1:A:184:MET:HE1	1:A:199:ILE:HD13	1.93	0.51
1:A:360:LEU:CD1	1:A:368:GLU:HG3	2.41	0.51
1:B:232:ALA:O	1:B:251:LYS:HG3	2.11	0.50
1:A:50:LYS:HG2	1:B:444:PRO:CG	2.41	0.50
1:A:314:LEU:HD23	3:A:465:3II:H9A	1.92	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:179:ALA:HB2	1:A:201:GLU:HB2	1.94	0.50
1:A:311:ASN:ND2	1:A:313:LEU:H	2.10	0.50
1:B:47:ILE:CG2	1:B:89:VAL:HG12	2.42	0.50
1:A:97:VAL:O	1:A:101:MET:HG2	2.12	0.50
1:B:313:LEU:HD12	1:B:323:GLN:HE22	1.77	0.49
1:B:296:TYR:HB3	1:B:330:THR:HG21	1.93	0.49
1:B:224:LYS:HB2	1:B:224:LYS:HZ2	1.75	0.49
1:B:16:TYR:O	1:B:20:ILE:HG13	2.12	0.49
1:B:27:LEU:HD11	1:B:332:ALA:HB2	1.95	0.49
1:B:171:PRO:HD3	4:B:706:HOH:O	2.11	0.49
1:B:179:ALA:HB1	1:B:184:MET:HE3	1.95	0.49
1:A:233:THR:HG22	1:A:249:VAL:HG22	1.94	0.49
1:A:301:VAL:HG11	1:A:304:ILE:HD12	1.95	0.48
1:A:362:GLU:HG2	1:A:366:ARG:HH11	1.78	0.48
1:B:244:GLN:C	1:B:244:GLN:HE21	2.16	0.48
1:B:158:VAL:HG22	1:B:263:VAL:CG1	2.44	0.48
1:B:52:LEU:CD1	1:B:86:SER:HB3	2.42	0.48
1:A:9:LEU:HD22	2:A:480:FAD:N1A	2.29	0.48
1:A:103:LYS:HE2	1:B:464:PHE:O	2.14	0.48
1:A:325:VAL:O	1:A:329:GLU:HG3	2.14	0.48
1:A:384:LYS:HD3	1:A:444:PRO:O	2.13	0.48
1:B:428:ARG:HB2	1:B:428:ARG:NH1	2.28	0.48
1:A:220:LYS:O	1:A:224:LYS:HG2	2.14	0.48
1:A:297:MET:HG3	1:A:323:GLN:HB3	1.96	0.47
1:B:146:THR:HG21	1:B:160:THR:HG21	1.97	0.47
1:A:393:GLY:HA3	1:A:412:VAL:O	2.15	0.47
1:B:380:THR:HG23	4:B:469:HOH:O	2.13	0.47
1:A:205:ARG:NH2	1:A:210:GLU:O	2.48	0.47
1:B:459:GLY:HA2	4:B:737:HOH:O	2.14	0.47
1:A:235:VAL:HA	1:A:249:VAL:HG23	1.95	0.47
1:A:420:LEU:N	1:A:421:PRO:CD	2.78	0.47
1:A:207:LEU:N	1:A:208:PRO:CD	2.78	0.47
1:B:233:THR:HG22	1:B:249:VAL:HB	1.97	0.46
1:B:395:VAL:HG11	1:B:450:LEU:HD23	1.97	0.46
1:A:271:PRO:HB3	1:A:310:VAL:HA	1.98	0.46
1:B:375:ALA:HB2	1:B:457:LEU:HD13	1.98	0.46
1:B:184:MET:CE	1:B:199:ILE:HG21	2.45	0.46
1:A:293:VAL:CG1	1:A:306:ALA:HB3	2.46	0.46
1:B:42:LEU:HD13	1:B:93:ARG:HB3	1.98	0.46
1:B:87:ARG:O	1:B:91:GLU:HG3	2.15	0.45
1:A:311:ASN:C	1:A:311:ASN:ND2	2.65	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:77:PHE:CD2	1:B:71:ILE:HG12	2.51	0.45
1:B:117:ASP:HB3	1:B:120:THR:HB	1.97	0.45
1:A:116:ALA:HB2	1:A:122:LEU:HD13	1.98	0.45
1:B:161:TYR:CD1	1:B:162:GLU:N	2.85	0.45
1:B:4:TYR:CE2	1:B:30:ALA:HB2	2.51	0.45
1:A:158:VAL:HG13	1:A:263:VAL:HG13	1.98	0.45
1:B:244:GLN:C	1:B:244:GLN:NE2	2.71	0.45
1:B:363:GLN:HA	1:B:363:GLN:OE1	2.17	0.45
1:B:420:LEU:N	1:B:421:PRO:CD	2.80	0.45
1:B:25:LEU:HD13	1:B:332:ALA:HB3	1.99	0.44
1:A:179:ALA:CB	1:A:201:GLU:HB2	2.48	0.44
1:B:293:VAL:HG22	1:B:297:MET:C	2.36	0.44
1:A:42:LEU:HD13	1:A:93:ARG:HB3	2.00	0.44
1:A:35:LYS:HE2	1:A:36:TYR:CZ	2.53	0.44
1:B:135:PHE:CE1	1:B:138:ALA:HB2	2.52	0.44
1:B:325:VAL:O	1:B:329:GLU:HG3	2.17	0.44
1:A:346:PRO:C	1:A:347:ARG:HG3	2.37	0.44
1:A:172:LYS:HD2	4:A:478:HOH:O	2.18	0.44
1:B:364:GLN:O	1:B:368:GLU:HG2	2.17	0.44
1:A:116:ALA:HB2	1:A:122:LEU:HD22	2.00	0.44
1:B:162:GLU:O	1:B:165:ILE:HG22	2.17	0.44
3:B:465:3II:H27	3:B:465:3II:H9	1.99	0.44
1:A:289:LYS:HE3	4:A:639:HOH:O	2.16	0.44
1:B:42:LEU:HA	1:B:47:ILE:HG13	2.00	0.44
1:A:32:VAL:HG22	1:A:109:ILE:HB	2.00	0.44
1:A:244:GLN:HE22	1:A:259:LYS:HZ2	1.59	0.44
1:A:307:ILE:O	1:A:310:VAL:HG12	2.18	0.44
1:B:34:PRO:HB3	1:B:111:GLY:C	2.38	0.44
1:B:42:LEU:HA	1:B:47:ILE:HG12	1.99	0.43
1:B:2:THR:O	1:B:133:VAL:HG13	2.18	0.43
1:A:50:LYS:HG2	1:B:444:PRO:HG2	2.00	0.43
1:A:368:GLU:HB2	1:A:370:TYR:CD1	2.52	0.43
1:B:115:PHE:HB2	1:B:277:GLY:O	2.18	0.43
1:A:44:VAL:CG2	1:A:161:TYR:HE1	2.30	0.43
1:A:308:GLY:HA3	2:A:480:FAD:O2P	2.19	0.43
1:B:59:VAL:HG22	1:B:77:PHE:CD2	2.52	0.43
1:A:184:MET:HE1	1:A:199:ILE:CD1	2.48	0.43
1:B:47:ILE:HG23	1:B:89:VAL:HG12	1.99	0.43
1:A:203:LEU:HB3	1:A:204:PRO:HD2	2.00	0.43
1:A:119:ASN:HA	1:A:135:PHE:CE1	2.54	0.43
1:B:61:ILE:O	1:B:63:THR:N	2.52	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:453:CYS:O	1:B:457:LEU:HB2	2.19	0.42
1:B:356:ALA:HB1	1:B:420:LEU:HD11	2.01	0.42
1:B:236:GLU:CG	1:B:250:THR:HG23	2.48	0.42
1:A:4:TYR:CE2	1:A:30:ALA:HB2	2.54	0.42
1:B:326:VAL:O	1:B:330:THR:HG23	2.20	0.42
1:B:159:VAL:HB	1:B:163:GLU:HB2	2.01	0.42
1:B:443:HIS:CE1	1:B:444:PRO:HB3	2.54	0.42
1:B:247:VAL:CG1	1:B:258:LEU:HB2	2.50	0.42
1:B:203:LEU:HD12	1:B:208:PRO:HG3	2.01	0.42
1:B:314:LEU:HB3	1:B:319:VAL:HG21	2.02	0.42
1:B:361:THR:HA	1:B:410:HIS:NE2	2.35	0.42
1:B:411:LEU:HD12	1:B:411:LEU:N	2.35	0.42
1:A:135:PHE:CE1	1:A:138:ALA:HB2	2.55	0.41
1:B:252:ASP:O	1:B:254:VAL:HG23	2.20	0.41
1:B:457:LEU:HA	1:B:457:LEU:HD12	1.92	0.41
1:B:77:PHE:N	1:B:77:PHE:CD1	2.88	0.41
1:A:62:PHE:CE2	1:A:75:VAL:HG22	2.55	0.41
1:B:385:ALA:HB1	1:B:391:PRO:HA	2.02	0.41
1:A:186:PHE:O	1:A:190:LEU:HG	2.19	0.41
1:A:426:ALA:CB	1:A:436:LEU:HD21	2.48	0.41
1:A:147:ARG:HE	3:A:465:3II:H38	1.85	0.41
1:A:362:GLU:O	1:A:366:ARG:HG3	2.20	0.41
1:B:244:GLN:HE21	1:B:245:VAL:N	2.18	0.41
1:B:29:THR:HG22	1:B:30:ALA:N	2.34	0.41
1:A:144:SER:HB2	1:A:270:ALA:O	2.20	0.41
1:A:59:VAL:HG22	1:A:77:PHE:CG	2.56	0.41
1:B:202:PHE:O	1:B:232:ALA:HA	2.19	0.41
1:A:202:PHE:O	1:A:232:ALA:HA	2.20	0.41
1:B:1:MET:CE	1:B:120:THR:HG23	2.51	0.41
1:B:1:MET:HE1	1:B:120:THR:HG23	2.01	0.41
1:B:311:ASN:C	1:B:311:ASN:ND2	2.62	0.41
1:A:50:LYS:HE3	1:A:53:LEU:CD1	2.45	0.41
1:A:184:MET:HE2	1:A:199:ILE:HD13	2.03	0.41
1:A:42:LEU:HA	1:A:47:ILE:HG12	2.02	0.41
1:A:4:TYR:O	1:A:135:PHE:HA	2.21	0.41
1:B:182:ILE:HD12	1:B:269:PHE:CE1	2.56	0.41
1:B:206:ALA:O	1:B:207:LEU:C	2.58	0.41
1:B:347:ARG:HB2	3:B:465:3II:H18B	2.03	0.40
1:B:79:TYR:CD1	1:B:170:LEU:HD11	2.56	0.40
1:B:426:ALA:HB2	1:B:436:LEU:HD21	2.03	0.40
1:A:215:SER:O	1:A:218:ILE:HG22	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:33:GLU:OE2	2:B:580:FAD:H1B	2.21	0.40
1:B:452:GLU:OE2	1:B:461:MET:HB3	2.21	0.40
1:A:158:VAL:HG22	1:A:263:VAL:CG1	2.52	0.40
1:A:115:PHE:HB2	1:A:277:GLY:O	2.22	0.40
1:A:341:ASP:HA	4:A:550:HOH:O	2.21	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	Percentiles	
1	A	461/466 (99%)	434 (94%)	25 (5%)	2 (0%)	39	54
1	B	462/466 (99%)	428 (93%)	32 (7%)	2 (0%)	39	54
All	All	923/932 (99%)	862 (93%)	57 (6%)	4 (0%)	39	54

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	62	PHE
1	B	302	GLY
1	A	46	CYS
1	A	459	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	A	359/361 (99%)	350 (98%)	9 (2%)	55	75
1	B	360/361 (100%)	349 (97%)	11 (3%)	47	68
All	All	719/722 (100%)	699 (97%)	20 (3%)	51	71

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

Mol	Chain	Res	Type
1	A	16	TYR
1	A	50	LYS
1	A	122	LEU
1	A	225	LEU
1	A	311	ASN
1	A	314	LEU
1	A	347	ARG
1	A	411	LEU
1	A	457	LEU
1	B	8	VAL
1	B	16	TYR
1	B	52	LEU
1	B	146	THR
1	B	244	GLN
1	B	263	VAL
1	B	293	VAL
1	B	311	ASN
1	B	314	LEU
1	B	316	LEU
1	B	347	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	244	GLN
1	A	272	ASN
1	A	311	ASN
1	A	403	HIS
1	A	439	ASN
1	A	451	GLN
1	A	455	HIS
1	B	244	GLN
1	B	272	ASN
1	B	311	ASN

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Mol	Chain	Res	Type
1	B	439	ASN
1	B	455	HIS
1	B	463	ASN

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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

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	3II	A	465	-	46,47,47	3.44	31 (67%)	60,67,67	1.76	16 (26%)
2	FAD	A	480	-	48,58,58	1.70	8 (16%)	54,89,89	1.96	6 (11%)
3	3II	B	465	-	46,47,47	3.40	31 (67%)	60,67,67	1.78	16 (26%)
2	FAD	B	580	-	48,58,58	1.62	8 (16%)	54,89,89	2.02	6 (11%)

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	3II	A	465	-	-	0/26/58/58	0/5/5/5
2	FAD	A	480	-	-	0/30/50/50	0/6/6/6
3	3II	B	465	-	-	1/26/58/58	0/5/5/5
2	FAD	B	580	-	-	0/30/50/50	0/6/6/6

All (78) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	465	3II	C25-C29	-10.20	1.36	1.52
3	B	465	3II	C25-C29	-9.44	1.37	1.52
3	B	465	3II	C34-N33	-3.24	1.38	1.46
3	A	465	3II	C34-N33	-3.23	1.38	1.46
3	B	465	3II	C35-C36	2.12	1.55	1.51
2	B	580	FAD	C9-C8	2.13	1.43	1.37
3	B	465	3II	C10-N8	2.15	1.50	1.47
2	A	480	FAD	C9-C8	2.18	1.43	1.37
3	A	465	3II	C3-N8	2.18	1.39	1.34
3	A	465	3II	C30-C28	2.22	1.43	1.38
3	A	465	3II	C10-N8	2.28	1.50	1.47
2	A	480	FAD	C4A-N3A	2.36	1.39	1.35
2	A	480	FAD	C2A-N1A	2.39	1.38	1.33
3	B	465	3II	C3-N8	2.40	1.40	1.34
2	B	580	FAD	C5X-N5	2.40	1.39	1.35
3	A	465	3II	C35-C36	2.41	1.56	1.51
3	B	465	3II	C30-C28	2.41	1.44	1.38
2	B	580	FAD	C2A-N1A	2.49	1.38	1.33
3	A	465	3II	C30-C31	2.59	1.44	1.38
3	A	465	3II	C1-C19	2.60	1.43	1.38
2	A	480	FAD	C5X-N5	2.65	1.39	1.35
3	B	465	3II	C30-C31	2.67	1.44	1.38
2	B	580	FAD	C4A-N3A	2.70	1.39	1.35
3	B	465	3II	C25-N21	2.70	1.49	1.45
3	A	465	3II	C2-C3	2.75	1.55	1.50
3	A	465	3II	C25-N21	2.77	1.49	1.45
3	A	465	3II	C28-C27	2.80	1.44	1.38
3	B	465	3II	C28-C27	2.82	1.44	1.38
3	B	465	3II	C1-C19	2.83	1.43	1.38
3	B	465	3II	C2-C3	2.88	1.55	1.50
3	B	465	3II	C31-C26	2.98	1.45	1.38
3	A	465	3II	C31-C26	3.07	1.45	1.38
3	B	465	3II	C24-N20	3.12	1.49	1.41
2	B	580	FAD	C9A-N10	3.19	1.43	1.38
3	A	465	3II	C24-N20	3.35	1.49	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	465	3II	C19-C16	3.36	1.45	1.38
3	B	465	3II	C12-N20	3.53	1.56	1.47
3	B	465	3II	C15-C4	3.53	1.45	1.38
3	B	465	3II	C19-C16	3.53	1.45	1.38
2	A	480	FAD	C9A-N10	3.54	1.43	1.38
3	A	465	3II	C2-C4	3.58	1.47	1.40
3	A	465	3II	C37-C38	3.58	1.45	1.38
3	B	465	3II	C15-C16	3.61	1.45	1.38
3	A	465	3II	C12-N20	3.63	1.57	1.47
3	A	465	3II	C38-C39	3.64	1.45	1.38
3	B	465	3II	C9-C12	3.65	1.58	1.54
3	A	465	3II	C1-C2	3.66	1.46	1.39
3	A	465	3II	C15-C16	3.66	1.45	1.38
3	B	465	3II	C1-C2	3.67	1.46	1.39
3	A	465	3II	C15-C4	3.68	1.45	1.38
2	B	580	FAD	C4-N3	3.71	1.40	1.33
3	B	465	3II	C38-C39	3.73	1.45	1.38
3	B	465	3II	C37-C38	3.78	1.45	1.38
3	B	465	3II	C2-C4	4.04	1.48	1.40
3	A	465	3II	C9-C12	4.05	1.58	1.54
3	A	465	3II	C41-C39	4.09	1.45	1.38
2	A	480	FAD	C4-N3	4.12	1.40	1.33
3	B	465	3II	C41-C39	4.16	1.45	1.38
3	B	465	3II	C14-N21	4.22	1.39	1.34
3	A	465	3II	C37-C36	4.32	1.47	1.39
3	B	465	3II	C37-C36	4.33	1.47	1.39
3	B	465	3II	C41-C42	4.40	1.45	1.38
2	B	580	FAD	C10-N10	4.56	1.44	1.39
3	A	465	3II	C41-C42	4.57	1.46	1.38
3	A	465	3II	C13-C12	4.64	1.59	1.54
3	A	465	3II	C27-C24	4.66	1.48	1.39
3	B	465	3II	C26-C24	4.67	1.48	1.39
3	A	465	3II	C26-C24	4.71	1.48	1.39
3	B	465	3II	C27-C24	4.74	1.48	1.39
3	A	465	3II	C14-N21	4.92	1.40	1.34
3	B	465	3II	C13-C12	5.06	1.60	1.54
3	B	465	3II	C42-C36	5.49	1.49	1.39
2	A	480	FAD	C10-N10	5.51	1.45	1.39
3	A	465	3II	C42-C36	5.59	1.49	1.39
2	A	480	FAD	C4X-N5	5.63	1.42	1.33
2	B	580	FAD	C4X-N5	5.72	1.42	1.33
3	A	465	3II	O5-C6	6.01	1.61	1.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	465	3II	O5-C6	6.15	1.61	1.42

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	480	FAD	C4X-C4-N3	-5.63	115.89	123.59
2	B	580	FAD	C4X-C10-N10	-5.60	117.22	120.52
2	A	480	FAD	C4X-C10-N10	-5.37	117.36	120.52
2	B	580	FAD	C4X-C4-N3	-5.30	116.34	123.59
2	B	580	FAD	C4-C4X-C10	-4.23	117.24	119.94
3	A	465	3II	C13-C12-N20	-3.90	109.14	112.71
3	B	465	3II	C15-C4-C2	-3.63	114.75	121.40
3	A	465	3II	C15-C4-C2	-3.60	114.80	121.40
3	B	465	3II	C25-N21-C23	-3.47	116.12	122.44
2	A	480	FAD	C4-C4X-C10	-3.42	117.75	119.94
3	B	465	3II	C13-C12-N20	-2.94	110.02	112.71
3	A	465	3II	C25-N21-C23	-2.90	117.16	122.44
2	B	580	FAD	N3A-C2A-N1A	-2.89	126.68	128.89
3	B	465	3II	C9-C12-N20	-2.82	110.13	112.71
2	A	480	FAD	N3A-C2A-N1A	-2.71	126.82	128.89
3	B	465	3II	O22-C14-N21	-2.63	123.75	125.84
3	B	465	3II	C27-C24-C26	-2.53	113.86	119.13
3	A	465	3II	C27-C24-C26	-2.48	113.98	119.13
3	A	465	3II	O22-C14-N21	-2.36	123.97	125.84
3	B	465	3II	C19-C16-C15	-2.23	117.44	120.56
3	A	465	3II	C19-C16-C15	-2.15	117.55	120.56
3	A	465	3II	C18-O17-C4	2.13	120.77	117.54
3	A	465	3II	C31-C26-C24	2.13	122.58	119.69
3	B	465	3II	C31-C26-C24	2.43	122.98	119.69
3	B	465	3II	C18-O17-C4	2.50	121.34	117.54
3	A	465	3II	C2-C3-N8	2.61	121.21	118.29
3	B	465	3II	C27-C24-N20	2.70	126.15	121.10
3	B	465	3II	C28-C27-C24	2.71	123.36	119.69
3	A	465	3II	C27-C24-N20	2.80	126.33	121.10
3	A	465	3II	C28-C27-C24	2.82	123.51	119.69
3	A	465	3II	C34-C35-C36	2.83	120.28	113.02
3	B	465	3II	C35-C34-N33	2.83	118.87	111.97
3	B	465	3II	C34-C35-C36	2.87	120.40	113.02
3	B	465	3II	C1-C2-C4	2.95	122.03	118.24
2	B	580	FAD	C1'-N10-C9A	3.03	122.27	118.86
2	A	480	FAD	C1'-N10-C9A	3.04	122.28	118.86
3	A	465	3II	C1-C2-C4	3.04	122.16	118.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	465	3II	C4-C15-C16	3.07	124.75	118.91
3	B	465	3II	C4-C15-C16	3.23	125.05	118.91
3	A	465	3II	C35-C34-N33	3.25	119.91	111.97
3	A	465	3II	O17-C4-C2	3.97	121.43	116.54
3	B	465	3II	O17-C4-C2	4.42	121.98	116.54
2	A	480	FAD	C4-N3-C2	9.23	123.23	115.25
2	B	580	FAD	C4-N3-C2	9.47	123.43	115.25

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	465	3II	C25-C29-N33-C34

There are no ring outliers.

4 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	465	3II	2	0
2	A	480	FAD	2	0
3	B	465	3II	2	0
2	B	580	FAD	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, 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	463/466 (99%)	-0.44	3 (0%) 90 90	15, 27, 43, 63	0
1	B	464/466 (99%)	-0.37	4 (0%) 85 85	16, 27, 43, 62	0
All	All	927/932 (99%)	-0.40	7 (0%) 87 87	15, 27, 43, 63	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	62	PHE	2.3
1	A	335	GLU	2.3
1	B	459	GLY	2.2
1	B	464	PHE	2.2
1	B	460	HIS	2.1
1	B	316	LEU	2.0
1	A	76	THR	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, 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	LLDF	B-factors(\AA^2)	Q<0.9
3	3II	A	465	43/43	0.82	0.21	1.33	31,35,48,48	0
3	3II	B	465	43/43	0.81	0.21	1.17	37,44,51,54	0
2	FAD	A	480	53/53	0.97	0.15	0.37	16,19,22,24	0
2	FAD	B	580	53/53	0.97	0.14	0.05	14,18,22,24	0

6.5 Other polymers [i](#)

There are no such residues in this entry.