



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 24, 2017 – 10:57 PM EDT

PDB ID : 3HGO
Title : Crystal structure of the F74Y/H244Y OPR3 double mutant from tomato
Authors : Clausen, T.; Breithaupt, C.
Deposited on : unknown
Resolution : 2.30 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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20030345
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
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) : rb-20030345

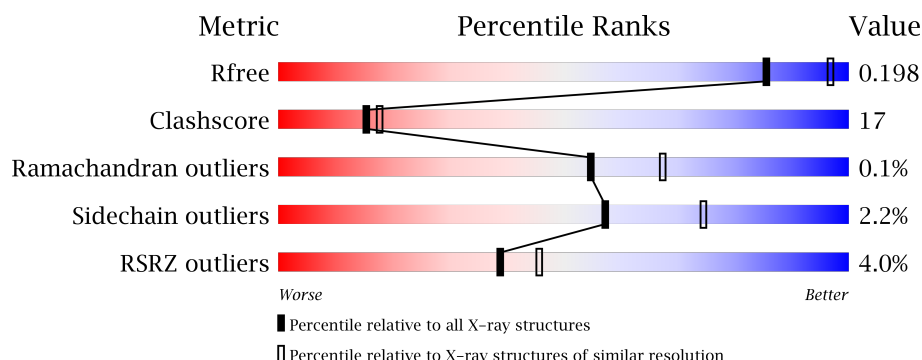
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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}	100719	4130 (2.30-2.30)
Clashscore	112137	4751 (2.30-2.30)
Ramachandran outliers	110173	4705 (2.30-2.30)
Sidechain outliers	110143	4704 (2.30-2.30)
RSRZ outliers	101464	4156 (2.30-2.30)

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	402	<div> <div>3%</div> <div>62%</div> <div>28%</div> <div>9%</div> </div>
1	B	402	<div> <div>4%</div> <div>66%</div> <div>24%</div> <div>10%</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6217 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.

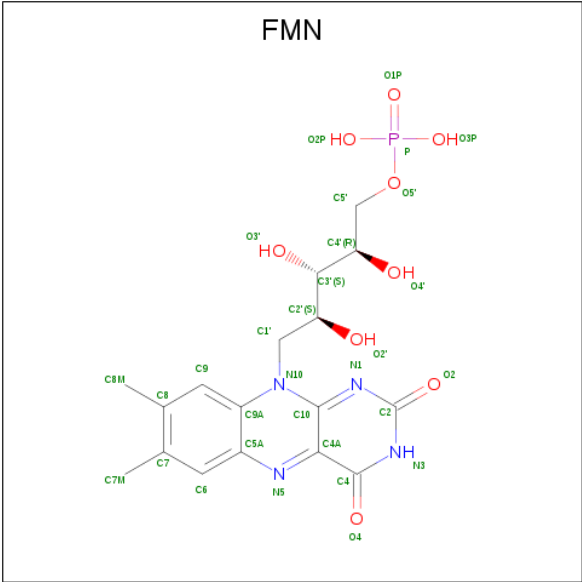
- Molecule 1 is a protein called 12-oxophytodienoate reductase 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	367	Total	C	N	O	S	0	0	0
			2838	1803	503	521	11			
1	B	363	Total	C	N	O	S	0	0	0
			2816	1791	498	516	11			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	HIS	-	EXPRESSION TAG	UNP Q9FEW9
A	-4	HIS	-	EXPRESSION TAG	UNP Q9FEW9
A	-3	HIS	-	EXPRESSION TAG	UNP Q9FEW9
A	-2	HIS	-	EXPRESSION TAG	UNP Q9FEW9
A	-1	HIS	-	EXPRESSION TAG	UNP Q9FEW9
A	0	HIS	-	EXPRESSION TAG	UNP Q9FEW9
A	74	TYR	PHE	ENGINEERED	UNP Q9FEW9
A	244	TYR	HIS	ENGINEERED	UNP Q9FEW9
B	-5	HIS	-	EXPRESSION TAG	UNP Q9FEW9
B	-4	HIS	-	EXPRESSION TAG	UNP Q9FEW9
B	-3	HIS	-	EXPRESSION TAG	UNP Q9FEW9
B	-2	HIS	-	EXPRESSION TAG	UNP Q9FEW9
B	-1	HIS	-	EXPRESSION TAG	UNP Q9FEW9
B	0	HIS	-	EXPRESSION TAG	UNP Q9FEW9
B	74	TYR	PHE	ENGINEERED	UNP Q9FEW9
B	244	TYR	HIS	ENGINEERED	UNP Q9FEW9

- Molecule 2 is FLAVIN MONONUCLEOTIDE (three-letter code: FMN) (formula: C₁₇H₂₁N₄O₉P).



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	49.14Å 92.86Å 89.62Å 90.00° 98.08° 90.00°	Depositor
Resolution (Å)	20.00 – 2.30 19.77 – 2.30	Depositor EDS
% Data completeness (in resolution range)	91.4 (20.00-2.30) 91.6 (19.77-2.30)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.14 (at 2.30Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.200 , 0.259 0.204 , 0.198	Depositor DCC
R_{free} test set	1622 reflections (5.01%)	DCC
Wilson B-factor (Å ²)	31.6	Xtriage
Anisotropy	0.400	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 52.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6217	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.54% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: FMN

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	0.33	0/2902	0.58	0/3938
1	B	0.35	0/2880	0.61	0/3907
All	All	0.34	0/5782	0.59	0/7845

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	2838	0	2807	100	0
1	B	2816	0	2795	89	0
2	A	31	0	19	3	0
2	B	31	0	19	1	0
3	A	225	0	0	13	0
3	B	276	0	0	22	0
All	All	6217	0	5640	188	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 17.

All (188) 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:253:LEU:HD13	1:B:304:LEU:HD22	1.39	1.05
1:B:267:GLN:NE2	1:B:274:LEU:H	1.69	0.90
1:A:348:ASN:HD21	1:A:362:ASN:H	1.21	0.88
1:B:267:GLN:HE21	1:B:274:LEU:H	1.21	0.87
1:B:295:LEU:HD23	1:B:298:GLU:HG2	1.58	0.86
1:A:267:GLN:NE2	1:A:274:LEU:H	1.73	0.85
1:A:295:LEU:HD23	1:A:298:GLU:HG2	1.64	0.80
1:B:19:GLY:HA3	3:B:539:HOH:O	1.81	0.80
1:B:348:ASN:HD21	1:B:362:ASN:H	1.32	0.76
1:A:8:GLY:C	1:A:10:ASN:H	1.89	0.75
1:B:253:LEU:HD13	1:B:304:LEU:CD2	2.17	0.75
1:B:325:GLU:HG2	3:B:504:HOH:O	1.86	0.73
1:A:368:THR:HG23	3:A:397:HOH:O	1.88	0.73
1:B:368:THR:HG23	3:B:404:HOH:O	1.88	0.72
1:A:243:ASP:HB2	3:A:559:HOH:O	1.90	0.71
1:A:159:GLU:HA	1:A:162:GLN:HE21	1.55	0.71
1:B:17:LYS:HD2	3:B:523:HOH:O	1.91	0.70
1:B:295:LEU:HD13	3:B:627:HOH:O	1.90	0.70
1:B:184:ILE:O	1:B:184:ILE:HG13	1.90	0.69
1:A:303:ARG:NH1	1:A:303:ARG:HB3	2.06	0.69
1:A:267:GLN:HE21	1:A:274:LEU:H	1.38	0.68
1:B:242:ILE:HD11	1:B:280:THR:HG22	1.75	0.68
1:A:253:LEU:HD13	1:A:304:LEU:CD2	2.25	0.67
1:A:348:ASN:ND2	1:A:362:ASN:H	1.91	0.67
1:A:97:HIS:HE1	1:A:180:ASP:OD2	1.78	0.67
1:A:83:LYS:HG2	3:A:456:HOH:O	1.95	0.66
1:A:348:ASN:HD21	1:A:362:ASN:N	1.92	0.66
1:A:10:ASN:OD1	1:A:12:LEU:HB2	1.96	0.66
1:A:263:LEU:O	1:A:267:GLN:HG3	1.95	0.65
1:B:317:ILE:HG12	1:B:338:LEU:HB2	1.79	0.64
1:B:240:PRO:HB2	3:B:534:HOH:O	1.97	0.64
1:A:82:THR:OG1	1:A:85:GLN:HG3	1.99	0.63
1:B:53:ARG:HG3	1:B:53:ARG:HH11	1.63	0.63
1:A:175:ILE:HD11	1:A:229:ILE:HD12	1.80	0.63
1:B:261:GLU:HG2	1:B:265:LYS:HE3	1.80	0.63
1:A:13:PHE:CZ	1:A:356:LYS:HB2	2.34	0.63
1:B:151:LYS:HE3	3:B:602:HOH:O	1.98	0.62
1:B:263:LEU:O	1:B:267:GLN:HG3	1.98	0.62
1:A:52:GLN:HE21	1:A:53:ARG:HH12	1.47	0.62
1:B:295:LEU:HG	1:B:297:SER:H	1.64	0.62
1:B:297:SER:O	1:B:301:GLU:HG2	1.99	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:218:ILE:O	1:B:222:VAL:HG23	1.98	0.62
1:A:130:GLU:HG2	1:A:151:LYS:NZ	2.14	0.62
1:A:303:ARG:HH11	1:A:303:ARG:HB3	1.63	0.61
1:A:248:MET:HB2	3:A:593:HOH:O	2.01	0.60
1:A:200:ILE:HD11	1:A:247:ALA:HA	1.83	0.60
1:A:48:GLU:HG2	1:A:379:THR:HG21	1.82	0.60
1:A:39:ASN:ND2	1:B:303:ARG:HB3	2.17	0.60
1:A:237:ARG:HA	1:A:278:HIS:O	2.01	0.60
1:B:66:MET:HB2	1:B:79:GLY:HA2	1.84	0.60
1:B:13:PHE:CZ	1:B:356:LYS:HB2	2.36	0.60
1:A:324:ARG:HG3	1:A:355:ILE:HG23	1.83	0.59
1:A:317:ILE:HG12	1:A:338:LEU:HB2	1.83	0.59
1:B:52:GLN:HE21	1:B:53:ARG:HH12	1.50	0.58
1:B:217:PHE:O	1:B:221:VAL:HG23	2.04	0.58
1:B:296:GLY:HA3	3:B:617:HOH:O	2.04	0.56
1:A:304:LEU:HD12	3:A:524:HOH:O	2.05	0.56
1:A:222:VAL:O	1:A:226:VAL:HG23	2.05	0.56
1:B:20:LYS:HG2	3:B:629:HOH:O	2.06	0.56
1:B:253:LEU:O	1:B:257:LEU:HD13	2.05	0.56
1:B:32:MET:HA	2:B:501:FMN:N5	2.21	0.56
1:B:237:ARG:HA	1:B:278:HIS:O	2.05	0.56
1:A:248:MET:HE3	3:A:521:HOH:O	2.07	0.55
1:B:97:HIS:HE1	1:B:180:ASP:OD2	1.90	0.54
1:B:350:ASP:O	1:B:354:ARG:HG3	2.07	0.54
1:A:239:SER:HB2	1:A:242:ILE:CG1	2.36	0.54
1:A:299:GLU:O	1:A:303:ARG:HG2	2.07	0.54
1:B:242:ILE:HG23	3:B:626:HOH:O	2.06	0.54
1:B:83:LYS:HG2	3:B:549:HOH:O	2.07	0.54
1:B:80:ILE:O	1:B:80:ILE:HG23	2.08	0.53
1:A:183:GLU:OE2	1:A:278:HIS:HD2	1.90	0.53
1:A:253:LEU:HD13	1:A:304:LEU:HD22	1.89	0.53
1:B:132:PRO:HB3	1:B:151:LYS:HA	1.90	0.53
1:B:304:LEU:O	1:B:308:LEU:HG	2.07	0.53
1:B:184:ILE:HD12	1:B:192:ILE:CD1	2.38	0.53
1:B:348:ASN:ND2	1:B:362:ASN:H	2.04	0.53
1:A:83:LYS:HG3	1:A:84:GLU:N	2.23	0.53
1:A:275:ALA:O	1:A:276:TYR:HB3	2.08	0.53
1:B:183:GLU:OE2	1:B:278:HIS:HD2	1.91	0.53
1:B:139:ILE:HG23	1:B:141:MET:HE2	1.91	0.53
1:B:282:PRO:O	1:B:283:ARG:HB2	2.08	0.53
1:A:11:PRO:HD2	1:A:328:ILE:HG23	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:130:GLU:HG2	1:A:151:LYS:HZ3	1.75	0.52
1:A:32:MET:HA	2:A:401:FMN:N5	2.25	0.52
1:A:241:ALA:O	1:A:282:PRO:HG3	2.10	0.52
1:A:325:GLU:CD	1:A:325:GLU:H	2.12	0.52
1:A:239:SER:HB2	1:A:242:ILE:HG13	1.94	0.50
1:B:275:ALA:O	1:B:276:TYR:HB3	2.11	0.50
1:B:112:ARG:HB3	1:B:127:SER:HB2	1.92	0.50
1:A:304:LEU:C	1:A:304:LEU:HD13	2.32	0.50
1:B:349:PRO:HG3	3:B:405:HOH:O	2.11	0.50
1:A:52:GLN:NE2	1:A:53:ARG:HH12	2.09	0.50
1:A:175:ILE:HG13	1:A:229:ILE:CG2	2.42	0.49
1:B:348:ASN:HD21	1:B:362:ASN:N	2.06	0.49
1:B:368:THR:HG21	3:B:401:HOH:O	2.11	0.49
1:B:347:SER:HA	1:B:381:TYR:CG	2.47	0.49
1:A:54:ALA:HB2	1:A:60:LEU:HD11	1.95	0.49
1:B:134:SER:OG	1:B:136:ARG:HG2	2.11	0.49
1:A:297:SER:O	1:A:301:GLU:HG2	2.13	0.49
1:A:80:ILE:HD11	1:A:174:ALA:HA	1.95	0.49
1:B:53:ARG:NH1	1:B:53:ARG:HG3	2.28	0.49
1:A:20:LYS:HD2	3:A:570:HOH:O	2.12	0.48
1:A:282:PRO:O	1:A:283:ARG:HB2	2.14	0.48
1:A:86:VAL:HG13	1:A:177:ALA:HB2	1.96	0.48
1:B:242:ILE:CD1	1:B:280:THR:HG22	2.41	0.48
1:B:304:LEU:C	1:B:304:LEU:HD13	2.33	0.48
1:B:262:ARG:O	1:B:266:ILE:HG13	2.14	0.48
1:A:128:SER:HB3	1:A:195:PHE:CE1	2.48	0.48
1:A:347:SER:HA	1:A:381:TYR:CG	2.48	0.48
1:A:33:THR:O	1:A:34:ARG:HD2	2.13	0.47
1:A:8:GLY:C	1:A:10:ASN:N	2.60	0.47
1:A:165:GLU:O	1:A:169:ARG:HG3	2.14	0.47
1:B:253:LEU:CD1	1:B:304:LEU:HD22	2.27	0.47
1:B:243:ASP:HB3	3:B:498:HOH:O	2.14	0.47
1:B:345:PHE:CD2	1:B:351:LEU:HD23	2.50	0.47
1:B:169:ARG:HD3	3:B:608:HOH:O	2.15	0.47
1:A:112:ARG:HB3	1:A:127:SER:HB2	1.97	0.46
1:B:130:GLU:O	1:B:152:PRO:HD2	2.15	0.46
1:A:97:HIS:HD2	3:A:583:HOH:O	1.98	0.46
1:A:237:ARG:HH11	1:A:278:HIS:CE1	2.34	0.46
1:A:83:LYS:HD2	3:A:565:HOH:O	2.16	0.46
1:B:319:SER:HB2	1:B:340:SER:OG	2.16	0.46
1:A:233:ARG:HD3	3:A:508:HOH:O	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:253:LEU:HD13	1:A:304:LEU:HD23	1.96	0.46
1:A:350:ASP:O	1:A:354:ARG:HG3	2.16	0.46
1:B:136:ARG:HD3	3:B:595:HOH:O	2.15	0.46
1:B:267:GLN:NE2	1:B:274:LEU:N	2.52	0.46
1:B:66:MET:HB2	1:B:79:GLY:CA	2.46	0.45
1:A:16:TYR:CE1	1:A:337:ASP:HB3	2.52	0.45
1:B:241:ALA:O	1:B:282:PRO:HG3	2.16	0.45
1:A:318:CYS:SG	1:A:339:VAL:HG22	2.56	0.45
1:A:253:LEU:HD12	1:A:253:LEU:O	2.17	0.45
1:B:120:PRO:HB3	3:B:503:HOH:O	2.17	0.45
1:A:87:ARG:NH1	3:A:565:HOH:O	2.49	0.45
1:A:256:GLY:O	1:A:260:VAL:HG23	2.16	0.44
1:A:321:GLY:HA2	2:A:401:FMN:O3P	2.16	0.44
1:B:376:VAL:HA	1:B:380:ASP:OD2	2.17	0.44
1:A:66:MET:HB2	1:A:79:GLY:HA2	1.99	0.44
1:B:267:GLN:HE21	1:B:274:LEU:N	2.01	0.44
1:A:134:SER:C	1:A:136:ARG:H	2.21	0.44
1:A:93:VAL:HG12	1:A:97:HIS:CD2	2.53	0.44
1:B:256:GLY:O	1:B:260:VAL:HG23	2.17	0.44
1:B:52:GLN:NE2	1:B:53:ARG:HH12	2.15	0.44
1:A:323:THR:H	1:A:326:LEU:HB2	1.82	0.44
1:B:329:GLU:HB2	3:B:470:HOH:O	2.16	0.44
1:B:130:GLU:O	1:B:151:LYS:HG3	2.18	0.44
1:A:370:TYR:CZ	2:A:401:FMN:HM72	2.52	0.44
1:A:278:HIS:HE1	1:A:319:SER:HB2	1.83	0.44
1:B:248:MET:HB2	3:B:541:HOH:O	2.17	0.44
1:B:129:THR:OG1	1:B:130:GLU:N	2.51	0.43
1:B:343:ARG:HB3	1:B:364:TYR:CE2	2.52	0.43
1:A:130:GLU:HG2	1:A:151:LYS:HZ1	1.82	0.43
1:A:182:ILE:C	1:A:182:ILE:HD12	2.39	0.43
1:B:182:ILE:HD12	1:B:182:ILE:C	2.39	0.43
1:A:99:LYS:HB3	1:A:99:LYS:NZ	2.34	0.43
1:A:276:TYR:CE1	1:A:278:HIS:HB2	2.53	0.43
1:B:325:GLU:OE1	1:B:325:GLU:HA	2.18	0.43
1:A:114:SER:O	1:A:150:PRO:HD2	2.19	0.43
1:A:203:ARG:HB2	1:A:208:GLY:HA3	2.00	0.43
1:A:257:LEU:HD12	1:A:257:LEU:HA	1.84	0.43
1:B:276:TYR:CD1	1:B:276:TYR:C	2.92	0.42
1:A:104:PHE:HA	1:A:181:GLY:O	2.19	0.42
1:A:216:LYS:HE3	1:A:216:LYS:HB2	1.84	0.42
1:A:237:ARG:HD3	1:A:278:HIS:CD2	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:354:ARG:O	1:A:358:ASN:N	2.52	0.42
1:A:140:LEU:HB3	3:A:407:HOH:O	2.19	0.42
1:A:80:ILE:HG23	1:A:80:ILE:O	2.19	0.42
1:A:8:GLY:O	1:A:10:ASN:N	2.51	0.42
1:B:28:VAL:O	1:B:340:SER:HA	2.20	0.42
1:B:48:GLU:HG2	3:B:654:HOH:O	2.19	0.42
1:B:15:PRO:HG3	3:B:499:HOH:O	2.18	0.42
1:B:165:GLU:HA	1:B:165:GLU:OE1	2.18	0.42
1:B:36:ARG:HB2	3:B:525:HOH:O	2.19	0.42
1:A:52:GLN:HE21	1:A:53:ARG:NH1	2.16	0.41
1:B:237:ARG:HG3	1:B:278:HIS:HB3	2.03	0.41
1:B:80:ILE:HD11	1:B:174:ALA:HA	2.01	0.41
1:A:198:ASP:OD1	1:A:249:ASP:HA	2.19	0.41
1:B:12:LEU:HD23	1:B:331:VAL:HG21	2.02	0.41
1:A:303:ARG:HH11	1:A:303:ARG:CB	2.30	0.41
1:B:139:ILE:HG23	1:B:141:MET:CE	2.49	0.41
1:A:99:LYS:HE3	3:A:616:HOH:O	2.20	0.41
1:A:216:LYS:O	1:A:220:GLN:HG3	2.21	0.41
1:B:171:ALA:O	1:B:175:ILE:HG12	2.21	0.41
1:A:128:SER:HB3	1:A:195:PHE:CD1	2.56	0.40
1:A:129:THR:OG1	1:A:130:GLU:N	2.54	0.40
1:A:132:PRO:HB3	1:A:151:LYS:HA	2.02	0.40
1:A:347:SER:HA	1:A:381:TYR:CD1	2.57	0.40
1:B:184:ILE:HD12	1:B:192:ILE:HD13	2.03	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	363/402 (90%)	343 (94%)	20 (6%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	359/402 (89%)	345 (96%)	13 (4%)	1 (0%)	44	55
All	All	722/804 (90%)	688 (95%)	33 (5%)	1 (0%)	55	67

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	80	ILE

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	295/328 (90%)	288 (98%)	7 (2%)	54	72
1	B	294/328 (90%)	288 (98%)	6 (2%)	60	77
All	All	589/656 (90%)	576 (98%)	13 (2%)	57	74

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	48	GLU
1	A	237	ARG
1	A	257	LEU
1	A	278	HIS
1	A	309	ARG
1	A	325	GLU
1	A	345	PHE
1	B	278	HIS
1	B	306	ARG
1	B	309	ARG
1	B	325	GLU
1	B	345	PHE
1	B	358	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	39	ASN
1	A	43	GLN
1	A	52	GLN
1	A	97	HIS
1	A	135	ASN
1	A	162	GLN
1	A	267	GLN
1	A	278	HIS
1	A	281	GLN
1	A	348	ASN
1	A	358	ASN
1	A	365	ASN
1	B	39	ASN
1	B	43	GLN
1	B	52	GLN
1	B	97	HIS
1	B	135	ASN
1	B	162	GLN
1	B	267	GLN
1	B	278	HIS
1	B	281	GLN
1	B	348	ASN
1	B	358	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 ⓘ

2 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	$\# Z > 2$
2	FMN	A	401	-	31,33,33	2.33	10 (32%)	38,50,50	2.90	12 (31%)
2	FMN	B	501	-	31,33,33	2.31	10 (32%)	38,50,50	2.92	11 (28%)

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
2	FMN	A	401	-	-	0/16/18/18	0/3/3/3
2	FMN	B	501	-	-	0/16/18/18	0/3/3/3

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	FMN	O4'-C4'	2.07	1.47	1.43
2	A	401	FMN	C9-C9A	2.30	1.45	1.40
2	A	401	FMN	C4-N3	2.41	1.37	1.33
2	B	501	FMN	C4-N3	2.61	1.37	1.33
2	A	401	FMN	C5A-N5	2.84	1.39	1.35
2	B	501	FMN	C9A-C5A	2.88	1.48	1.42
2	A	401	FMN	C9A-C5A	2.93	1.48	1.42
2	B	501	FMN	C5A-N5	3.16	1.40	1.35
2	B	501	FMN	C2-N3	3.41	1.44	1.38
2	A	401	FMN	C2-N3	3.50	1.45	1.38
2	B	501	FMN	C4-C4A	3.84	1.48	1.41
2	A	401	FMN	C4-C4A	4.12	1.49	1.41
2	A	401	FMN	C8-C7	4.16	1.51	1.41
2	A	401	FMN	C6-C5A	4.28	1.48	1.41
2	B	501	FMN	C6-C5A	4.57	1.48	1.41
2	B	501	FMN	C8-C7	4.58	1.52	1.41
2	B	501	FMN	C9A-N10	4.92	1.45	1.38
2	B	501	FMN	C4A-C10	5.20	1.50	1.41
2	A	401	FMN	C4A-C10	5.32	1.50	1.41
2	A	401	FMN	C9A-N10	5.43	1.45	1.38

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	FMN	C4A-C10-N10	-5.99	116.36	120.52
2	A	401	FMN	C4A-C10-N10	-5.94	116.39	120.52
2	B	501	FMN	C4A-C4-N3	-5.36	115.86	123.48
2	A	401	FMN	C4A-C4-N3	-5.30	115.93	123.48
2	B	501	FMN	C4-C4A-C10	-4.79	116.09	119.96
2	A	401	FMN	C4-C4A-C10	-4.76	116.11	119.96
2	A	401	FMN	O2'-C2'-C3'	-3.37	100.74	109.09
2	A	401	FMN	C5A-C9A-N10	-3.26	115.24	117.66
2	B	501	FMN	O2'-C2'-C3'	-3.25	101.03	109.09
2	B	501	FMN	C5A-C9A-N10	-3.12	115.34	117.66
2	A	401	FMN	O2'-C2'-C1'	2.03	114.48	109.79
2	A	401	FMN	C1'-N10-C9A	2.04	120.22	118.35
2	B	501	FMN	C1'-N10-C9A	2.12	120.29	118.35
2	B	501	FMN	O5'-C5'-C4'	2.20	115.22	109.36
2	A	401	FMN	O5'-C5'-C4'	2.28	115.45	109.36
2	B	501	FMN	C4-C4A-N5	2.30	121.20	118.68
2	A	401	FMN	C4-C4A-N5	2.32	121.22	118.68
2	B	501	FMN	C4A-N5-C5A	3.05	119.98	116.76
2	A	401	FMN	C4A-N5-C5A	3.15	120.09	116.76
2	B	501	FMN	C1'-C2'-C3'	3.47	119.73	109.82
2	A	401	FMN	C1'-C2'-C3'	3.49	119.80	109.82
2	A	401	FMN	C4-N3-C2	12.15	125.78	115.16
2	B	501	FMN	C4-N3-C2	12.24	125.86	115.16

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	FMN	3	0
2	B	501	FMN	1	0

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	367/402 (91%)	0.08	14 (3%) 41 48	20, 34, 51, 97	0
1	B	363/402 (90%)	-0.04	15 (4%) 38 45	18, 29, 48, 93	0
All	All	730/804 (90%)	0.02	29 (3%) 39 46	18, 32, 50, 97	0

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	7	ASP	12.6
1	A	8	GLY	8.8
1	A	295	LEU	8.2
1	B	295	LEU	8.2
1	A	296	GLY	5.2
1	B	296	GLY	5.1
1	B	297	SER	4.9
1	A	9	ASN	4.5
1	A	297	SER	4.4
1	B	298	GLU	4.2
1	B	299	GLU	3.2
1	B	282	PRO	3.1
1	A	298	GLU	2.9
1	A	121	ALA	2.9
1	B	204	THR	2.8
1	B	136	ARG	2.7
1	B	248	MET	2.5
1	B	283	ARG	2.5
1	A	136	ARG	2.5
1	A	268	LEU	2.4
1	B	121	ALA	2.4
1	B	301	GLU	2.3
1	A	202	ASP	2.3
1	B	83	LYS	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	87	ARG	2.2
1	B	268	LEU	2.1
1	A	366	ARG	2.1
1	A	122	GLY	2.1
1	A	248	MET	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(Å ²)	Q<0.9
2	FMN	A	401	31/31	0.96	0.12	0.52	30,35,37,38	0
2	FMN	B	501	31/31	0.97	0.10	-0.10	22,25,28,31	0

6.5 Other polymers [i](#)

There are no such residues in this entry.