



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 09:43 PM UTC

PDB ID : 3KKJ / pdb_00003kkj
Title : X-ray structure of P. syringae q888a4 Oxidoreductase at resolution 2.5Å, Northeast Structural Genomics Consortium target PsR10
Authors : Kuzin, A.P.; Chen, Y.; Forouhar, F.; Vorobiev, S.; Acton, T.; Ma, L.C.; Xiao, R.; Montelione, G.T.; Hunt, J.F.; Tong, T.; Northeast Structural Genomics Consortium (NESG)
Deposited on : 2009-11-05
Resolution : 2.50 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtrriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

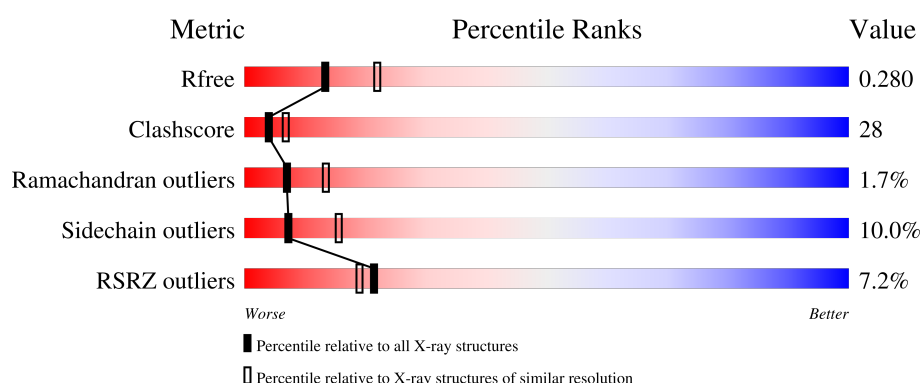
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	336	4% 56% 34% 7% .
1	B	336	9% 52% 38% 7% ..

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5328 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.

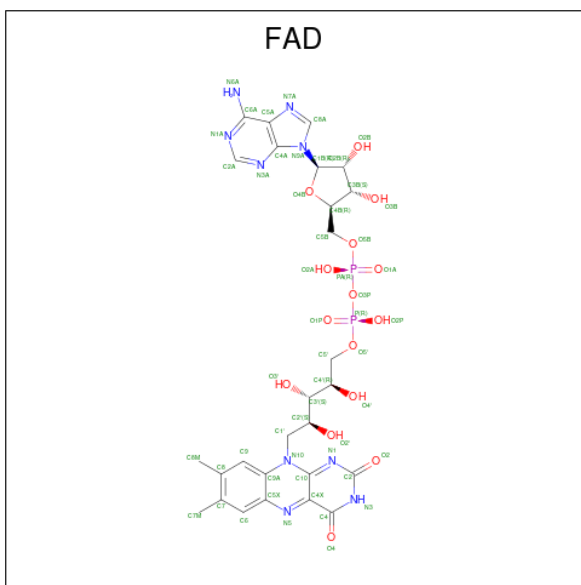
- Molecule 1 is a protein called Amine oxidase, flavin-containing.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	328	Total	C	N	O	S	Se	0	0	0
			2524	1590	461	460	5	8			
1	B	328	Total	C	N	O	S	Se	0	0	0
			2524	1590	461	460	5	8			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	-	initiating methionine	UNP Q888A4
A	329	LEU	-	expression tag	UNP Q888A4
A	330	GLU	-	expression tag	UNP Q888A4
A	331	HIS	-	expression tag	UNP Q888A4
A	332	HIS	-	expression tag	UNP Q888A4
A	333	HIS	-	expression tag	UNP Q888A4
A	334	HIS	-	expression tag	UNP Q888A4
A	335	HIS	-	expression tag	UNP Q888A4
A	336	HIS	-	expression tag	UNP Q888A4
B	1	MSE	-	initiating methionine	UNP Q888A4
B	329	LEU	-	expression tag	UNP Q888A4
B	330	GLU	-	expression tag	UNP Q888A4
B	331	HIS	-	expression tag	UNP Q888A4
B	332	HIS	-	expression tag	UNP Q888A4
B	333	HIS	-	expression tag	UNP Q888A4
B	334	HIS	-	expression tag	UNP Q888A4
B	335	HIS	-	expression tag	UNP Q888A4
B	336	HIS	-	expression tag	UNP Q888A4

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



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

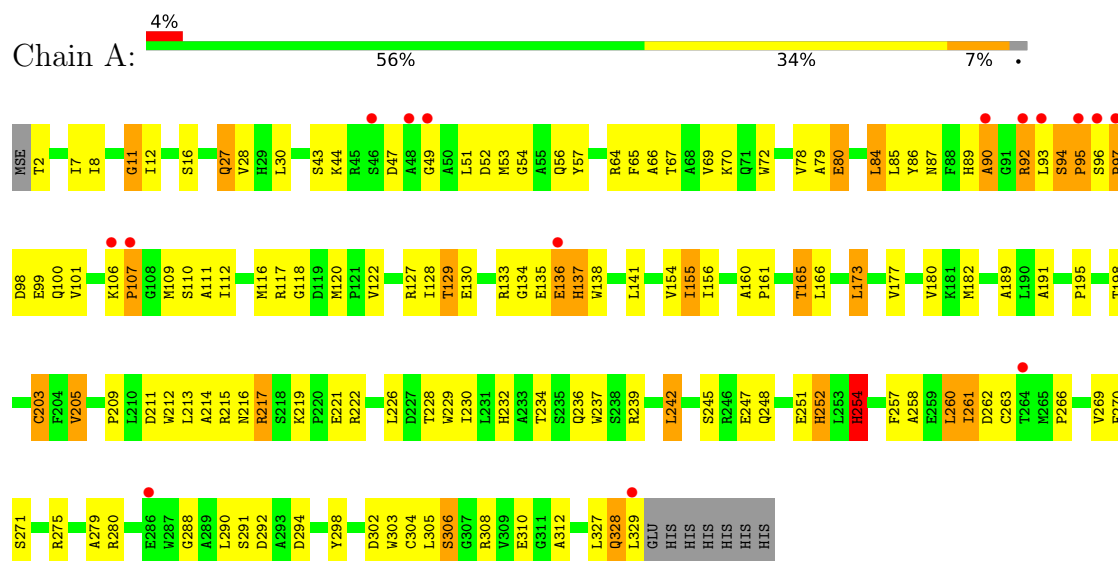
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	86	Total O 86 86	0	0
3	B	88	Total O 88 88	0	0

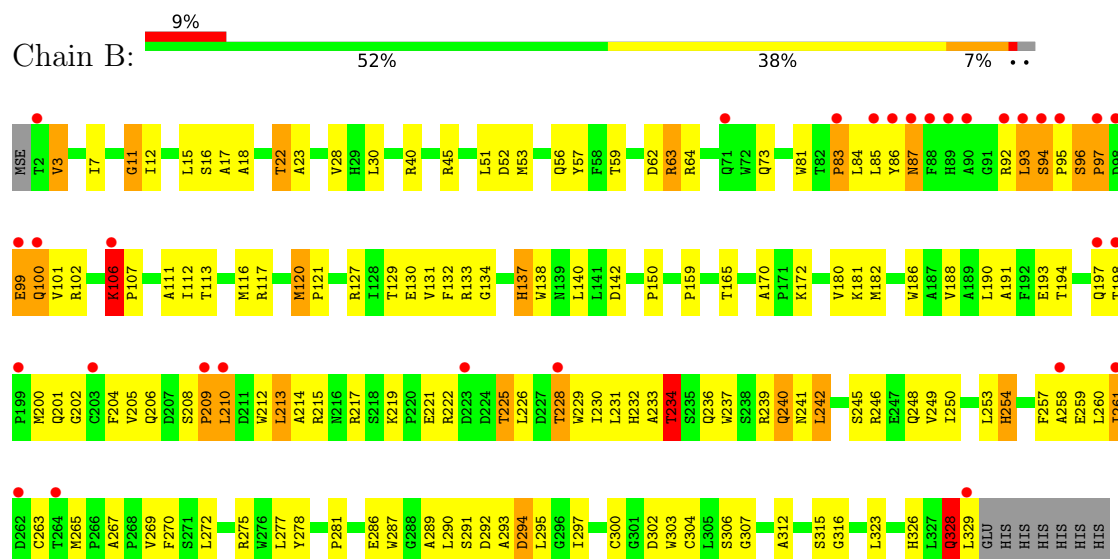
3 Residue-property plots

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

- Molecule 1: Amine oxidase, flavin-containing



- Molecule 1: Amine oxidase, flavin-containing



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	120.70Å 120.70Å 114.79Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 2.50 20.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	96.8 (20.00-2.50) 97.0 (20.00-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	7.08 (at 2.51Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.4_115)	Depositor
R, R_{free}	0.227 , 0.281 0.227 , 0.280	Depositor DCC
R_{free} test set	3189 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	38.8	Xtriage
Anisotropy	0.194	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 46.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.018 for -h,-k,l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	5328	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.20% 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: 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	0.64	0/2585	1.07	19/3507 (0.5%)
1	B	0.67	0/2585	1.15	30/3507 (0.9%)
All	All	0.66	0/5170	1.11	49/7014 (0.7%)

There are no bond length outliers.

All (49) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	205	VAL	N-CA-C	10.38	123.92	108.46
1	B	328	GLN	N-CA-C	-9.61	99.39	112.94
1	A	11	GLY	N-CA-C	-8.71	99.35	112.41
1	A	291	SER	N-CA-C	7.74	119.91	108.14
1	B	106	LYS	CA-C-N	-7.52	110.65	118.85
1	B	106	LYS	C-N-CA	-7.52	110.65	118.85
1	B	137	HIS	N-CA-C	7.19	121.02	110.59
1	A	205	VAL	N-CA-C	6.92	117.79	108.11
1	B	291	SER	N-CA-C	6.46	118.37	108.42
1	B	96	SER	N-CA-C	6.44	118.60	108.55
1	B	194	THR	CA-C-N	6.35	126.72	119.93
1	B	194	THR	C-N-CA	6.35	126.72	119.93
1	B	94	SER	CA-C-N	6.33	128.01	120.23
1	B	94	SER	C-N-CA	6.33	128.01	120.23
1	A	252	HIS	N-CA-C	6.28	117.92	111.14
1	A	209	PRO	N-CA-C	-6.26	105.19	113.65
1	A	137	HIS	N-CA-C	6.11	118.49	109.69
1	A	135	GLU	N-CA-C	-6.02	104.64	111.14
1	B	293	ALA	N-CA-C	6.00	119.43	111.75
1	B	306	SER	N-CA-C	5.97	117.87	111.36
1	A	294	ASP	N-CA-C	-5.87	104.80	111.07
1	B	307	GLY	N-CA-C	-5.78	107.00	115.72

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	269	VAL	N-CA-C	-5.74	104.77	110.62
1	B	267	ALA	N-CA-C	-5.72	102.19	110.08
1	B	11	GLY	N-CA-C	-5.54	103.69	112.66
1	A	292	ASP	N-CA-C	-5.54	99.38	109.56
1	A	304	CYS	N-CA-C	-5.44	106.60	113.18
1	A	216	ASN	N-CA-C	5.42	117.88	111.33
1	A	99	GLU	N-CA-C	-5.35	106.59	113.01
1	A	203	CYS	N-CA-C	5.33	116.73	107.93
1	B	112	ILE	N-CA-C	5.33	115.53	110.53
1	B	106	LYS	N-CA-C	5.29	121.51	109.81
1	A	254	HIS	N-CA-C	-5.29	105.51	111.28
1	B	304	CYS	N-CA-C	-5.28	106.84	113.28
1	A	134	GLY	N-CA-C	-5.22	105.14	112.60
1	B	83	PRO	N-CA-C	5.22	118.69	110.50
1	B	170	ALA	CA-C-N	5.19	126.33	119.84
1	B	170	ALA	C-N-CA	5.19	126.33	119.84
1	B	219	LYS	CA-C-N	5.19	125.19	119.90
1	B	219	LYS	C-N-CA	5.19	125.19	119.90
1	A	263	CYS	N-CA-C	5.17	117.48	110.35
1	A	97	PRO	N-CA-C	-5.16	101.83	112.47
1	A	306	SER	N-CA-C	5.15	119.27	112.89
1	B	254	HIS	N-CA-C	-5.12	105.39	111.69
1	B	209	PRO	N-CA-C	-5.08	107.51	113.86
1	B	63	ARG	N-CA-C	5.07	117.70	111.82
1	B	198	THR	N-CA-C	-5.02	101.64	108.11
1	B	17	ALA	N-CA-C	-5.01	105.71	111.07
1	B	134	GLY	N-CA-C	-5.00	105.35	112.55

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	2524	0	2448	138	0
1	B	2524	0	2448	150	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	53	0	31	7	0
2	B	53	0	31	6	0
3	A	86	0	0	5	0
3	B	88	0	0	4	0
All	All	5328	0	4958	282	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 28.

All (282) 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:A:129:THR:HG22	1:A:130:GLU:HG2	1.37	1.05
1:B:215:ARG:HH21	1:B:217:ARG:NH2	1.61	0.98
1:B:57:TYR:CD2	1:B:102:ARG:HD3	2.06	0.91
1:B:87:ASN:HB3	1:B:95:PRO:HG2	1.50	0.91
1:B:234:THR:HG22	1:B:237:TRP:H	1.34	0.91
1:A:85:LEU:O	1:A:95:PRO:HD2	1.74	0.88
1:A:92:ARG:H	1:A:92:ARG:HE	1.22	0.87
1:B:97:PRO:C	1:B:99:GLU:H	1.79	0.87
1:B:215:ARG:NH2	1:B:217:ARG:HH21	1.73	0.86
1:B:215:ARG:HH21	1:B:217:ARG:HH21	0.89	0.86
1:B:234:THR:CG2	1:B:237:TRP:H	1.89	0.85
1:B:85:LEU:HD13	1:B:204:PHE:CE2	2.12	0.84
1:B:215:ARG:HH11	1:B:229:TRP:HZ2	1.26	0.84
1:B:106:LYS:HB3	1:B:107:PRO:HD3	1.60	0.82
1:B:93:LEU:C	1:B:95:PRO:HD3	2.06	0.81
1:B:94:SER:N	1:B:95:PRO:HD3	1.96	0.80
1:B:258:ALA:HA	1:B:261:ILE:HD11	1.62	0.80
1:A:327:LEU:C	1:A:329:LEU:H	1.85	0.80
1:A:328:GLN:HE21	1:A:328:GLN:N	1.82	0.78
1:B:254:HIS:NE2	1:B:265:MSE:HE3	1.99	0.77
1:B:245:SER:OG	1:B:248:GLN:HG3	1.85	0.77
1:B:85:LEU:HD13	1:B:204:PHE:HE2	1.50	0.76
1:A:92:ARG:HD3	1:A:260:LEU:O	1.84	0.75
1:B:87:ASN:N	1:B:95:PRO:HD2	2.01	0.74
1:A:7:ILE:HB	1:A:30:LEU:HD23	1.69	0.74
1:B:100:GLN:HE21	1:B:100:GLN:CA	2.01	0.73
1:A:96:SER:C	1:A:98:ASP:H	1.97	0.73
1:A:52:ASP:O	1:A:230:ILE:HD13	1.87	0.73
1:A:236:GLN:H	1:A:236:GLN:NE2	1.87	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:258:ALA:HB2	1:B:265:MSE:CE	2.19	0.72
1:A:261:ILE:HD13	1:A:261:ILE:H	1.55	0.72
1:B:100:GLN:HE21	1:B:100:GLN:HA	1.54	0.72
1:A:236:GLN:H	1:A:236:GLN:CD	1.98	0.71
1:A:86:TYR:HA	1:A:95:PRO:HD3	1.73	0.71
1:A:52:ASP:OD2	1:A:109:MSE:HB2	1.91	0.70
1:A:288:GLY:HA2	1:A:305:LEU:HD21	1.74	0.70
1:B:97:PRO:C	1:B:99:GLU:N	2.42	0.70
1:A:129:THR:HG22	1:A:130:GLU:CG	2.19	0.69
1:B:107:PRO:HD2	1:B:111:ALA:HB2	1.75	0.69
1:A:112:ILE:O	1:A:116:MSE:HG3	1.93	0.68
1:B:215:ARG:HD2	1:B:229:TRP:CZ2	2.29	0.67
1:B:87:ASN:HB3	1:B:95:PRO:CG	2.22	0.67
1:A:239:ARG:HG2	1:A:239:ARG:HH11	1.61	0.66
1:B:84:LEU:HD23	1:B:201:GLN:OE1	1.96	0.66
1:A:106:LYS:HB3	1:A:107:PRO:HD3	1.78	0.65
1:A:155:ILE:HD12	1:A:298:TYR:HB2	1.78	0.65
1:B:237:TRP:NE1	1:B:241:ASN:ND2	2.43	0.65
1:A:160:ALA:HB3	1:A:182:MSE:HE3	1.78	0.65
1:A:257:PHE:O	1:A:261:ILE:HD13	1.96	0.65
1:B:258:ALA:HB2	1:B:265:MSE:HE2	1.78	0.65
1:A:258:ALA:HA	1:A:261:ILE:HD11	1.78	0.64
1:B:237:TRP:CE2	1:B:241:ASN:ND2	2.65	0.64
1:A:248:GLN:O	1:A:252:HIS:HD2	1.81	0.63
1:A:234:THR:HG23	1:A:237:TRP:H	1.64	0.62
1:A:165:THR:OG1	1:B:165:THR:HG21	1.99	0.62
1:B:40:ARG:NE	2:B:402:FAD:O2A	2.29	0.62
1:B:292:ASP:OD2	1:B:295:LEU:HB2	2.00	0.61
1:A:79:ALA:HB2	1:A:106:LYS:CD	2.31	0.61
1:B:215:ARG:NH2	1:B:217:ARG:NH2	2.40	0.61
1:B:133:ARG:HD2	1:B:138:TRP:CZ2	2.36	0.61
1:A:302:ASP:HB3	1:A:312:ALA:HB2	1.82	0.61
1:B:56:GLN:HE21	1:B:232:HIS:CE1	2.19	0.61
1:A:92:ARG:H	1:A:92:ARG:NE	1.96	0.61
1:B:294:ASP:OD2	1:B:294:ASP:N	2.24	0.61
1:A:95:PRO:HG2	1:A:96:SER:H	1.66	0.61
1:A:161:PRO:O	1:B:165:THR:HG22	2.00	0.61
1:A:212:TRP:CZ2	1:A:214:ALA:HB2	2.37	0.60
1:A:154:VAL:C	1:A:155:ILE:HD13	2.26	0.60
1:A:130:GLU:HB2	1:A:141:LEU:HB2	1.83	0.59
1:A:234:THR:CG2	1:A:237:TRP:H	2.15	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:56:GLN:HG3	1:B:212:TRP:NE1	2.18	0.59
1:B:302:ASP:HB3	2:B:402:FAD:H5'2	1.85	0.59
1:A:260:LEU:C	1:A:260:LEU:HD12	2.28	0.59
1:B:215:ARG:CD	1:B:229:TRP:CZ2	2.85	0.59
1:A:87:ASN:O	1:A:92:ARG:HA	2.03	0.58
1:B:180:VAL:HG12	1:B:182:MSE:HE3	1.86	0.58
1:B:236:GLN:CD	1:B:236:GLN:H	2.10	0.58
1:A:67:THR:HA	1:A:70:LYS:HD2	1.86	0.57
1:A:203:CYS:HB3	1:A:213:LEU:HG	1.86	0.57
1:A:86:TYR:CA	1:A:95:PRO:HD3	2.34	0.57
1:A:65:PHE:O	1:A:69:VAL:HG23	2.04	0.57
1:A:133:ARG:HD2	1:A:138:TRP:CZ2	2.40	0.57
1:B:117:ARG:O	1:B:120:MSE:HB2	2.04	0.57
1:B:261:ILE:HB	1:B:263:CYS:SG	2.45	0.57
1:B:200:MSE:C	1:B:201:GLN:HG3	2.29	0.56
1:A:57:TYR:C	1:A:57:TYR:CD1	2.84	0.56
1:B:326:HIS:C	1:B:328:GLN:H	2.14	0.56
1:B:209:PRO:HB3	1:B:237:TRP:CG	2.40	0.56
1:B:182:MSE:HE1	1:B:281:PRO:HB3	1.86	0.56
1:A:80:GLU:OE2	1:A:101:VAL:HG13	2.05	0.56
1:A:261:ILE:HD13	1:A:261:ILE:N	2.20	0.56
1:B:221:GLU:O	1:B:221:GLU:HG2	2.05	0.55
1:A:180:VAL:HG12	1:A:182:MSE:HE2	1.86	0.55
1:A:96:SER:C	1:A:98:ASP:N	2.63	0.55
1:B:56:GLN:NE2	1:B:232:HIS:CE1	2.74	0.55
1:A:92:ARG:HE	1:A:92:ARG:N	2.00	0.55
1:B:300:CYS:HB2	1:B:315:SER:OG	2.07	0.55
1:B:106:LYS:HB3	1:B:107:PRO:CD	2.35	0.55
1:A:127:ARG:HD3	3:A:416:HOH:O	2.06	0.55
1:A:327:LEU:C	1:A:329:LEU:N	2.56	0.55
1:B:100:GLN:HE21	1:B:101:VAL:N	2.05	0.54
1:A:7:ILE:CD1	1:A:28:VAL:HB	2.37	0.54
1:A:47:ASP:HB2	1:A:270:PHE:CE2	2.42	0.54
1:A:130:GLU:OE2	1:B:278:TYR:HE2	1.91	0.54
1:B:237:TRP:CZ3	1:B:253:LEU:HD21	2.43	0.54
1:B:7:ILE:HB	1:B:30:LEU:HD23	1.90	0.53
1:A:44:LYS:HD3	1:A:53:MSE:HE1	1.91	0.53
1:B:209:PRO:HB3	1:B:237:TRP:CD1	2.43	0.53
1:B:257:PHE:O	1:B:261:ILE:HG13	2.09	0.53
1:A:89:HIS:O	1:A:90:ALA:CB	2.56	0.53
1:B:85:LEU:HD12	1:B:97:PRO:CD	2.38	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:215:ARG:HD2	1:B:229:TRP:CE2	2.43	0.53
1:A:155:ILE:HD13	1:A:155:ILE:N	2.22	0.53
1:A:89:HIS:O	1:A:90:ALA:HB3	2.09	0.53
1:B:56:GLN:HE21	1:B:232:HIS:HE1	1.57	0.53
1:B:59:THR:HG22	1:B:101:VAL:O	2.08	0.53
1:B:100:GLN:HA	1:B:100:GLN:NE2	2.22	0.52
1:B:182:MSE:HE2	1:B:281:PRO:HA	1.90	0.52
1:A:191:ALA:HB2	1:A:228:THR:HG22	1.90	0.52
1:A:217:ARG:HG2	3:A:455:HOH:O	2.10	0.52
1:B:86:TYR:C	1:B:95:PRO:HD2	2.34	0.52
1:A:86:TYR:HA	1:A:95:PRO:CD	2.37	0.52
1:B:57:TYR:CE2	1:B:102:ARG:HD3	2.44	0.52
1:B:131:VAL:HG22	1:B:140:LEU:HD22	1.92	0.52
1:B:259:GLU:O	1:B:259:GLU:HG2	2.10	0.52
1:B:182:MSE:HE1	1:B:303:TRP:CB	2.40	0.52
1:A:219:LYS:HB2	1:A:222:ARG:HG3	1.92	0.52
1:B:3:VAL:HG13	1:B:150:PRO:HG2	1.92	0.52
1:B:57:TYR:O	2:B:402:FAD:N3	2.43	0.52
1:A:11:GLY:HA3	2:A:401:FAD:O1A	2.10	0.52
1:A:67:THR:O	1:A:70:LYS:HB2	2.09	0.51
1:A:328:GLN:HE21	1:A:328:GLN:H	1.57	0.51
1:B:100:GLN:CA	1:B:100:GLN:NE2	2.73	0.51
1:B:81:TRP:CZ2	1:B:83:PRO:HB3	2.45	0.51
1:B:242:LEU:HD23	1:B:275:ARG:HD2	1.92	0.51
1:B:323:LEU:C	1:B:323:LEU:HD23	2.35	0.51
1:B:242:LEU:CD2	1:B:275:ARG:HD2	2.41	0.51
1:A:86:TYR:C	1:A:95:PRO:HD3	2.36	0.51
1:B:202:GLY:HA2	1:B:213:LEU:O	2.10	0.51
1:A:95:PRO:HG2	1:A:96:SER:N	2.25	0.51
1:B:234:THR:HG23	1:B:236:GLN:N	2.26	0.51
1:A:107:PRO:HD2	1:A:111:ALA:CB	2.41	0.51
1:A:95:PRO:HG2	1:A:97:PRO:HD3	1.93	0.50
1:A:254:HIS:C	1:A:254:HIS:HD1	2.18	0.50
1:B:97:PRO:O	1:B:99:GLU:N	2.45	0.50
1:A:79:ALA:HB2	1:A:106:LYS:HE2	1.94	0.50
1:B:233:ALA:HB2	1:B:253:LEU:HD11	1.93	0.50
1:B:254:HIS:CE1	1:B:265:MSE:HE3	2.46	0.50
1:B:186:TRP:CE3	1:B:249:VAL:HG11	2.47	0.50
1:B:85:LEU:HD13	1:B:204:PHE:CZ	2.45	0.49
1:A:280:ARG:HD2	3:A:462:HOH:O	2.11	0.49
1:B:172:LYS:HB2	3:B:409:HOH:O	2.11	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:92:ARG:HH12	1:A:262:ASP:CG	2.20	0.49
1:A:239:ARG:HG2	1:A:239:ARG:NH1	2.26	0.49
1:B:193:GLU:HB2	1:B:269:VAL:CG2	2.43	0.49
1:B:100:GLN:HE21	1:B:100:GLN:C	2.21	0.48
1:B:182:MSE:HE1	1:B:303:TRP:HB3	1.94	0.48
1:B:16:SER:HB3	1:B:316:GLY:HA3	1.95	0.48
1:B:191:ALA:HB2	1:B:228:THR:HG23	1.95	0.48
1:A:117:ARG:HA	1:A:120:MSE:HE1	1.95	0.48
1:A:245:SER:OG	1:A:248:GLN:HG3	2.13	0.48
1:B:85:LEU:CD1	1:B:204:PHE:HE2	2.24	0.48
1:B:180:VAL:CG1	1:B:182:MSE:HE3	2.43	0.48
1:A:257:PHE:O	1:A:261:ILE:CD1	2.60	0.48
1:A:27:GLN:HA	3:A:409:HOH:O	2.12	0.48
1:A:232:HIS:CE1	2:A:401:FAD:H6	2.49	0.48
1:B:107:PRO:HD2	1:B:111:ALA:CB	2.43	0.48
1:B:302:ASP:HB3	1:B:312:ALA:HB2	1.95	0.48
1:B:190:LEU:HD13	1:B:254:HIS:HD2	1.78	0.47
1:A:182:MSE:HE1	1:A:303:TRP:CB	2.45	0.47
1:A:49:GLY:O	1:A:51:LEU:HD12	2.15	0.47
1:A:160:ALA:CB	1:A:182:MSE:HE3	2.42	0.47
1:B:63:ARG:HH11	1:B:63:ARG:HB3	1.80	0.47
1:A:173:LEU:O	1:A:177:VAL:HG23	2.15	0.47
1:B:94:SER:N	1:B:95:PRO:CD	2.73	0.47
1:A:165:THR:CG2	1:B:165:THR:OG1	2.62	0.47
1:B:258:ALA:HB2	1:B:265:MSE:HE1	1.95	0.47
1:B:11:GLY:HA3	2:B:402:FAD:H52A	1.96	0.47
1:B:45:ARG:HG3	1:B:45:ARG:HH11	1.79	0.47
1:A:106:LYS:NZ	1:A:221:GLU:HB2	2.30	0.46
1:A:211:ASP:HB2	1:A:234:THR:HA	1.97	0.46
1:A:232:HIS:CE1	2:A:401:FAD:HM73	2.50	0.46
1:A:189:ALA:O	1:A:271:SER:HA	2.15	0.46
1:A:130:GLU:OE2	1:B:278:TYR:CE2	2.68	0.46
1:A:136:GLU:HB3	1:A:137:HIS:CD2	2.51	0.46
1:B:127:ARG:HD3	3:B:415:HOH:O	2.16	0.46
1:A:64:ARG:O	1:A:67:THR:HB	2.16	0.46
1:A:203:CYS:HB3	1:A:213:LEU:CD1	2.45	0.46
1:A:288:GLY:CA	1:A:305:LEU:HD21	2.45	0.46
1:A:306:SER:HB2	1:A:308:ARG:HD3	1.98	0.46
1:A:191:ALA:CB	1:A:228:THR:HG22	2.46	0.46
1:A:12:ILE:O	1:A:16:SER:HB2	2.17	0.45
1:B:85:LEU:HD12	1:B:97:PRO:HD2	1.97	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:254:HIS:C	1:A:254:HIS:ND1	2.74	0.45
1:B:208:SER:OG	1:B:209:PRO:N	2.50	0.45
1:A:217:ARG:HD2	3:A:442:HOH:O	2.15	0.45
1:B:210:LEU:HD13	1:B:213:LEU:HD23	1.98	0.45
1:B:12:ILE:O	1:B:16:SER:HB2	2.17	0.45
1:A:66:ALA:O	1:A:70:LYS:HG3	2.17	0.45
1:B:83:PRO:CG	1:B:85:LEU:HD21	2.46	0.45
1:A:234:THR:HG22	1:A:237:TRP:HB2	1.99	0.45
1:A:95:PRO:CG	1:A:96:SER:H	2.28	0.45
1:A:195:PRO:HG3	1:A:217:ARG:NH2	2.32	0.45
1:A:117:ARG:O	1:A:118:GLY:C	2.58	0.44
1:B:270:PHE:CD1	1:B:270:PHE:C	2.94	0.44
1:A:56:GLN:HG3	1:A:212:TRP:NE1	2.32	0.44
1:B:18:ALA:O	1:B:22:THR:HB	2.17	0.44
1:B:215:ARG:HE	1:B:217:ARG:HE	1.65	0.44
1:B:236:GLN:O	1:B:240:GLN:HG2	2.17	0.44
1:A:72:TRP:HB3	1:A:78:VAL:HG23	1.99	0.44
1:A:133:ARG:HB3	1:A:138:TRP:CE3	2.53	0.44
1:B:83:PRO:HG2	1:B:85:LEU:HD21	2.00	0.44
1:A:234:THR:HG22	1:A:237:TRP:CB	2.48	0.44
1:B:52:ASP:OD2	1:B:222:ARG:NH2	2.42	0.44
1:B:323:LEU:HD23	1:B:323:LEU:O	2.18	0.44
1:A:79:ALA:HB2	1:A:106:LYS:HD3	2.00	0.44
1:B:212:TRP:CZ2	1:B:214:ALA:HB2	2.52	0.44
1:A:129:THR:O	1:A:130:GLU:OE1	2.36	0.44
1:B:116:MSE:HE3	3:B:485:HOH:O	2.17	0.44
1:A:242:LEU:HD13	1:B:132:PHE:CD2	2.53	0.44
1:B:191:ALA:CB	1:B:228:THR:HG23	2.48	0.44
1:B:287:TRP:C	1:B:289:ALA:H	2.26	0.44
1:A:215:ARG:HD2	1:A:229:TRP:CZ2	2.52	0.43
1:A:7:ILE:HD13	1:A:28:VAL:HB	2.00	0.43
1:B:96:SER:HA	1:B:97:PRO:HD3	1.75	0.43
1:B:329:LEU:N	1:B:329:LEU:HD12	2.34	0.43
1:A:96:SER:N	1:A:97:PRO:HD3	2.32	0.43
1:A:106:LYS:HZ3	1:A:221:GLU:HB2	1.84	0.43
1:B:137:HIS:CG	1:B:150:PRO:HB2	2.53	0.43
1:B:53:MSE:HG2	1:B:230:ILE:HD12	2.00	0.43
1:B:142:ASP:OD1	1:B:142:ASP:C	2.62	0.43
1:A:165:THR:HG21	1:B:165:THR:OG1	2.19	0.43
1:A:203:CYS:HB3	1:A:213:LEU:CG	2.49	0.43
1:A:43:SER:HB2	1:A:110:SER:OG	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:233:ALA:O	1:B:234:THR:C	2.61	0.43
1:B:215:ARG:HD3	1:B:229:TRP:CZ2	2.54	0.43
1:B:87:ASN:O	1:B:92:ARG:HA	2.18	0.43
1:A:242:LEU:HD22	1:A:275:ARG:HD2	2.01	0.42
1:A:95:PRO:CG	1:A:96:SER:N	2.82	0.42
1:A:8:ILE:HD12	1:A:156:ILE:HD13	2.01	0.42
1:B:200:MSE:C	1:B:201:GLN:CG	2.91	0.42
1:A:165:THR:HG23	1:B:165:THR:HG23	2.01	0.42
1:B:113:THR:HG23	3:B:412:HOH:O	2.18	0.42
1:B:63:ARG:HB3	1:B:63:ARG:NH1	2.34	0.42
1:A:79:ALA:HB2	1:A:106:LYS:CE	2.50	0.42
1:A:98:ASP:C	1:A:100:GLN:N	2.77	0.42
1:B:258:ALA:O	1:B:261:ILE:HD12	2.19	0.42
1:A:93:LEU:HD12	1:A:94:SER:O	2.20	0.42
1:B:129:THR:OG1	1:B:130:GLU:HG2	2.20	0.42
1:B:234:THR:HG22	1:B:237:TRP:N	2.17	0.42
1:B:326:HIS:C	1:B:328:GLN:N	2.77	0.42
1:A:65:PHE:CE2	1:A:310:GLU:HB2	2.55	0.42
1:A:232:HIS:HE1	2:A:401:FAD:H6	1.83	0.42
1:A:156:ILE:HD12	1:A:166:LEU:HB3	2.02	0.42
1:B:209:PRO:HB3	1:B:237:TRP:CD2	2.55	0.42
2:B:402:FAD:HM71	2:B:402:FAD:HM83	1.91	0.42
1:A:182:MSE:HE1	1:A:303:TRP:CG	2.55	0.41
2:A:401:FAD:H9	2:A:401:FAD:H1'1	1.81	0.41
1:B:11:GLY:CA	2:B:402:FAD:H52A	2.50	0.41
1:B:15:LEU:HA	1:B:18:ALA:HB3	2.01	0.41
1:B:188:VAL:HA	1:B:272:LEU:O	2.19	0.41
1:B:246:ARG:O	1:B:250:ILE:HG13	2.18	0.41
1:B:239:ARG:HH11	1:B:239:ARG:HG2	1.85	0.41
1:A:180:VAL:CG1	1:A:182:MSE:HE2	2.50	0.41
1:A:195:PRO:CG	1:A:217:ARG:HH22	2.34	0.41
1:B:28:VAL:O	1:B:121:PRO:HG2	2.20	0.41
1:B:62:ASP:OD1	1:B:64:ARG:N	2.53	0.41
1:B:63:ARG:O	1:B:63:ARG:HG2	2.20	0.41
1:A:84:LEU:HD13	1:A:86:TYR:HE1	1.85	0.41
1:A:92:ARG:NH1	1:A:262:ASP:CG	2.78	0.41
1:B:234:THR:CG2	1:B:236:GLN:HB2	2.50	0.41
1:B:22:THR:HG22	1:B:23:ALA:N	2.35	0.41
1:A:327:LEU:HD23	1:A:327:LEU:HA	1.83	0.41
1:B:261:ILE:HD12	1:B:261:ILE:O	2.19	0.41
1:A:54:GLY:HA2	2:A:401:FAD:HM72	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:279:ALA:HA	2:A:401:FAD:HM81	2.03	0.41
1:A:328:GLN:HE21	1:A:328:GLN:CA	2.28	0.41
1:A:106:LYS:HB3	1:A:107:PRO:CD	2.50	0.41
1:A:156:ILE:HD12	1:A:166:LEU:CB	2.51	0.40
1:B:182:MSE:CE	1:B:281:PRO:HA	2.50	0.40
1:A:30:LEU:O	1:A:122:VAL:HA	2.21	0.40
1:B:208:SER:OG	1:B:209:PRO:CD	2.70	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	326/336 (97%)	304 (93%)	16 (5%)	6 (2%)	6	12
1	B	326/336 (97%)	297 (91%)	24 (7%)	5 (2%)	8	16
All	All	652/672 (97%)	601 (92%)	40 (6%)	11 (2%)	7	13

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	136	GLU
1	A	95	PRO
1	B	97	PRO
1	B	99	GLU
1	A	90	ALA
1	A	94	SER
1	B	234	THR
1	B	225	THR
1	A	266	PRO
1	B	106	LYS
1	A	107	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	256/255 (100%)	234 (91%)	22 (9%)	10	21
1	B	256/255 (100%)	227 (89%)	29 (11%)	5	12
All	All	512/510 (100%)	461 (90%)	51 (10%)	7	15

All (51) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	THR
1	A	27	GLN
1	A	80	GLU
1	A	84	LEU
1	A	92	ARG
1	A	128	ILE
1	A	129	THR
1	A	155	ILE
1	A	165	THR
1	A	173	LEU
1	A	198	THR
1	A	205	VAL
1	A	217	ARG
1	A	226	LEU
1	A	242	LEU
1	A	247	GLU
1	A	251	GLU
1	A	254	HIS
1	A	260	LEU
1	A	261	ILE
1	A	290	LEU
1	A	328	GLN
1	B	3	VAL
1	B	22	THR
1	B	51	LEU
1	B	73	GLN
1	B	87	ASN

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Mol	Chain	Res	Type
1	B	93	LEU
1	B	100	GLN
1	B	120	MSE
1	B	159	PRO
1	B	181	LYS
1	B	197	GLN
1	B	206	GLN
1	B	210	LEU
1	B	213	LEU
1	B	225	THR
1	B	226	LEU
1	B	228	THR
1	B	231	LEU
1	B	234	THR
1	B	240	GLN
1	B	242	LEU
1	B	260	LEU
1	B	261	ILE
1	B	277	LEU
1	B	286	GLU
1	B	290	LEU
1	B	294	ASP
1	B	297	ILE
1	B	328	GLN

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

Mol	Chain	Res	Type
1	A	26	HIS
1	A	73	GLN
1	A	77	HIS
1	A	89	HIS
1	A	100	GLN
1	A	137	HIS
1	A	206	GLN
1	A	236	GLN
1	A	240	GLN
1	A	252	HIS
1	A	326	HIS
1	A	328	GLN
1	B	75	GLN
1	B	77	HIS

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Mol	Chain	Res	Type
1	B	100	GLN
1	B	137	HIS
1	B	139	ASN
1	B	153	HIS
1	B	232	HIS
1	B	241	ASN
1	B	248	GLN
1	B	252	HIS
1	B	254	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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	FAD	A	401	-	58,58,58	1.24	5 (8%)	85,89,89	1.69	19 (22%)
2	FAD	B	402	-	58,58,58	1.22	4 (6%)	85,89,89	1.68	19 (22%)

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	FAD	A	401	-	-	2/34/50/50	0/6/6/6
2	FAD	B	402	-	-	2/34/50/50	0/6/6/6

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	FAD	C4X-N5	4.57	1.40	1.30
2	B	402	FAD	C4X-N5	4.47	1.40	1.30
2	B	402	FAD	C10-N1	3.59	1.40	1.33
2	A	401	FAD	C10-N1	3.54	1.40	1.33
2	A	401	FAD	C5A-N7A	-3.11	1.33	1.39
2	B	402	FAD	C5A-N7A	-3.00	1.33	1.39
2	A	401	FAD	C8A-N9A	-2.34	1.33	1.37
2	B	402	FAD	C8A-N9A	-2.31	1.33	1.37
2	A	401	FAD	P-O3P	2.01	1.61	1.59

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	402	FAD	C5A-C4A-N3A	-5.45	119.22	126.72
2	A	401	FAD	C5A-C4A-N3A	-5.36	119.34	126.72
2	B	402	FAD	N3A-C2A-N1A	-4.79	121.33	128.58
2	A	401	FAD	N3A-C2A-N1A	-4.62	121.59	128.58
2	A	401	FAD	C4B-O4B-C1B	-4.18	100.24	109.47
2	B	402	FAD	N3A-C4A-N9A	3.78	133.59	127.17
2	B	402	FAD	C2A-N3A-C4A	3.77	121.03	111.83
2	B	402	FAD	C4B-O4B-C1B	-3.74	101.21	109.47
2	A	401	FAD	C2A-N3A-C4A	3.65	120.74	111.83
2	A	401	FAD	N3A-C4A-N9A	3.43	133.00	127.17
2	A	401	FAD	C4-N3-C2	-3.27	119.84	125.64
2	B	402	FAD	N9A-C8A-N7A	-3.24	109.33	113.94
2	A	401	FAD	C4A-C5A-N7A	-3.18	106.94	110.58
2	B	402	FAD	O4B-C1B-C2B	-3.16	99.85	106.62
2	B	402	FAD	C4-N3-C2	-3.16	120.03	125.64
2	B	402	FAD	C4A-C5A-N7A	-3.05	107.10	110.58
2	B	402	FAD	C5A-N7A-C8A	3.04	108.23	103.45
2	A	401	FAD	C5A-N7A-C8A	2.93	108.05	103.45
2	A	401	FAD	N9A-C8A-N7A	-2.92	109.79	113.94
2	B	402	FAD	C9A-C5X-N5	-2.80	119.48	122.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	FAD	C9A-C5X-N5	-2.77	119.52	122.45
2	A	401	FAD	C4X-C4-N3	2.73	120.20	113.25
2	A	401	FAD	O4B-C1B-C2B	-2.72	100.79	106.62
2	A	401	FAD	O4B-C1B-N9A	2.65	113.18	108.09
2	B	402	FAD	C4X-C4-N3	2.65	119.99	113.25
2	A	401	FAD	O4-C4-C4X	-2.59	119.68	126.53
2	B	402	FAD	O4-C4-C4X	-2.49	119.96	126.53
2	A	401	FAD	C5X-C9A-N10	2.41	120.14	117.97
2	A	401	FAD	C10-C4X-N5	-2.37	119.96	124.81
2	A	401	FAD	O4B-C4B-C5B	2.35	116.86	109.33
2	A	401	FAD	C5'-C4'-C3'	-2.33	107.82	112.22
2	B	402	FAD	O4B-C1B-N9A	2.32	112.55	108.09
2	B	402	FAD	C10-C4X-N5	-2.32	120.06	124.81
2	B	402	FAD	C5X-C9A-N10	2.32	120.06	117.97
2	B	402	FAD	C4A-N9A-C8A	2.30	108.16	105.74
2	A	401	FAD	C4X-C10-N10	2.27	119.73	116.48
2	B	402	FAD	O4B-C4B-C5B	2.24	116.50	109.33
2	B	402	FAD	C4X-C10-N10	2.10	119.50	116.48

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	401	FAD	C5B-O5B-PA-O3P
2	B	402	FAD	PA-O3P-P-O5'
2	A	401	FAD	P-O3P-PA-O2A
2	B	402	FAD	P-O3P-PA-O2A

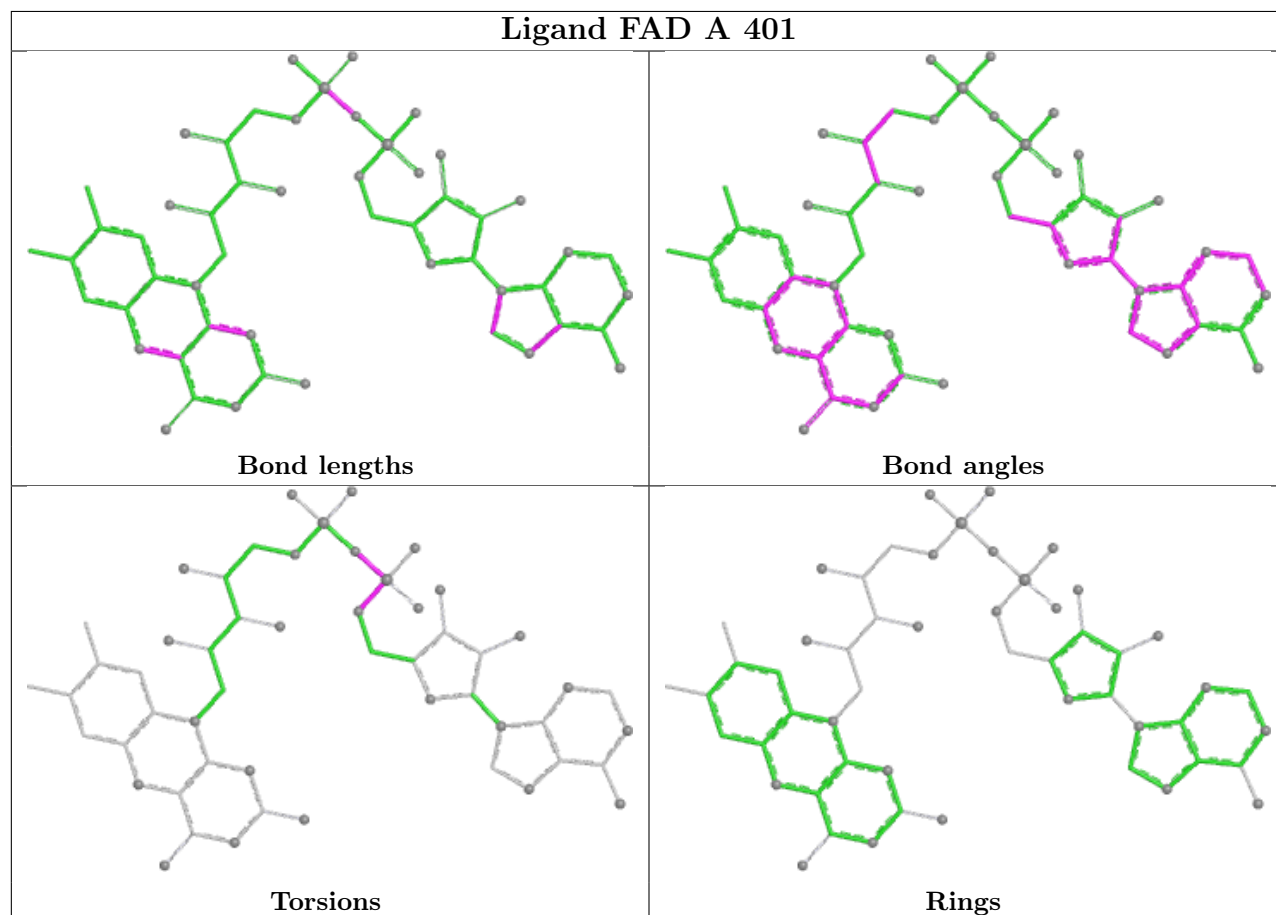
There are no ring outliers.

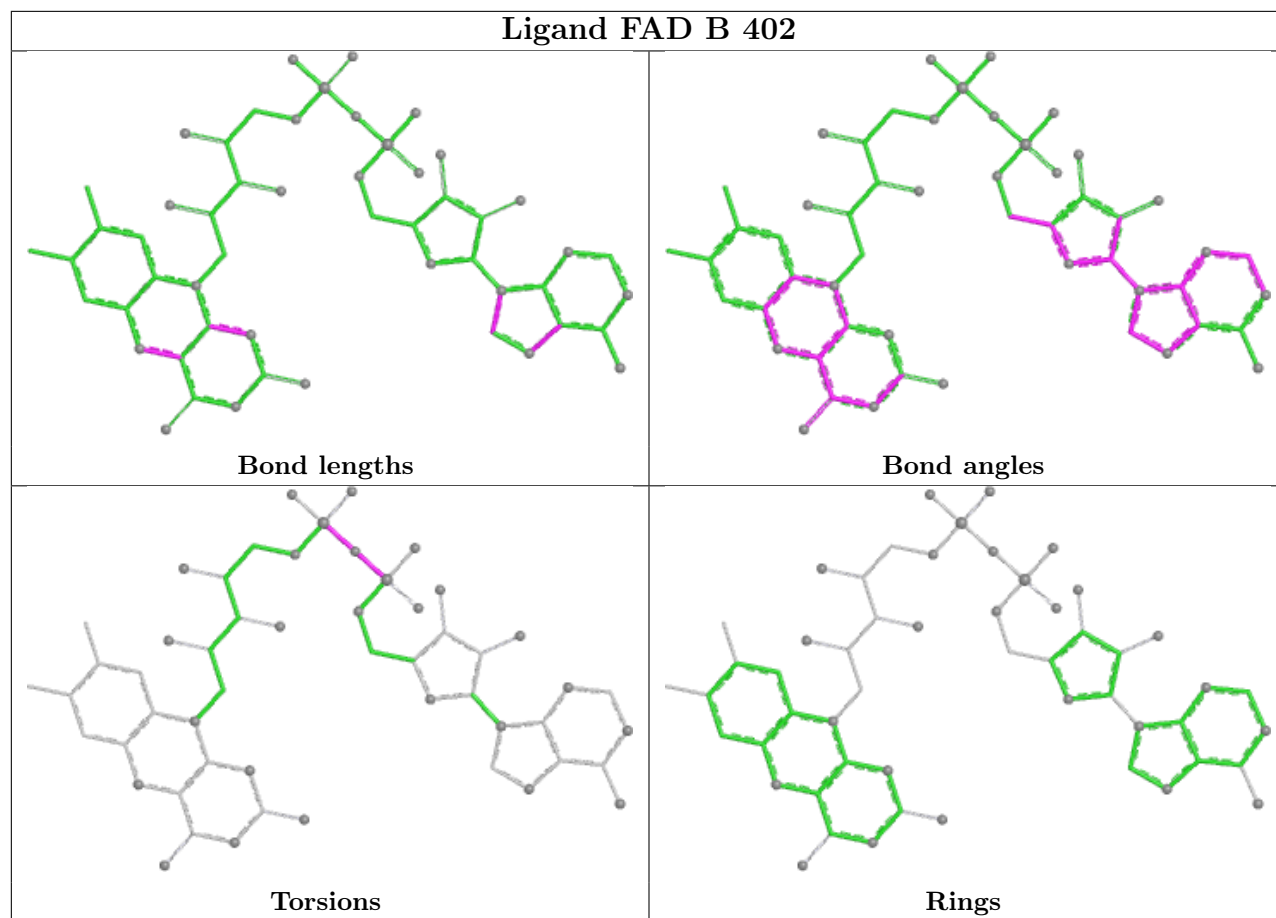
2 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	FAD	7	0
2	B	402	FAD	6	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring

in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	320/336 (95%)	0.26	15 (4%) 36 32	17, 37, 83, 121	0
1	B	320/336 (95%)	0.48	31 (9%) 13 11	19, 38, 78, 119	0
All	All	640/672 (95%)	0.37	46 (7%) 21 19	17, 38, 80, 121	0

All (46) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	95	PRO	7.3
1	B	90	ALA	7.0
1	B	329	LEU	6.7
1	A	329	LEU	6.5
1	A	97	PRO	6.2
1	B	94	SER	6.1
1	B	97	PRO	5.6
1	B	93	LEU	5.2
1	B	92	ARG	4.6
1	B	99	GLU	4.5
1	A	106	LYS	4.3
1	B	87	ASN	4.3
1	B	264	THR	4.2
1	B	88	PHE	3.5
1	B	106	LYS	3.5
1	A	93	LEU	3.4
1	B	85	LEU	3.4
1	B	210	LEU	3.3
1	B	100	GLN	3.3
1	A	92	ARG	3.3
1	A	264	THR	3.3
1	A	136	GLU	3.0
1	B	89	HIS	3.0
1	A	96	SER	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	209	PRO	2.8
1	B	199	PRO	2.8
1	B	86	TYR	2.7
1	B	261	ILE	2.6
1	A	286	GLU	2.6
1	A	48	ALA	2.5
1	B	83	PRO	2.5
1	B	2	THR	2.4
1	B	197	GLN	2.4
1	A	49	GLY	2.4
1	B	198	THR	2.4
1	B	262	ASP	2.3
1	A	107	PRO	2.3
1	B	71	GLN	2.3
1	A	90	ALA	2.2
1	B	258	ALA	2.2
1	A	46	SER	2.2
1	A	95	PRO	2.2
1	B	203	CYS	2.1
1	B	223	ASP	2.1
1	B	98	ASP	2.1
1	B	228	THR	2.1

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	FAD	A	401	53/53	0.96	0.07	15,28,43,44	0

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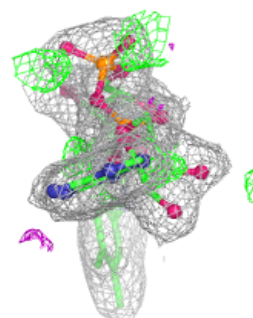
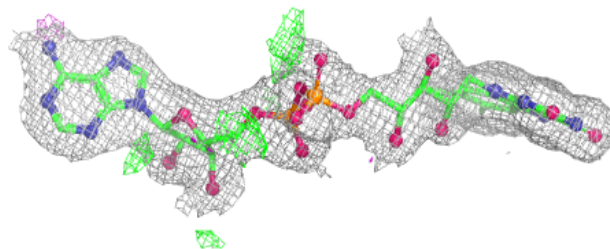
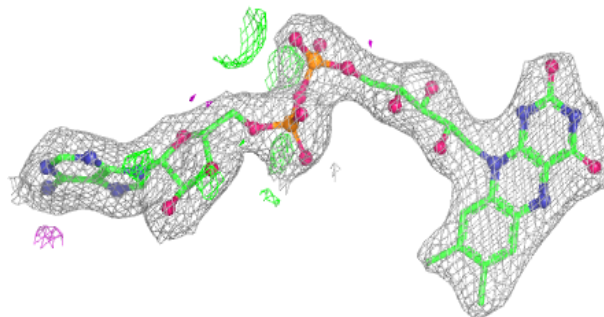
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	FAD	B	402	53/53	0.97	0.07	10,26,43,45	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

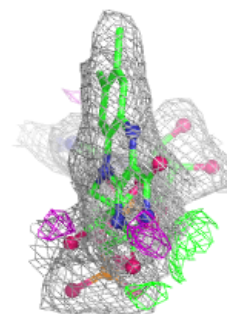
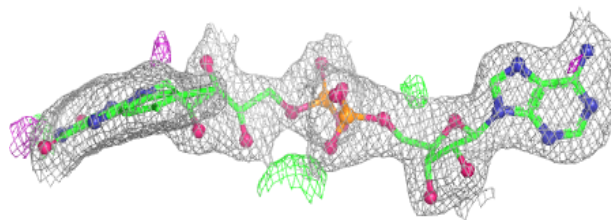
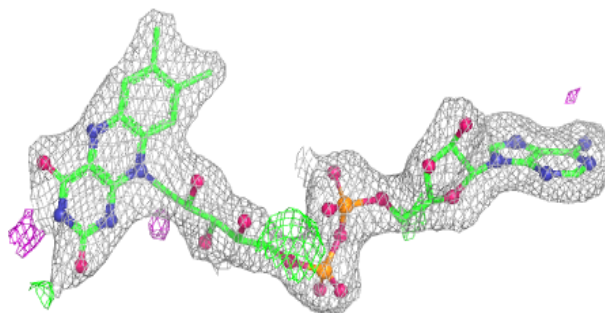
Electron density around FAD A 401:

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



Electron density around FAD B 402:

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



6.5 Other polymers [i](#)

There are no such residues in this entry.