



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 15, 2017 – 03:11 am GMT

PDB ID : 3P4R
Title : Crystal structure of Menaquinol:fumarate oxidoreductase in complex with glutarate
Authors : Tomasiak, T.M.; Archuleta, T.L.; Andrell, J.; Luna-Chavez, C.; Davis, T.A.; Sarwar, M.; Ham, A.J.; McDonald, W.H.; Yankowskaya, V.; Stern, H.A.; Johnston, J.N.; Maklashina, E.; Cecchini, G.; Iverson, T.M.
Deposited on : 2010-10-07
Resolution : 3.05 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 : trunk28620
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) : recalc28949

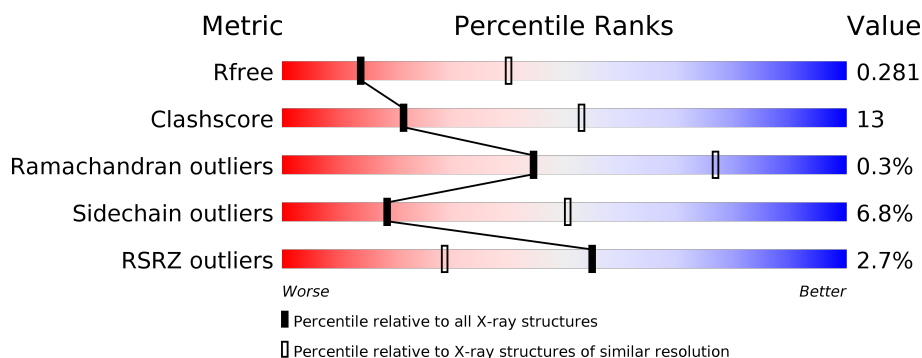
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.05 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	1348 (3.10-3.02)
Clashscore	112137	1462 (3.10-3.02)
Ramachandran outliers	110173	1410 (3.10-3.02)
Sidechain outliers	110143	1410 (3.10-3.02)
RSRZ outliers	101464	1355 (3.10-3.02)

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	577	<div> <div>78%</div> <div>20%</div> <div>.</div> </div>
1	M	577	<div> <div>8%</div> <div>63%</div> <div>33%</div> <div>.</div> </div>
2	B	243	<div> <div>82%</div> <div>15%</div> <div>.</div> </div>
2	N	243	<div> <div>3%</div> <div>67%</div> <div>32%</div> <div>.</div> </div>
3	C	130	<div> <div>%</div> <div>77%</div> <div>22%</div> <div>.</div> </div>
3	O	130	<div> <div>%</div> <div>79%</div> <div>19%</div> <div>.</div> </div>

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Mol	Chain	Length	Quality of chain
4	D	119	<div> <div> <div></div> <div>61%</div> <div>29%</div> <div>6%</div> <div></div> </div> <div> <div>2%</div> <div></div> <div></div> <div></div> <div></div> </div> </div>
4	P	119	<div> <div> <div></div> <div>63%</div> <div>24%</div> <div>8%</div> <div>5%</div> </div> <div> <div>2%</div> <div></div> <div></div> <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
6	F3S	B	245	-	-	X	-
6	F3S	N	245	-	-	X	-
7	SF4	B	246	-	-	X	-
7	SF4	N	246	-	-	X	-

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 16804 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 Fumarate reductase flavoprotein subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	577	Total	C	N	O	S	0	0	0
			4449	2775	802	841	31			
1	M	577	Total	C	N	O	S	0	0	0
			4449	2775	802	841	31			

- Molecule 2 is a protein called Fumarate reductase iron-sulfur subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	243	Total	C	N	O	S	0	0	0
			1888	1189	323	357	19			
2	N	243	Total	C	N	O	S	0	0	0
			1888	1189	323	357	19			

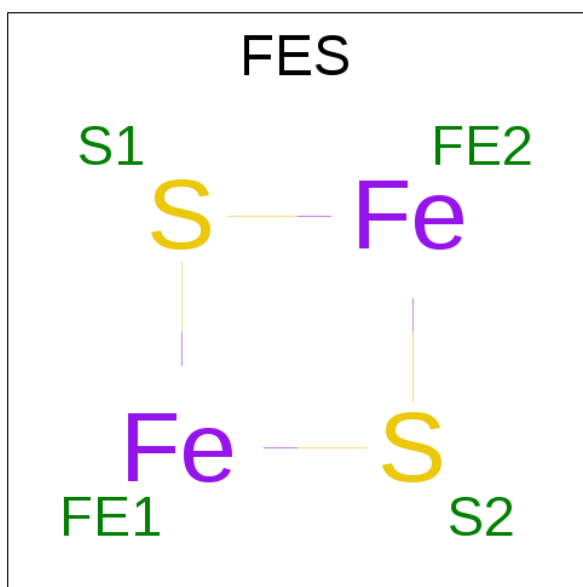
- Molecule 3 is a protein called Fumarate reductase subunit C.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	130	Total	C	N	O	S	0	0	0
			1058	720	166	169	3			
3	O	130	Total	C	N	O	S	0	0	0
			1058	720	166	169	3			

- Molecule 4 is a protein called Fumarate reductase subunit D.

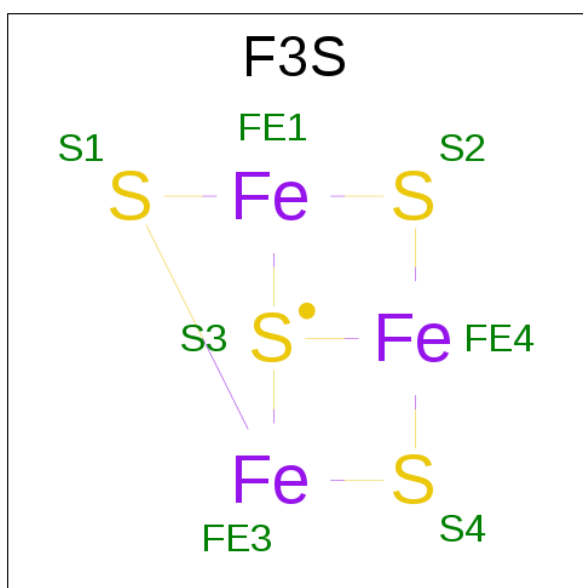
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	119	Total	C	N	O	S	0	0	0
			926	626	151	142	7			
4	P	119	Total	C	N	O	S	0	0	0
			926	626	151	142	7			

- Molecule 5 is FE2/S2 (INORGANIC) CLUSTER (three-letter code: FES) (formula: Fe₂S₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	Fe	S	0	0
			4	2	2		
5	N	1	Total	Fe	S	0	0
			4	2	2		

- Molecule 6 is FE3-S4 CLUSTER (three-letter code: F3S) (formula: Fe₃S₄).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	B	1	Total	Fe	S	0	0
			7	3	4		
6	N	1	Total	Fe	S	0	0
			7	3	4		

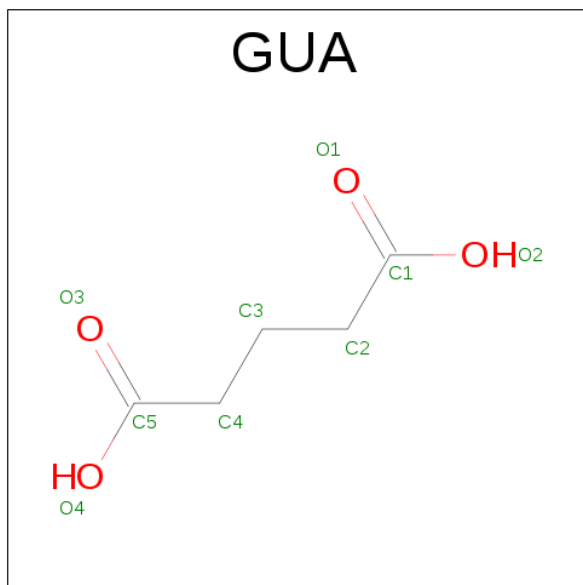
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- The diagram illustrates the coordination of four Fe atoms to the sulfur atom in SF₄. The sulfur atom is represented by a yellow dot, and the four Fe atoms are represented by purple dots. The sulfur atom is labeled 'S' in green, and the Fe atoms are labeled 'Fe' in green. The sulfur atom is bonded to four Fe atoms, forming a tetrahedral geometry. The Fe atoms are labeled 'FE1', 'FE2', 'FE3', and 'FE4' in green. The sulfur atom is labeled 'S1', 'S2', 'S3', and 'S4' in green. The Fe atoms are labeled 'FE1', 'FE2', 'FE3', and 'FE4' in green. The sulfur atom is labeled 'S1', 'S2', 'S3', and 'S4' in green.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	B	1	Total 8	Fe 4	S 4	0	0
7	N	1	Total 8	Fe 4	S 4	0	0

- # FAD
-
- The image displays the chemical structure of Flavin Adenine Dinucleotide (FAD). It consists of a riboflavin (isoalloxazine) ring system, which is a bicyclic structure with a nitrogen atom at the top and a carbonyl group at the bottom. This ring is linked via a pyrophosphate bridge to an adenine ring system, which is a purine derivative. The structure is shown in a 3D representation with various atoms labeled with their respective symbols (C, N, O, P, H) and bonds indicating stereochemistry. The overall structure is a complex molecule with multiple functional groups and a high degree of symmetry.

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

- Molecule 9 is GLUTARIC ACID (three-letter code: GUA) (formula: $C_5H_8O_4$).

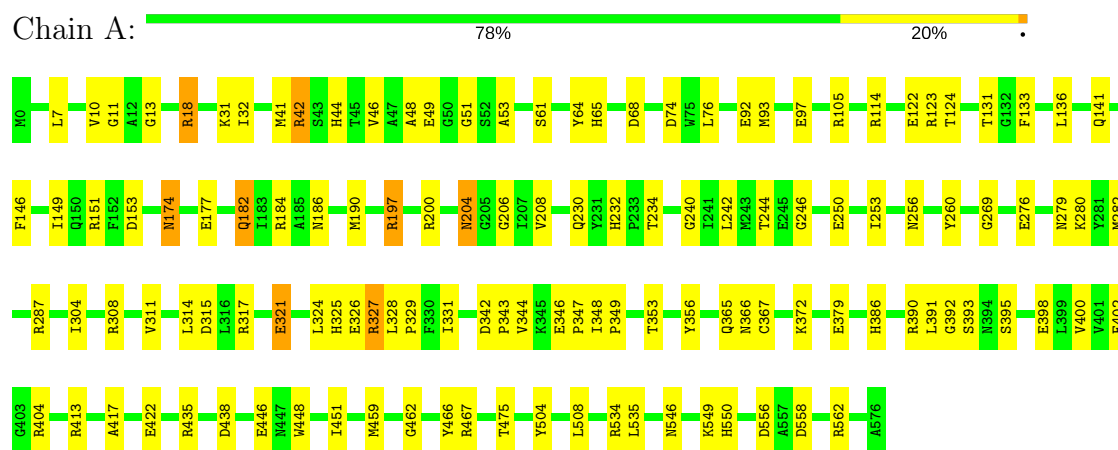


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	A	1	Total	C	O	0	0
			9	5	4		
9	M	1	Total	C	O	0	0
			9	5	4		

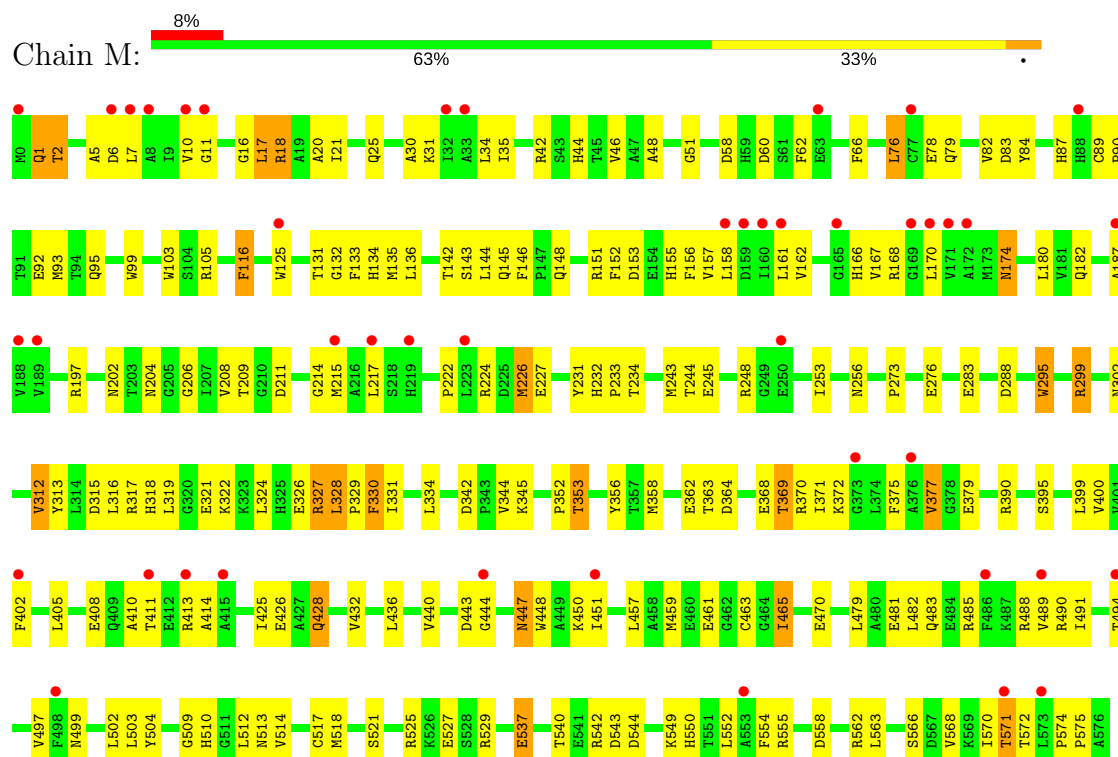
3 Residue-property plots

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

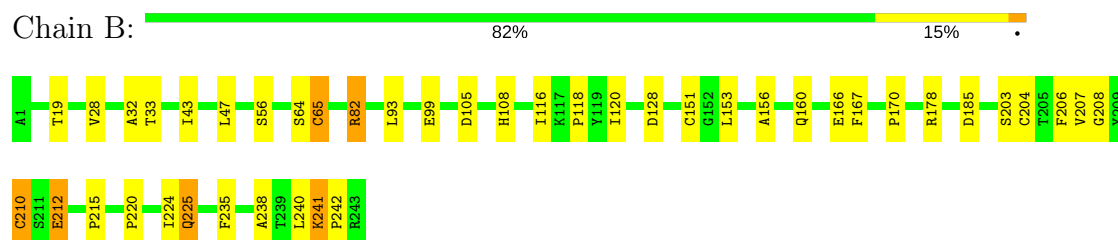
• Molecule 1: Fumarate reductase flavoprotein subunit



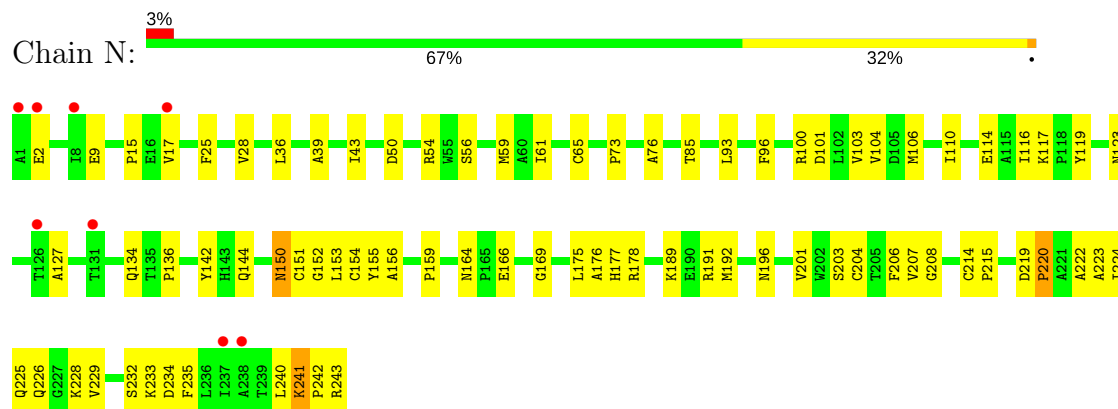
• Molecule 1: Fumarate reductase flavoprotein subunit



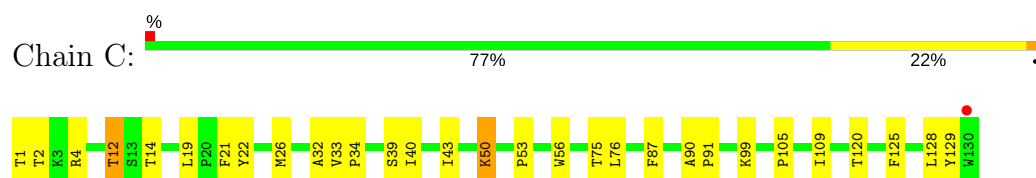
- Molecule 2: Fumarate reductase iron-sulfur subunit



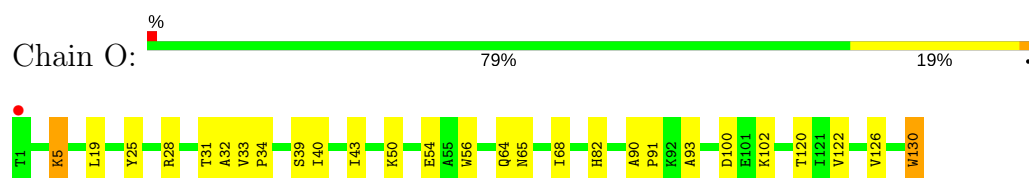
- Molecule 2: Fumarate reductase iron-sulfur subunit



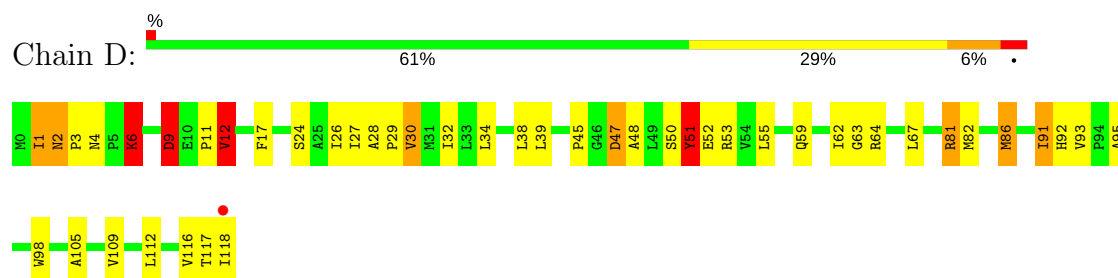
- Molecule 3: Fumarate reductase subunit C



- Molecule 3: Fumarate reductase subunit C

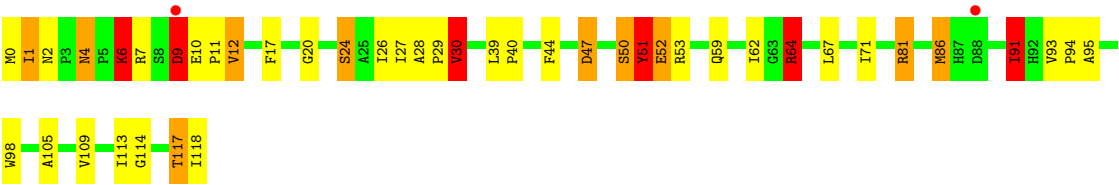


- Molecule 4: Fumarate reductase subunit D



- Molecule 4: Fumarate reductase subunit D





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	96.68Å 138.02Å 269.56Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.00 – 3.05 91.00 – 3.05	Depositor EDS
% Data completeness (in resolution range)	95.3 (40.00-3.05) 94.6 (91.00-3.05)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.05 (at 3.07Å)	Xtriage
Refinement program	REFMAC5	Depositor
R, R_{free}	0.248 , 0.284 0.247 , 0.281	Depositor DCC
R_{free} test set	1194 reflections (1.85%)	DCC
Wilson B-factor (Å ²)	78.8	Xtriage
Anisotropy	0.452	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 41.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	16804	wwPDB-VP
Average B, all atoms (Å ²)	95.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.71% 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: GUA, F3S, SF4, FAD, FES

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	0.39	0/4541	0.56	0/6139
1	M	0.34	0/4541	0.54	0/6139
2	B	0.40	1/1931 (0.1%)	0.54	0/2617
2	N	0.35	0/1931	0.52	0/2617
3	C	0.39	0/1094	0.53	0/1496
3	O	0.39	0/1094	0.52	0/1496
4	D	1.54	16/956 (1.7%)	1.50	18/1303 (1.4%)
4	P	1.55	20/956 (2.1%)	1.46	18/1303 (1.4%)
All	All	0.63	37/17044 (0.2%)	0.71	36/23110 (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
4	D	0	2
4	P	0	2
All	All	0	4

All (37) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	51	TYR	CE1-CZ	-20.60	1.11	1.38
4	P	51	TYR	CE2-CZ	-19.42	1.13	1.38
4	D	51	TYR	CE2-CZ	-18.05	1.15	1.38
4	D	51	TYR	CG-CD1	-15.85	1.18	1.39
4	P	9	ASP	CG-OD1	15.19	1.60	1.25
4	P	51	TYR	CZ-OH	15.17	1.63	1.37
4	P	4	ASN	CB-CG	12.51	1.79	1.51
4	D	51	TYR	CG-CD2	-11.29	1.24	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	P	51	TYR	CB-CG	-11.26	1.34	1.51
4	D	6	LYS	CD-CE	10.80	1.78	1.51
4	P	47	ASP	CG-OD2	10.74	1.50	1.25
4	P	81	ARG	CZ-NH1	10.18	1.46	1.33
4	D	51	TYR	CZ-OH	9.69	1.54	1.37
4	P	4	ASN	CG-ND2	8.29	1.53	1.32
4	D	4	ASN	CB-CG	7.84	1.69	1.51
4	D	47	ASP	CG-OD1	7.68	1.43	1.25
4	P	64	ARG	CZ-NH1	7.61	1.43	1.33
2	B	65	CYS	CB-SG	-7.57	1.69	1.82
4	P	6	LYS	CD-CE	7.22	1.69	1.51
4	D	81	ARG	CZ-NH1	6.85	1.42	1.33
4	P	47	ASP	CA-CB	6.75	1.68	1.53
4	P	51	TYR	CD2-CE2	6.43	1.49	1.39
4	P	52	GLU	CG-CD	-6.35	1.42	1.51
4	P	51	TYR	CE1-CZ	-6.09	1.30	1.38
4	D	12	VAL	CB-CG1	6.06	1.65	1.52
4	P	6	LYS	CE-NZ	5.96	1.64	1.49
4	D	17	PHE	CE1-CZ	-5.92	1.26	1.37
4	P	17	PHE	CE1-CZ	-5.76	1.26	1.37
4	D	52	GLU	CD-OE1	5.74	1.31	1.25
4	D	81	ARG	CZ-NH2	5.66	1.40	1.33
4	P	30	VAL	CB-CG2	-5.50	1.41	1.52
4	D	17	PHE	CG-CD2	-5.46	1.30	1.38
4	P	17	PHE	CG-CD2	-5.40	1.30	1.38
4	P	50	SER	CB-OG	5.34	1.49	1.42
4	P	51	TYR	CG-CD2	-5.17	1.32	1.39
4	D	51	TYR	CD1-CE1	5.11	1.47	1.39
4	D	51	TYR	N-CA	-5.06	1.36	1.46

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	51	TYR	CB-CG-CD2	28.09	137.85	121.00
4	P	47	ASP	CB-CG-OD2	22.25	138.33	118.30
4	D	51	TYR	CD1-CG-CD2	-15.67	100.66	117.90
4	P	47	ASP	CB-CG-OD1	-15.17	104.64	118.30
4	P	51	TYR	CD1-CG-CD2	-14.84	101.58	117.90
4	D	51	TYR	CD1-CE1-CZ	14.24	132.62	119.80
4	P	51	TYR	CG-CD2-CE2	14.14	132.61	121.30
4	D	51	TYR	CZ-CE2-CD2	13.79	132.21	119.80
4	P	51	TYR	CD1-CE1-CZ	13.13	131.62	119.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	P	51	TYR	CB-CG-CD1	12.95	128.77	121.00
4	D	51	TYR	CE1-CZ-CE2	-12.67	99.52	119.80
4	P	81	ARG	NE-CZ-NH2	-11.68	114.46	120.30
4	D	51	TYR	N-CA-CB	-9.03	94.34	110.60
4	D	81	ARG	NE-CZ-NH2	-8.25	116.17	120.30
4	D	51	TYR	CG-CD1-CE1	8.23	127.89	121.30
4	P	91	ILE	CB-CG1-CD1	7.91	136.04	113.90
4	P	47	ASP	CA-CB-CG	-7.17	97.61	113.40
4	P	6	LYS	CD-CE-NZ	6.95	127.69	111.70
4	D	6	LYS	CD-CE-NZ	6.87	127.49	111.70
4	D	51	TYR	CB-CG-CD1	-6.59	117.05	121.00
4	P	62	ILE	CG1-CB-CG2	-6.55	96.99	111.40
4	P	9	ASP	CB-CG-OD1	-6.11	112.80	118.30
4	P	81	ARG	NE-CZ-NH1	6.07	123.33	120.30
4	D	52	GLU	CG-CD-OE1	5.85	129.99	118.30
4	P	64	ARG	NE-CZ-NH1	5.83	123.22	120.30
4	P	51	TYR	CE1-CZ-CE2	-5.79	110.53	119.80
4	P	52	GLU	CG-CD-OE2	-5.77	106.77	118.30
4	D	51	TYR	CB-CA-C	5.68	121.76	110.40
4	P	81	ARG	CD-NE-CZ	-5.67	115.67	123.60
4	D	47	ASP	CB-CG-OD1	-5.62	113.25	118.30
4	D	47	ASP	CA-CB-CG	-5.58	101.11	113.40
4	D	9	ASP	CB-CG-OD1	-5.33	113.50	118.30
4	D	30	VAL	CG1-CB-CG2	-5.19	102.59	110.90
4	D	4	ASN	OD1-CG-ND2	-5.12	110.12	121.90
4	D	51	TYR	CG-CD2-CE2	5.10	125.38	121.30
4	P	52	GLU	CG-CD-OE1	5.09	128.47	118.30

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
4	D	51	TYR	Sidechain
4	D	9	ASP	Sidechain
4	P	51	TYR	Sidechain
4	P	9	ASP	Sidechain

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	4449	0	4335	91	0
1	M	4449	0	4335	163	0
2	B	1888	0	1841	37	0
2	N	1888	0	1838	63	0
3	C	1058	0	1108	25	0
3	O	1058	0	1108	30	0
4	D	926	0	971	44	0
4	P	926	0	971	47	0
5	B	4	0	0	0	0
5	N	4	0	0	0	0
6	B	7	0	0	6	0
6	N	7	0	0	3	0
7	B	8	0	0	2	0
7	N	8	0	0	6	0
8	A	53	0	31	8	0
8	M	53	0	30	13	0
9	A	9	0	0	1	0
9	M	9	0	0	2	0
All	All	16804	0	16568	449	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 13.

All (449) 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)
4:D:6:LYS:CD	4:D:6:LYS:CE	1.78	1.58
4:D:91:ILE:CG1	4:D:91:ILE:CD1	1.76	1.57
4:P:4:ASN:CG	4:P:4:ASN:CB	1.79	1.50
4:P:51:TYR:OH	4:P:51:TYR:CZ	1.63	1.50
1:M:44:HIS:NE2	8:M:601:FAD:HM82	1.44	1.32
1:M:11:GLY:HA2	8:M:601:FAD:H1B	1.25	1.18
2:N:214:CYS:SG	7:N:246:SF4:S1	2.43	1.17
4:D:92:HIS:HB3	2:N:243:ARG:HB3	1.28	1.15
1:A:44:HIS:NE2	8:A:601:FAD:HM82	1.63	1.12
1:M:44:HIS:NE2	8:M:601:FAD:C8M	2.19	1.05
1:A:44:HIS:NE2	8:A:601:FAD:C8M	2.20	1.05
1:A:372:LYS:HE3	1:A:413:ARG:HH12	1.25	1.00
2:B:204:CYS:HG	6:B:245:F3S:FE1	0.79	0.97

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:151:CYS:HG	7:B:246:SF4:FE1	0.78	0.97
4:D:55:LEU:HG	4:D:59:GLN:HE21	1.31	0.94
2:N:116:ILE:HG21	2:N:176:ALA:HB2	1.47	0.94
4:D:55:LEU:HG	4:D:59:GLN:NE2	1.83	0.94
2:N:214:CYS:SG	7:N:246:SF4:FE4	1.61	0.92
1:A:44:HIS:CE1	8:A:601:FAD:HM82	2.06	0.90
3:O:120:THR:HG23	4:P:30:VAL:HG23	1.51	0.90
1:M:11:GLY:HA2	8:M:601:FAD:C1B	2.05	0.86
1:M:158:LEU:HD13	1:M:436:LEU:HD11	1.57	0.85
4:P:26:ILE:HG22	4:P:27:ILE:HG13	1.56	0.85
1:M:537:GLU:CD	1:M:537:GLU:H	1.80	0.85
4:P:47:ASP:HB2	4:P:50:SER:OG	1.76	0.84
2:B:32:ALA:O	2:B:82:ARG:NH1	2.10	0.83
4:D:26:ILE:HG22	4:D:27:ILE:HG13	1.60	0.83
7:N:246:SF4:FE3	7:N:246:SF4:S2	1.69	0.83
1:A:92:GLU:OE1	1:A:92:GLU:HA	1.79	0.82
2:B:210:CYS:SG	6:B:245:F3S:FE3	1.71	0.82
4:P:47:ASP:N	4:P:47:ASP:OD1	2.07	0.81
1:M:44:HIS:CE1	8:M:601:FAD:HM82	2.15	0.81
2:N:214:CYS:HG	7:N:246:SF4:FE4	0.96	0.81
1:A:208:VAL:HG12	1:A:208:VAL:O	1.81	0.80
2:N:116:ILE:CG2	2:N:176:ALA:HB2	2.10	0.80
1:M:390:ARG:HD2	1:M:395:SER:HB2	1.64	0.80
1:A:390:ARG:HH22	9:A:577:GUA:C1	1.95	0.79
2:N:222:ALA:O	2:N:226:GLN:HG3	1.82	0.78
4:D:6:LYS:CE	4:D:6:LYS:CG	2.62	0.78
1:M:217:LEU:HG	1:M:555:ARG:HB3	1.65	0.77
4:D:64:ARG:HG2	4:D:64:ARG:HH11	1.49	0.77
3:O:5:LYS:O	3:O:5:LYS:HG3	1.85	0.76
1:M:1:GLN:HE21	1:M:2:THR:H	1.34	0.76
2:B:204:CYS:SG	6:B:245:F3S:FE1	1.76	0.75
4:P:51:TYR:HE1	4:P:52:GLU:OE1	1.69	0.75
2:N:196:ASN:ND2	2:N:234:ASP:OD2	2.20	0.75
3:O:28:ARG:O	3:O:31:THR:HB	1.87	0.75
7:N:246:SF4:S1	7:N:246:SF4:FE4	1.77	0.75
1:M:217:LEU:HG	1:M:555:ARG:CB	2.17	0.74
2:N:154:CYS:SG	7:N:246:SF4:S2	2.85	0.74
1:A:232:HIS:CG	1:A:242:LEU:HD11	2.22	0.74
3:O:56:TRP:CD1	4:P:51:TYR:HB3	2.23	0.74
1:A:321:GLU:OE1	1:A:344:VAL:HG21	1.89	0.73
1:M:2:THR:HA	1:M:182:GLN:O	1.88	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:O:40:ILE:HA	3:O:43:ILE:HD12	1.71	0.72
2:B:151:CYS:SG	7:B:246:SF4:FE1	1.81	0.72
1:M:322:LYS:O	1:M:326:GLU:HB2	1.89	0.72
3:O:31:THR:HG21	3:O:82:HIS:HB2	1.71	0.72
3:O:33:VAL:HB	3:O:34:PRO:HD3	1.71	0.72
1:A:324:LEU:HD23	1:A:328:LEU:HD12	1.71	0.71
1:A:44:HIS:NE2	8:A:601:FAD:HM81	2.02	0.71
4:P:51:TYR:CE1	4:P:52:GLU:HG2	2.25	0.71
2:B:225:GLN:HA	2:B:225:GLN:HE21	1.55	0.71
2:B:210:CYS:HG	6:B:245:F3S:FE3	1.07	0.70
2:N:204:CYS:SG	2:N:224:ILE:HG21	2.31	0.70
1:A:197:ARG:HD2	1:A:206:GLY:HA2	1.73	0.69
2:N:54:ARG:HE	2:N:103:VAL:HG13	1.57	0.69
1:A:324:LEU:HD13	1:A:344:VAL:HG22	1.74	0.69
2:N:28:VAL:HG22	2:N:43:ILE:HD11	1.74	0.69
4:D:92:HIS:HB3	2:N:243:ARG:CB	2.17	0.69
1:M:48:ALA:HB3	1:M:132:GLY:HA3	1.75	0.68
1:M:152:PHE:HB3	1:M:155:HIS:ND1	2.08	0.68
3:C:2:THR:OG1	3:C:4:ARG:HG2	1.92	0.68
1:A:356:TYR:CE1	1:A:379:GLU:HG3	2.28	0.68
1:M:7:LEU:HD21	1:M:411:THR:HA	1.75	0.68
1:M:224:ARG:NH2	1:M:363:THR:O	2.26	0.67
1:A:18:ARG:HG2	1:A:400:VAL:HA	1.77	0.67
1:M:152:PHE:HD2	1:M:155:HIS:CE1	2.12	0.67
2:N:225:GLN:HE21	3:O:93:ALA:HB2	1.58	0.67
1:A:253:ILE:HG13	1:A:315:ASP:HB3	1.76	0.67
2:N:164:ASN:OD1	2:N:166:GLU:HG2	1.95	0.67
1:M:162:VAL:HG21	1:M:371:ILE:HD13	1.76	0.67
1:A:204:ASN:HD22	1:A:204:ASN:N	1.93	0.66
1:M:155:HIS:CD2	1:M:174:ASN:HA	2.31	0.66
2:B:28:VAL:HG22	2:B:43:ILE:HD11	1.77	0.66
3:C:12:THR:HG22	3:C:14:THR:H	1.60	0.66
1:M:525:ARG:HD2	1:M:527:GLU:HG2	1.78	0.66
4:P:44:PHE:HE1	4:P:47:ASP:HA	1.61	0.66
1:M:315:ASP:OD1	1:M:317:ARG:HG3	1.95	0.66
1:A:11:GLY:HA2	8:A:601:FAD:H1B	1.77	0.65
3:C:50:LYS:HB3	4:D:118:ILE:CG2	2.26	0.65
3:C:33:VAL:HB	3:C:34:PRO:HD3	1.79	0.65
1:M:42:ARG:NH2	2:N:150:ASN:O	2.30	0.65
1:A:46:VAL:HG23	1:A:133:PHE:HA	1.78	0.65
4:D:86:MET:HG3	4:D:93:VAL:HG21	1.78	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:P:86:MET:HE3	4:P:91:ILE:HG21	1.79	0.65
1:M:116:PHE:CE2	1:M:245:GLU:HB3	2.32	0.64
1:M:529:ARG:HH21	1:M:542:ARG:HE	1.44	0.64
1:M:92:GLU:HB3	1:M:400:VAL:HB	1.78	0.64
4:D:92:HIS:CB	2:N:243:ARG:HB3	2.17	0.64
2:B:238:ALA:HA	2:B:241:LYS:HD2	1.79	0.64
3:C:50:LYS:CB	4:D:118:ILE:HG22	2.29	0.63
1:M:224:ARG:HH21	1:M:363:THR:H	1.46	0.63
1:A:372:LYS:HE3	1:A:413:ARG:NH1	2.06	0.63
3:C:50:LYS:HB3	4:D:118:ILE:HG22	1.81	0.63
1:M:35:ILE:HD13	1:M:157:VAL:HG23	1.81	0.63
2:N:65:CYS:HB2	2:N:76:ALA:HB3	1.80	0.63
4:D:86:MET:CE	4:D:91:ILE:HG21	2.29	0.63
4:P:86:MET:CE	4:P:91:ILE:HG21	2.29	0.62
1:M:232:HIS:HD2	1:M:234:THR:O	1.83	0.62
1:A:390:ARG:HD2	1:A:395:SER:HB2	1.82	0.62
1:A:398:GLU:HG2	1:A:402:PHE:HD1	1.64	0.62
1:A:208:VAL:CG1	1:A:208:VAL:O	2.48	0.62
1:M:316:LEU:HB3	1:M:319:LEU:HD12	1.82	0.61
1:M:331:ILE:HA	1:M:334:LEU:HD12	1.83	0.61
1:A:546:ASN:O	1:A:549:LYS:HE2	2.00	0.61
1:M:233:PRO:HG2	1:M:248:ARG:HH22	1.64	0.61
1:A:61:SER:HB3	1:A:64:TYR:CD2	2.36	0.61
3:C:120:THR:HA	4:D:30:VAL:HG21	1.83	0.61
3:C:105:PRO:O	3:C:109:ILE:HG12	2.00	0.60
1:A:174:ASN:HD22	1:A:174:ASN:C	2.05	0.60
3:O:50:LYS:HG3	4:P:118:ILE:HA	1.83	0.60
1:A:232:HIS:CD2	1:A:242:LEU:HD11	2.36	0.60
1:M:105:ARG:HD3	2:N:134:GLN:O	2.00	0.60
1:M:222:PRO:HG2	1:M:362:GLU:HB3	1.82	0.60
3:C:120:THR:HG23	4:D:30:VAL:HG22	1.84	0.60
1:M:158:LEU:HD22	1:M:436:LEU:HD21	1.83	0.60
1:M:6:ASP:OD2	1:M:30:ALA:HA	2.03	0.59
3:O:28:ARG:HD2	4:P:81:ARG:NH2	2.18	0.59
1:M:155:HIS:HD2	1:M:174:ASN:HA	1.68	0.59
3:C:120:THR:HA	4:D:30:VAL:CG2	2.32	0.59
1:A:346:GLU:HB2	1:A:347:PRO:CD	2.33	0.59
1:A:462:GLY:HA3	1:A:475:THR:OG1	2.03	0.59
2:B:153:LEU:HD12	2:B:215:PRO:HD3	1.83	0.58
1:M:197:ARG:HD2	1:M:206:GLY:HA2	1.83	0.58
1:A:65:HIS:HA	1:A:123:ARG:NH1	2.18	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:18:ARG:NH1	1:M:92:GLU:OE1	2.36	0.58
1:M:84:TYR:OH	1:M:405:LEU:HD11	2.03	0.58
1:A:269:GLY:HA2	1:A:282:MET:HE1	1.85	0.58
1:A:321:GLU:HA	1:A:321:GLU:OE1	2.03	0.58
1:M:95:GLN:HE21	2:N:127:ALA:HB2	1.69	0.58
1:M:180:LEU:H	1:M:180:LEU:HD12	1.69	0.57
1:A:174:ASN:ND2	1:A:177:GLU:H	2.01	0.57
2:B:120:ILE:CD1	2:B:185:ASP:HB2	2.34	0.57
1:M:44:HIS:CE1	1:M:204:ASN:HA	2.39	0.57
1:M:510:HIS:O	1:M:514:VAL:HG23	2.04	0.57
3:O:31:THR:CG2	3:O:82:HIS:HB2	2.34	0.57
1:A:7:LEU:HD21	1:A:32:ILE:HG12	1.86	0.57
1:M:46:VAL:HG23	1:M:133:PHE:HA	1.87	0.57
3:O:90:ALA:N	3:O:91:PRO:HD2	2.19	0.57
1:M:399:LEU:HD21	8:M:601:FAD:O2P	2.05	0.56
2:N:134:GLN:NE2	2:N:142:TYR:HE2	2.03	0.56
4:D:86:MET:HE2	4:D:91:ILE:HG21	1.85	0.56
2:N:36:LEU:HD23	2:N:76:ALA:HA	1.88	0.56
1:A:232:HIS:HD2	1:A:234:THR:H	1.52	0.56
2:B:242:PRO:HB3	4:P:93:VAL:C	2.26	0.56
4:P:51:TYR:CD1	4:P:52:GLU:N	2.74	0.56
2:N:207:VAL:HG23	6:N:245:F3S:S2	2.46	0.56
2:B:210:CYS:HB2	2:B:220:PRO:HG2	1.87	0.56
3:C:75:THR:HG22	4:D:32:ILE:HD13	1.88	0.56
4:P:47:ASP:CB	4:P:50:SER:OG	2.49	0.56
1:A:246:GLY:O	1:A:250:GLU:HG2	2.06	0.55
2:N:93:LEU:HG	2:N:156:ALA:HB2	1.88	0.55
1:M:214:GLY:HA3	1:M:510:HIS:CD2	2.41	0.55
2:B:212:GLU:HG3	3:C:21:PHE:CE2	2.42	0.55
3:C:53:PRO:HA	4:D:51:TYR:CE1	2.42	0.55
3:O:65:ASN:HB3	3:O:68:ILE:HG12	1.87	0.55
1:M:35:ILE:HG12	8:M:601:FAD:H2A	1.87	0.54
4:P:105:ALA:O	4:P:109:VAL:HG23	2.08	0.54
1:M:405:LEU:HA	1:M:408:GLU:HB2	1.89	0.54
1:A:197:ARG:HB2	1:A:208:VAL:O	2.07	0.54
3:C:87:PHE:O	3:C:109:ILE:HD13	2.07	0.54
1:M:5:ALA:HA	1:M:31:LYS:HB3	1.89	0.54
1:A:279:ASN:O	1:A:280:LYS:HB2	2.07	0.54
1:A:42:ARG:HG2	2:B:64:SER:HB3	1.89	0.54
1:A:61:SER:HB3	1:A:64:TYR:CE2	2.42	0.54
1:M:375:PHE:CZ	1:M:410:ALA:HA	2.42	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:46:VAL:CG2	1:M:133:PHE:HA	2.38	0.54
1:A:46:VAL:HB	1:A:136:LEU:HD23	1.88	0.54
3:C:125:PHE:O	3:C:129:TYR:HB2	2.07	0.54
1:M:166:HIS:CD2	1:M:372:LYS:HB3	2.43	0.54
1:A:314:LEU:HD23	1:A:348:ILE:HD11	1.89	0.54
1:M:529:ARG:NH2	1:M:544:ASP:OD1	2.41	0.54
4:D:105:ALA:O	4:D:109:VAL:HG23	2.08	0.54
1:M:377:VAL:HG11	1:M:402:PHE:O	2.08	0.54
2:N:116:ILE:HG22	2:N:191:ARG:HD3	1.89	0.54
3:O:39:SER:OG	4:P:71:ILE:O	2.25	0.54
3:O:120:THR:HA	4:P:30:VAL:HG21	1.90	0.53
1:M:18:ARG:HG2	1:M:400:VAL:HA	1.89	0.53
1:M:571:THR:HG22	1:M:572:THR:N	2.23	0.53
2:B:225:GLN:HA	2:B:225:GLN:NE2	2.23	0.53
1:M:214:GLY:HA3	1:M:510:HIS:HD2	1.73	0.53
1:M:157:VAL:HG22	1:M:170:LEU:HD13	1.91	0.53
1:M:1:GLN:HE21	1:M:2:THR:N	2.02	0.53
1:M:358:MET:SD	1:M:390:ARG:N	2.82	0.53
1:M:451:ILE:HG23	1:M:482:LEU:HD22	1.91	0.53
2:N:116:ILE:HG21	2:N:176:ALA:CB	2.31	0.52
4:P:44:PHE:CE1	4:P:47:ASP:HA	2.42	0.52
2:B:240:LEU:O	2:B:242:PRO:HD3	2.09	0.52
1:M:168:ARG:HD3	1:M:425:ILE:HD11	1.89	0.52
3:O:32:ALA:HB2	4:P:81:ARG:HD3	1.91	0.52
1:M:35:ILE:HD13	1:M:157:VAL:CG2	2.39	0.52
1:A:51:GLY:HA2	1:A:131:THR:HG21	1.92	0.52
1:M:44:HIS:HE1	1:M:204:ASN:HA	1.74	0.52
1:A:390:ARG:HH11	1:A:395:SER:HB3	1.74	0.52
2:N:159:PRO:HG2	2:N:207:VAL:HG21	1.91	0.52
3:O:50:LYS:HD2	4:P:117:THR:HB	1.92	0.52
4:D:62:ILE:CG2	4:D:63:GLY:N	2.73	0.52
1:A:466:TYR:CD2	1:A:535:LEU:HD21	2.44	0.52
1:M:224:ARG:HH21	1:M:363:THR:N	2.08	0.52
2:N:155:TYR:CE2	2:N:169:GLY:HA3	2.44	0.52
1:M:79:GLN:HB2	1:M:571:THR:HB	1.91	0.51
2:N:106:MET:O	2:N:110:ILE:HG12	2.11	0.51
1:M:330:PHE:HZ	2:N:61:ILE:HB	1.75	0.51
2:N:225:GLN:NE2	3:O:93:ALA:HB2	2.23	0.51
2:B:170:PRO:HA	2:B:224:ILE:HD11	1.93	0.51
1:M:35:ILE:CD1	1:M:157:VAL:HG23	2.41	0.51
1:M:217:LEU:HD21	1:M:517:CYS:SG	2.51	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:N:153:LEU:HD12	2:N:215:PRO:HD3	1.92	0.51
1:A:326:GLU:OE2	1:A:327:ARG:NH1	2.44	0.51
1:M:226:MET:HB3	1:M:518:MET:HA	1.92	0.51
1:A:7:LEU:CD2	1:A:32:ILE:HG12	2.40	0.50
1:M:217:LEU:HG	1:M:555:ARG:HB2	1.90	0.50
1:M:369:THR:HG22	1:M:370:ARG:H	1.77	0.50
3:C:56:TRP:CD1	4:D:51:TYR:HB2	2.46	0.50
4:P:10:GLU:N	4:P:11:PRO:HD2	2.26	0.50
1:M:46:VAL:HB	1:M:136:LEU:HD23	1.94	0.50
1:M:168:ARG:HD3	1:M:425:ILE:CD1	2.42	0.50
1:M:152:PHE:CD2	1:M:155:HIS:CE1	2.97	0.50
1:M:103:TRP:CZ3	1:M:125:TRP:HB3	2.46	0.49
4:P:4:ASN:CA	4:P:4:ASN:CG	2.72	0.49
1:A:13:GLY:H	8:A:601:FAD:H4B	1.76	0.49
4:D:6:LYS:HB3	4:P:6:LYS:HG2	1.94	0.49
1:M:326:GLU:HB3	1:M:327:ARG:HD3	1.94	0.49
1:M:79:GLN:HG3	1:M:570:ILE:HA	1.95	0.49
1:A:287:ARG:NH2	1:A:390:ARG:O	2.46	0.49
1:A:308:ARG:HD2	2:B:33:THR:HG21	1.94	0.49
4:D:112:LEU:O	4:D:116:VAL:HG23	2.11	0.49
4:D:28:ALA:N	4:D:29:PRO:HD2	2.28	0.49
1:M:358:MET:HG2	1:M:390:ARG:HB2	1.93	0.49
1:A:97:GLU:OE1	1:A:105:ARG:NH2	2.37	0.49
1:M:161:LEU:O	1:M:167:VAL:HA	2.13	0.49
1:M:549:LYS:HA	1:M:568:VAL:HG23	1.94	0.49
1:M:99:TRP:CD2	1:M:142:THR:HG21	2.48	0.49
3:O:40:ILE:O	3:O:43:ILE:HB	2.12	0.49
2:N:151:CYS:SG	2:N:152:GLY:N	2.85	0.49
1:A:256:ASN:HD21	1:A:260:TYR:HB3	1.77	0.49
1:A:392:GLY:O	1:A:393:SER:OG	2.27	0.49
2:B:120:ILE:HD12	2:B:185:ASP:HB2	1.95	0.49
2:N:208:GLY:N	6:N:245:F3S:S1	2.85	0.49
1:A:141:GLN:HB3	2:B:118:PRO:O	2.13	0.49
3:C:125:PHE:CE1	3:C:129:TYR:HD1	2.31	0.49
4:D:86:MET:HE3	4:D:91:ILE:HG21	1.95	0.49
1:M:11:GLY:CA	8:M:601:FAD:C1B	2.86	0.49
4:P:86:MET:HG3	4:P:93:VAL:HG21	1.93	0.49
1:M:144:LEU:HD13	2:N:114:GLU:HG2	1.94	0.48
3:O:130:TRP:HB3	4:P:53:ARG:HH21	1.78	0.48
1:A:48:ALA:HA	8:A:601:FAD:C6	2.43	0.48
1:A:366:ASN:O	1:A:367:CYS:HB2	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:390:ARG:NH1	1:A:395:SER:HB3	2.27	0.48
3:C:90:ALA:N	3:C:91:PRO:HD2	2.27	0.48
1:A:448:TRP:HE3	1:A:508:LEU:HD22	1.79	0.48
1:M:48:ALA:HB3	1:M:132:GLY:CA	2.43	0.48
1:M:491:ILE:HD13	1:M:502:LEU:HD12	1.94	0.48
1:M:549:LYS:HB2	1:M:566:SER:O	2.12	0.48
1:A:68:ASP:HB3	1:A:391:LEU:HD21	1.96	0.48
1:A:435:ARG:O	1:A:438:ASP:HB2	2.14	0.48
4:D:62:ILE:HG22	4:D:63:GLY:N	2.29	0.48
2:B:167:PHE:CD1	2:B:203:SER:HB2	2.49	0.48
1:M:497:VAL:HG21	2:N:15:PRO:HG2	1.96	0.48
2:N:225:GLN:HG3	3:O:93:ALA:HA	1.96	0.48
2:N:201:VAL:O	2:N:228:LYS:HE3	2.14	0.48
4:D:95:ALA:HB1	4:D:98:TRP:HB2	1.95	0.48
1:M:157:VAL:HB	8:M:601:FAD:N1A	2.28	0.48
2:N:225:GLN:HE21	3:O:93:ALA:CB	2.24	0.48
2:N:155:TYR:CZ	2:N:169:GLY:HA3	2.49	0.48
1:M:273:PRO:HG2	1:M:276:GLU:HG3	1.96	0.47
1:M:116:PHE:HE2	9:M:577:GUA:O3	1.96	0.47
4:P:51:TYR:CD2	4:P:51:TYR:N	2.81	0.47
2:N:134:GLN:NE2	2:N:142:TYR:CE2	2.83	0.47
1:M:483:GLN:NE2	1:M:555:ARG:HH12	2.12	0.47
3:C:125:PHE:CD1	3:C:129:TYR:CD1	3.02	0.47
1:M:133:PHE:HE2	1:M:134:HIS:CE1	2.31	0.47
1:M:21:ILE:O	1:M:25:GLN:HG3	2.14	0.47
2:N:50:ASP:O	2:N:100:ARG:NH1	2.46	0.47
1:M:328:LEU:N	1:M:329:PRO:CD	2.77	0.47
3:O:64:GLN:NE2	4:P:40:PRO:O	2.43	0.47
4:P:9:ASP:O	4:P:12:VAL:HG23	2.15	0.47
3:O:25:TYR:HD1	3:O:28:ARG:NH2	2.12	0.47
1:A:346:GLU:HB2	1:A:347:PRO:HD3	1.96	0.47
2:B:225:GLN:CA	2:B:225:GLN:HE21	2.24	0.46
1:M:116:PHE:CE2	9:M:577:GUA:O3	2.68	0.46
1:M:253:ILE:HG12	1:M:318:HIS:CE1	2.50	0.46
1:M:550:HIS:HB2	1:M:566:SER:OG	2.15	0.46
2:N:39:ALA:O	2:N:43:ILE:HG12	2.15	0.46
3:C:32:ALA:HB2	4:D:81:ARG:HD3	1.97	0.46
3:C:33:VAL:HA	4:D:82:MET:HE2	1.96	0.46
1:M:256:ASN:HB2	1:M:302:ASN:O	2.15	0.46
1:A:114:ARG:O	1:A:124:THR:HB	2.16	0.46
1:A:328:LEU:N	1:A:329:PRO:CD	2.78	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:N:9:GLU:HG3	2:N:25:PHE:CE2	2.51	0.46
1:A:304:ILE:HD12	1:A:304:ILE:H	1.81	0.46
1:M:35:ILE:HD12	1:M:155:HIS:HB3	1.98	0.46
1:A:151:ARG:NH1	1:A:153:ASP:OD2	2.48	0.46
1:M:142:THR:O	1:M:145:GLN:HG2	2.16	0.46
2:N:136:PRO:HB2	3:O:100:ASP:OD1	2.16	0.46
1:M:499:ASN:O	1:M:503:LEU:HG	2.15	0.46
2:B:108:HIS:ND1	4:D:1:ILE:HD11	2.31	0.46
1:M:312:VAL:HG13	1:M:313:TYR:N	2.31	0.46
1:M:509:GLY:HA2	1:M:512:LEU:HD12	1.98	0.46
1:A:10:VAL:HG21	1:A:190:MET:HE1	1.98	0.45
1:M:83:ASP:O	1:M:87:HIS:HD2	1.99	0.45
1:M:10:VAL:HG13	8:M:601:FAD:C2A	2.46	0.45
1:M:440:VAL:O	1:M:490:ARG:NH1	2.50	0.45
1:M:521:SER:HA	1:M:563:LEU:HD21	1.98	0.45
2:N:59:MET:HG3	2:N:61:ILE:HG22	1.99	0.45
4:P:95:ALA:HB1	4:P:98:TRP:HB2	1.98	0.45
1:M:10:VAL:HG22	1:M:157:VAL:HG21	1.98	0.45
1:M:428:GLN:O	1:M:432:VAL:HG23	2.17	0.45
2:N:177:HIS:CE1	2:N:226:GLN:HB2	2.52	0.45
1:M:151:ARG:NH1	1:M:153:ASP:OD2	2.50	0.45
1:M:17:LEU:O	1:M:21:ILE:HG13	2.17	0.45
1:A:390:ARG:CD	1:A:395:SER:HB2	2.47	0.45
1:M:78:GLU:O	1:M:82:VAL:HG23	2.17	0.45
1:M:103:TRP:O	1:M:105:ARG:HG3	2.17	0.45
2:B:120:ILE:HD13	2:B:185:ASP:HB2	1.98	0.44
1:M:232:HIS:O	1:M:352:PRO:HA	2.17	0.44
1:M:211:ASP:O	1:M:215:MET:HG3	2.17	0.44
1:M:222:PRO:HB3	1:M:554:PHE:CZ	2.53	0.44
2:N:235:PHE:HA	4:P:7:ARG:NH2	2.31	0.44
2:N:241:LYS:H	2:N:242:PRO:HD2	1.83	0.44
1:A:398:GLU:HG2	1:A:402:PHE:CD1	2.48	0.44
4:D:2:ASN:ND2	4:D:3:PRO:HD2	2.32	0.44
2:N:144:GLN:NE2	2:N:219:ASP:HB2	2.33	0.44
2:N:229:VAL:O	2:N:233:LYS:HG3	2.18	0.44
4:D:9:ASP:O	4:D:12:VAL:HG23	2.17	0.44
1:M:356:TYR:CZ	1:M:379:GLU:HG3	2.52	0.44
1:M:35:ILE:CD1	1:M:155:HIS:HB3	2.48	0.44
1:M:324:LEU:HD23	1:M:328:LEU:HD12	1.99	0.44
4:P:28:ALA:N	4:P:29:PRO:HD2	2.32	0.44
4:P:59:GLN:HA	4:P:64:ARG:NH1	2.33	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:204:ASN:N	1:A:204:ASN:ND2	2.64	0.44
1:M:233:PRO:HG2	1:M:248:ARG:NH2	2.30	0.44
1:M:231:TYR:HA	1:M:353:THR:O	2.18	0.44
1:M:66:PHE:HZ	1:M:79:GLN:HB3	1.81	0.44
1:M:364:ASP:OD1	1:M:368:GLU:N	2.46	0.44
2:N:229:VAL:HG12	2:N:233:LYS:NZ	2.33	0.44
2:B:166:GLU:HA	2:B:166:GLU:OE1	2.18	0.43
2:B:235:PHE:CD1	4:D:11:PRO:HG2	2.53	0.43
3:C:76:LEU:HD13	4:D:32:ILE:HB	1.98	0.43
2:N:224:ILE:HD13	6:N:245:F3S:S3	2.58	0.43
1:A:146:PHE:HB2	1:A:149:ILE:HD12	1.99	0.43
1:A:174:ASN:ND2	1:A:174:ASN:C	2.71	0.43
2:N:175:LEU:O	2:N:178:ARG:HB3	2.18	0.43
1:M:197:ARG:HB2	1:M:208:VAL:O	2.19	0.43
3:O:120:THR:CG2	4:P:30:VAL:HG23	2.37	0.43
4:P:9:ASP:C	4:P:11:PRO:HD2	2.38	0.43
1:M:448:TRP:CH2	1:M:504:TYR:HB3	2.52	0.43
4:P:51:TYR:CD1	4:P:51:TYR:C	2.89	0.43
1:M:459:MET:O	1:M:463:CYS:HB2	2.18	0.43
1:M:552:LEU:HD11	1:M:566:SER:HB3	2.00	0.43
3:O:122:VAL:O	3:O:126:VAL:HG23	2.18	0.43
1:M:399:LEU:CD2	8:M:601:FAD:O2P	2.67	0.43
1:M:58:ASP:C	1:M:60:ASP:H	2.21	0.43
1:A:53:ALA:HB2	1:A:393:SER:OG	2.19	0.43
3:C:40:ILE:HA	3:C:43:ILE:HD12	2.00	0.43
1:M:479:LEU:HA	1:M:482:LEU:HD12	2.00	0.43
4:P:20:GLY:O	4:P:24:SER:HB3	2.19	0.43
1:A:230:GLN:NE2	1:A:390:ARG:HH21	2.17	0.43
1:M:11:GLY:O	1:M:16:GLY:HA3	2.19	0.43
1:A:230:GLN:HE22	1:A:390:ARG:HH21	1.66	0.42
1:M:485:ARG:O	1:M:489:VAL:HG23	2.19	0.42
2:B:160:GLN:OE1	2:B:160:GLN:HA	2.18	0.42
1:M:51:GLY:HA2	1:M:131:THR:HG21	2.01	0.42
1:A:365:GLN:CD	1:A:365:GLN:H	2.23	0.42
2:B:116:ILE:H	2:B:116:ILE:HG13	1.73	0.42
1:M:481:GLU:O	1:M:485:ARG:HG3	2.19	0.42
4:P:52:GLU:HG2	4:P:52:GLU:H	1.54	0.42
1:A:356:TYR:HE1	8:A:601:FAD:O3'	2.03	0.42
1:M:517:CYS:O	1:M:521:SER:OG	2.27	0.42
2:B:99:GLU:OE2	2:B:105:ASP:HB2	2.20	0.42
1:M:217:LEU:HD21	1:M:517:CYS:HG	1.83	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:62:PHE:CE2	1:M:87:HIS:HA	2.55	0.42
4:D:34:LEU:HD23	4:D:38:LEU:HD12	2.02	0.42
1:M:447:ASN:HB3	1:M:450:LYS:HG2	2.01	0.42
1:M:76:LEU:HG	1:M:529:ARG:HD3	2.01	0.42
3:O:120:THR:HG23	4:P:30:VAL:CG2	2.35	0.42
3:O:130:TRP:HB3	4:P:53:ARG:NH2	2.34	0.42
3:O:28:ARG:HD2	4:P:81:ARG:HH21	1.85	0.42
1:A:92:GLU:OE2	1:A:404:ARG:HD2	2.20	0.42
2:B:208:GLY:N	6:B:245:F3S:S1	2.92	0.42
1:M:356:TYR:CE2	1:M:390:ARG:HD3	2.54	0.42
1:M:11:GLY:CA	8:M:601:FAD:H1B	2.19	0.42
2:N:159:PRO:CG	2:N:207:VAL:HG21	2.49	0.42
1:A:244:THR:HG22	1:A:331:ILE:HG13	2.01	0.42
1:M:1:GLN:NE2	1:M:2:THR:N	2.68	0.42
8:M:601:FAD:H9	8:M:601:FAD:H1'1	1.63	0.42
2:N:229:VAL:HG12	2:N:233:LYS:HZ2	1.85	0.42
1:A:446:GLU:HG2	1:A:451:ILE:HD11	2.01	0.41
1:M:465:ILE:H	1:M:465:ILE:HG12	1.62	0.41
1:M:217:LEU:HD23	1:M:513:ASN:HB3	2.01	0.41
2:B:43:ILE:HD12	2:B:47:LEU:HD12	2.01	0.41
4:D:47:ASP:HB3	4:D:50:SER:OG	2.20	0.41
1:M:574:PRO:HA	1:M:575:PRO:HD3	1.92	0.41
1:A:342:ASP:HA	1:A:343:PRO:HD2	1.92	0.41
1:A:467:ARG:O	1:A:534:ARG:HA	2.20	0.41
1:A:182:GLN:OE1	1:A:184:ARG:NH2	2.50	0.41
4:P:113:ILE:HG22	4:P:114:GLY:N	2.35	0.41
4:D:48:ALA:HA	4:D:53:ARG:HD3	2.02	0.41
1:M:20:ALA:HB2	1:M:34:LEU:HD13	2.03	0.41
1:M:5:ALA:HB1	1:M:31:LYS:O	2.21	0.41
1:A:311:VAL:HG21	1:A:349:PRO:HB3	2.01	0.41
2:B:242:PRO:HB3	4:P:94:PRO:N	2.36	0.41
1:M:342:ASP:HB3	1:M:345:LYS:HB3	2.02	0.41
1:M:89:CYS:HB2	1:M:90:PRO:HD3	2.03	0.41
4:P:0:MET:SD	4:P:1:ILE:HD13	2.61	0.41
1:A:186:ASN:HB3	1:A:417:ALA:HB2	2.02	0.41
3:C:128:LEU:O	4:D:45:PRO:HG2	2.21	0.41
1:M:443:ASP:HA	1:M:490:ARG:HD3	2.03	0.41
4:D:9:ASP:C	4:D:11:PRO:HD2	2.40	0.41
1:M:485:ARG:HA	1:M:488:ARG:HG2	2.03	0.41
1:A:356:TYR:CZ	1:A:379:GLU:HG3	2.56	0.41
1:M:146:PHE:HB3	1:M:148:GLN:OE1	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:N:116:ILE:CG2	2:N:176:ALA:CB	2.89	0.41
1:A:386:HIS:ND1	1:A:390:ARG:HG3	2.36	0.41
1:A:448:TRP:CH2	1:A:504:TYR:HB3	2.56	0.41
1:A:556:ASP:OD2	1:A:562:ARG:HD2	2.20	0.41
1:M:46:VAL:O	1:M:133:PHE:N	2.54	0.41
2:N:96:PHE:HB3	2:N:104:VAL:HB	2.03	0.41
3:C:19:LEU:HD23	3:C:22:TYR:CE2	2.56	0.40
1:M:131:THR:O	1:M:135:MET:HG3	2.21	0.40
1:M:145:GLN:HB3	2:N:119:TYR:CE2	2.56	0.40
1:M:187:ALA:HB2	1:M:414:ALA:HB2	2.03	0.40
1:M:116:PHE:HE2	1:M:245:GLU:HB3	1.82	0.40
1:M:554:PHE:HB2	1:M:562:ARG:HB2	2.03	0.40
1:A:346:GLU:CB	1:A:347:PRO:CD	2.99	0.40
1:A:240:GLY:HA2	1:A:353:THR:HG21	2.03	0.40
2:N:220:PRO:O	2:N:223:ALA:HB3	2.22	0.40
2:N:117:LYS:O	2:N:191:ARG:NH2	2.46	0.40
1:M:145:GLN:HB3	2:N:119:TYR:CZ	2.57	0.40
2:N:73:PRO:HB2	2:N:153:LEU:HD22	2.03	0.40
1:A:76:LEU:O	1:A:550:HIS:HE1	2.04	0.40
2:B:93:LEU:HG	2:B:156:ALA:HB2	2.03	0.40
1:M:76:LEU:HD11	1:M:525:ARG:HH22	1.85	0.40
4:P:86:MET:HE2	4:P:91:ILE:HG21	2.04	0.40
2:B:207:VAL:HG23	6:B:245:F3S:S2	2.62	0.40
2:B:108:HIS:CE1	4:D:1:ILE:HD11	2.57	0.40
4:D:55:LEU:CG	4:D:59:GLN:NE2	2.69	0.40
1:M:295:TRP:O	1:M:299:ARG:HG2	2.22	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	575/577 (100%)	551 (96%)	24 (4%)	0	100	100
1	M	575/577 (100%)	545 (95%)	28 (5%)	2 (0%)	44	77
2	B	241/243 (99%)	229 (95%)	10 (4%)	2 (1%)	22	58
2	N	241/243 (99%)	225 (93%)	13 (5%)	3 (1%)	15	47
3	C	128/130 (98%)	125 (98%)	3 (2%)	0	100	100
3	O	128/130 (98%)	123 (96%)	5 (4%)	0	100	100
4	D	117/119 (98%)	103 (88%)	14 (12%)	0	100	100
4	P	117/119 (98%)	104 (89%)	13 (11%)	0	100	100
All	All	2122/2138 (99%)	2005 (94%)	110 (5%)	7 (0%)	44	77

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	N	56	SER
2	B	56	SER
1	M	444	GLY
2	N	101	ASP
2	B	82	ARG
1	M	328	LEU
2	N	241	LYS

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	460/460 (100%)	439 (95%)	21 (5%)	31	66
1	M	460/460 (100%)	418 (91%)	42 (9%)	11	36
2	B	205/205 (100%)	196 (96%)	9 (4%)	33	67
2	N	205/205 (100%)	193 (94%)	12 (6%)	23	56
3	C	111/111 (100%)	105 (95%)	6 (5%)	26	59
3	O	111/111 (100%)	106 (96%)	5 (4%)	32	66
4	D	97/97 (100%)	86 (89%)	11 (11%)	7	25

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
4	P	97/97 (100%)	85 (88%)	12 (12%)	5	20
All	All	1746/1746 (100%)	1628 (93%)	118 (7%)	18	50

All (118) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	ARG
1	A	31	LYS
1	A	41	MET
1	A	42	ARG
1	A	49	GLU
1	A	74	ASP
1	A	93	MET
1	A	122	GLU
1	A	174	ASN
1	A	182	GLN
1	A	197	ARG
1	A	200	ARG
1	A	204	ASN
1	A	276	GLU
1	A	317	ARG
1	A	321	GLU
1	A	325	HIS
1	A	327	ARG
1	A	422	GLU
1	A	459	MET
1	A	558	ASP
2	B	19	THR
2	B	65	CYS
2	B	128	ASP
2	B	178	ARG
2	B	206	PHE
2	B	210	CYS
2	B	212	GLU
2	B	225	GLN
2	B	241	LYS
3	C	1	THR
3	C	12	THR
3	C	26	MET
3	C	39	SER
3	C	50	LYS
3	C	99	LYS

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Mol	Chain	Res	Type
4	D	1	ILE
4	D	2	ASN
4	D	6	LYS
4	D	9	ASP
4	D	12	VAL
4	D	24	SER
4	D	39	LEU
4	D	67	LEU
4	D	86	MET
4	D	91	ILE
4	D	117	THR
1	M	1	GLN
1	M	2	THR
1	M	17	LEU
1	M	18	ARG
1	M	76	LEU
1	M	93	MET
1	M	116	PHE
1	M	143	SER
1	M	156	PHE
1	M	174	ASN
1	M	202	ASN
1	M	209	THR
1	M	226	MET
1	M	227	GLU
1	M	243	MET
1	M	244	THR
1	M	283	GLU
1	M	288	ASP
1	M	295	TRP
1	M	299	ARG
1	M	312	VAL
1	M	321	GLU
1	M	327	ARG
1	M	330	PHE
1	M	344	VAL
1	M	353	THR
1	M	369	THR
1	M	377	VAL
1	M	413	ARG
1	M	426	GLU
1	M	428	GLN

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Mol	Chain	Res	Type
1	M	447	ASN
1	M	457	LEU
1	M	461	GLU
1	M	465	ILE
1	M	470	GLU
1	M	494	THR
1	M	537	GLU
1	M	540	THR
1	M	543	ASP
1	M	558	ASP
1	M	571	THR
2	N	2	GLU
2	N	17	VAL
2	N	85	THR
2	N	123	ASN
2	N	150	ASN
2	N	189	LYS
2	N	192	MET
2	N	203	SER
2	N	206	PHE
2	N	220	PRO
2	N	232	SER
2	N	240	LEU
3	O	5	LYS
3	O	19	LEU
3	O	54	GLU
3	O	102	LYS
3	O	130	TRP
4	P	1	ILE
4	P	2	ASN
4	P	6	LYS
4	P	12	VAL
4	P	24	SER
4	P	30	VAL
4	P	39	LEU
4	P	64	ARG
4	P	67	LEU
4	P	86	MET
4	P	91	ILE
4	P	117	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (41) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	137	HIS
1	A	141	GLN
1	A	150	GLN
1	A	174	ASN
1	A	230	GLN
1	A	232	HIS
1	A	279	ASN
1	A	409	GLN
1	A	434	GLN
1	A	442	GLN
2	B	134	GLN
2	B	225	GLN
3	C	72	ASN
4	D	4	ASN
4	D	59	GLN
4	D	80	HIS
4	D	92	HIS
1	M	1	GLN
1	M	57	GLN
1	M	59	HIS
1	M	65	HIS
1	M	87	HIS
1	M	95	GLN
1	M	134	HIS
1	M	166	HIS
1	M	174	ASN
1	M	182	GLN
1	M	232	HIS
1	M	428	GLN
1	M	447	ASN
1	M	483	GLN
2	N	95	ASN
2	N	123	ASN
2	N	138	GLN
2	N	177	HIS
2	N	225	GLN
2	N	226	GLN
3	O	51	ASN
3	O	72	ASN
4	P	59	GLN
4	P	80	HIS

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 ⓘ

10 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$
9	GUA	A	577	-	2,8,8	0.19	0	1,9,9	0.58	0
8	FAD	A	601	-	51,58,58	0.88	1 (1%)	54,89,89	2.15	11 (20%)
5	FES	B	244	2	0,4,4	0.00	-	0,4,4	0.00	-
6	F3S	B	245	2	0,9,9	0.00	-	0,15,15	0.00	-
7	SF4	B	246	2	0,12,12	0.00	-	0,24,24	0.00	-
9	GUA	M	577	-	2,8,8	0.28	0	1,9,9	0.20	0
8	FAD	M	601	-	51,58,58	1.21	4 (7%)	54,89,89	2.13	8 (14%)
5	FES	N	244	2	0,4,4	0.00	-	0,4,4	0.00	-
6	F3S	N	245	2	0,9,9	0.00	-	0,15,15	0.00	-
7	SF4	N	246	2	0,12,12	0.00	-	0,24,24	0.00	-

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
9	GUA	A	577	-	-	0/2/6/6	0/0/0/0
8	FAD	A	601	-	-	0/28/50/50	0/6/6/6
5	FES	B	244	2	-	0/0/4/4	0/1/1/1
6	F3S	B	245	2	-	0/0/24/24	0/0/3/3
7	SF4	B	246	2	-	0/0/48/48	0/6/5/5
9	GUA	M	577	-	-	0/2/6/6	0/0/0/0
8	FAD	M	601	-	-	0/28/50/50	0/6/6/6
5	FES	N	244	2	-	0/0/4/4	0/1/1/1
6	F3S	N	245	2	-	0/0/24/24	0/0/3/3
7	SF4	N	246	2	-	0/0/48/48	0/6/5/5

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	M	601	FAD	O2B-C2B	-3.38	1.35	1.43
8	M	601	FAD	C9A-N10	2.01	1.41	1.38
8	A	601	FAD	O4B-C1B	2.03	1.44	1.41
8	M	601	FAD	C10-N1	2.28	1.36	1.33
8	M	601	FAD	O3B-C3B	4.62	1.53	1.43

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	A	601	FAD	N3A-C2A-N1A	-10.10	120.06	128.86
8	M	601	FAD	N3A-C2A-N1A	-9.48	120.60	128.86
8	M	601	FAD	C4X-C10-N10	-2.78	118.59	120.52
8	M	601	FAD	C1B-N9A-C4A	-2.60	122.15	126.64
8	A	601	FAD	C4X-C4-N3	-2.56	119.84	123.48
8	A	601	FAD	C1B-N9A-C4A	-2.48	122.35	126.64
8	M	601	FAD	C4X-C4-N3	-2.41	120.06	123.48
8	A	601	FAD	C4A-C5A-N7A	-2.30	107.19	109.41
8	A	601	FAD	C4-C4X-C10	-2.18	118.20	119.96
8	A	601	FAD	C1'-C2'-C3'	2.33	116.47	109.82
8	A	601	FAD	C4-C4X-N5	2.54	121.46	118.68
8	A	601	FAD	C5X-C9A-N10	2.59	119.58	117.66
8	M	601	FAD	C4X-N5-C5X	3.83	120.81	116.76
8	A	601	FAD	C4X-N5-C5X	3.94	120.93	116.76
8	M	601	FAD	C5X-C9A-N10	3.96	120.60	117.66
8	A	601	FAD	C1'-N10-C9A	4.17	122.17	118.35
8	M	601	FAD	C1'-N10-C10	5.77	124.42	118.50
8	M	601	FAD	C4-N3-C2	5.86	120.28	115.16
8	A	601	FAD	C4-N3-C2	6.45	120.80	115.16

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

8 monomers are involved in 41 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
9	A	577	GUA	1	0
8	A	601	FAD	8	0
6	B	245	F3S	6	0
7	B	246	SF4	2	0
9	M	577	GUA	2	0
8	M	601	FAD	13	0
6	N	245	F3S	3	0
7	N	246	SF4	6	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

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	577/577 (100%)	-0.04	0 100 100	44, 61, 96, 134	0
1	M	577/577 (100%)	0.50	44 (7%) 15 5	72, 135, 169, 186	0
2	B	243/243 (100%)	-0.08	0 100 100	46, 67, 104, 154	0
2	N	243/243 (100%)	0.17	8 (3%) 47 23	65, 106, 151, 167	0
3	C	130/130 (100%)	-0.08	1 (0%) 86 69	58, 80, 106, 155	0
3	O	130/130 (100%)	0.04	1 (0%) 86 69	72, 102, 147, 197	0
4	D	119/119 (100%)	-0.05	1 (0%) 86 69	59, 83, 124, 158	0
4	P	119/119 (100%)	0.09	2 (1%) 70 47	72, 91, 178, 254	0
All	All	2138/2138 (100%)	0.13	57 (2%) 55 29	44, 87, 158, 254	0

All (57) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	M	169	GLY	5.2
1	M	8	ALA	5.2
1	M	411	THR	4.7
3	O	1	THR	4.6
1	M	172	ALA	4.6
1	M	188	VAL	4.5
1	M	373	GLY	4.5
4	D	118	ILE	4.4
2	N	1	ALA	4.4
1	M	376	ALA	4.3
1	M	6	ASP	4.1
1	M	219	HIS	4.0
1	M	159	ASP	3.9
1	M	77	CYS	3.8
1	M	250	GLU	3.6
1	M	217	LEU	3.3

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Mol	Chain	Res	Type	RSRZ
1	M	494	THR	3.2
1	M	171	VAL	3.1
1	M	571	THR	3.1
1	M	413	ARG	3.0
2	N	238	ALA	2.9
1	M	160	ILE	2.8
4	P	9	ASP	2.8
1	M	215	MET	2.8
1	M	33	ALA	2.8
1	M	189	VAL	2.7
1	M	7	LEU	2.7
1	M	223	LEU	2.6
1	M	158	LEU	2.6
2	N	17	VAL	2.5
1	M	10	VAL	2.5
3	C	130	TRP	2.4
1	M	498	PHE	2.4
1	M	32	ILE	2.4
1	M	187	ALA	2.4
2	N	2	GLU	2.4
1	M	553	ALA	2.3
2	N	237	ILE	2.3
1	M	486	PHE	2.2
1	M	170	LEU	2.2
1	M	0	MET	2.2
1	M	573	LEU	2.2
1	M	11	GLY	2.2
2	N	126	THR	2.2
1	M	415	ALA	2.1
1	M	402	PHE	2.1
2	N	131	THR	2.1
1	M	88	HIS	2.1
1	M	125	TRP	2.1
1	M	444	GLY	2.1
1	M	489	VAL	2.1
4	P	88	ASP	2.1
2	N	8	ILE	2.1
1	M	161	LEU	2.1
1	M	165	GLY	2.0
1	M	451	ILE	2.0
1	M	63	GLU	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
6	F3S	B	245	7/7	0.97	0.25	1.36	75,77,79,80	0
9	GUA	A	577	9/9	0.82	0.28	1.23	61,62,65,65	0
7	SF4	B	246	8/8	0.98	0.26	0.78	71,74,76,76	0
6	F3S	N	245	7/7	0.97	0.23	0.68	93,95,96,96	0
5	FES	B	244	4/4	0.99	0.23	0.25	50,50,54,55	0
7	SF4	N	246	8/8	0.94	0.20	-0.19	88,88,89,89	0
5	FES	N	244	4/4	0.99	0.20	-0.32	72,73,73,74	0
9	GUA	M	577	9/9	0.78	0.22	-0.47	86,87,88,88	0
8	FAD	M	601	53/53	0.89	0.21	-0.67	65,84,96,96	0
8	FAD	A	601	53/53	0.98	0.20	-0.70	35,43,47,48	0

6.5 Other polymers [i](#)

There are no such residues in this entry.