



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 26, 2014 – 08:40 PM GMT

PDB ID : 2F5Z
Title : Crystal Structure of Human Dihydrolipoamide Dehydrogenase (E3) Complexed to the E3-Binding Domain of Human E3-Binding Protein
Authors : Brautigam, C.A.; Chuang, J.L.; Wynn, R.M.; Tomchick, D.R.; Machius, M.; Chuang, D.T.
Deposited on : 2005-11-28
Resolution : 2.18 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

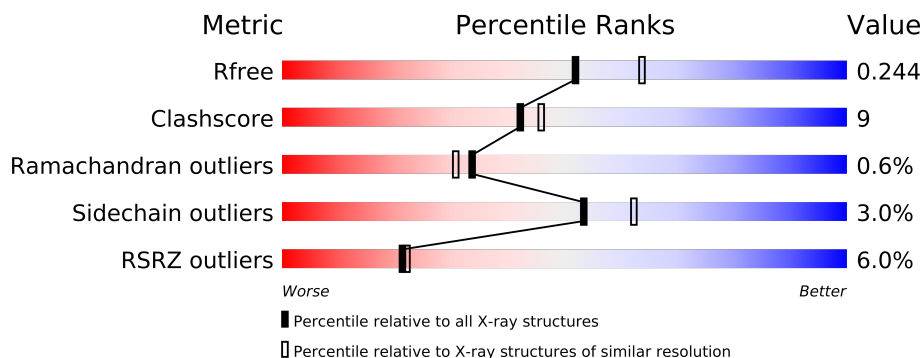
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.18 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}	66092	3841 (2.20-2.16)
Clashscore	79885	4835 (2.20-2.16)
Ramachandran outliers	78287	4740 (2.20-2.16)
Sidechain outliers	78261	4741 (2.20-2.16)
RSRZ outliers	66119	3842 (2.20-2.16)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	474	
1	B	474	
1	C	474	
1	D	474	
1	E	474	
1	F	474	
1	G	474	
1	H	474	
1	I	474	
1	J	474	
2	K	64	
2	L	64	
2	M	64	
2	N	64	

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Mol	Chain	Length	Quality of chain
2	O	64	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
3	SO4	A	2483	-	X
3	SO4	B	2487	-	X
3	SO4	B	2488	-	X
3	SO4	C	2490	-	X
3	SO4	C	2491	-	X
3	SO4	D	2496	-	X
3	SO4	E	2499	-	X
3	SO4	E	2500	-	X
3	SO4	F	2503	-	X
3	SO4	F	2504	-	X
3	SO4	G	2508	-	X
3	SO4	H	2511	-	X
3	SO4	H	2512	-	X
3	SO4	I	2515	-	X
3	SO4	J	2519	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 40176 atoms, of which 0 are hydrogen and 0 are deuterium.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Dihydrolipoyl dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	472	Total	C	N	O	S	0	0	0
			3505	2208	606	672	19			
1	B	471	Total	C	N	O	S	0	0	0
			3496	2203	604	670	19			
1	C	472	Total	C	N	O	S	0	0	0
			3505	2208	606	672	19			
1	D	471	Total	C	N	O	S	0	0	0
			3496	2203	604	670	19			
1	E	472	Total	C	N	O	S	0	0	0
			3505	2208	606	672	19			
1	F	471	Total	C	N	O	S	0	0	0
			3496	2203	604	670	19			
1	G	472	Total	C	N	O	S	0	0	0
			3505	2208	606	672	19			
1	H	472	Total	C	N	O	S	0	0	0
			3505	2208	606	672	19			
1	I	470	Total	C	N	O	S	0	0	0
			3496	2204	604	669	19			
1	J	472	Total	C	N	O	S	0	0	0
			3505	2208	606	672	19			

- Molecule 2 is a protein called Pyruvate dehydrogenase protein X component.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	K	43	Total	C	N	O	0	0	0
			301	186	57	58			
2	L	43	Total	C	N	O	0	43	0
			590	364	112	114			
2	M	43	Total	C	N	O	0	43	0
			590	364	112	114			
2	N	43	Total	C	N	O	0	43	0
			590	364	112	114			

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	O	43	Total	C	N	O	0	43	0
			590	364	112	114			

There are 40 discrepancies between the modelled and reference sequences:

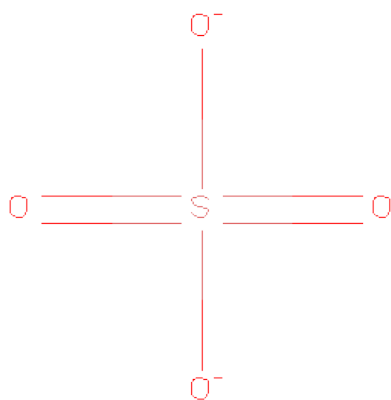
Chain	Residue	Modelled	Actual	Comment	Reference
K	120	GLY	LYS	ENGINEERED	UNP O00330
K	176	LEU	THR	ENGINEERED	UNP O00330
K	178	HIS	-	EXPRESSION TAG	UNP O00330
K	179	HIS	-	EXPRESSION TAG	UNP O00330
K	180	HIS	-	EXPRESSION TAG	UNP O00330
K	181	HIS	-	EXPRESSION TAG	UNP O00330
K	182	HIS	-	EXPRESSION TAG	UNP O00330
K	183	HIS	-	EXPRESSION TAG	UNP O00330
L	120	GLY	LYS	ENGINEERED	UNP O00330
L	176	LEU	THR	ENGINEERED	UNP O00330
L	178	HIS	-	EXPRESSION TAG	UNP O00330
L	179	HIS	-	EXPRESSION TAG	UNP O00330
L	180	HIS	-	EXPRESSION TAG	UNP O00330
L	181	HIS	-	EXPRESSION TAG	UNP O00330
L	182	HIS	-	EXPRESSION TAG	UNP O00330
L	183	HIS	-	EXPRESSION TAG	UNP O00330
M	120	GLY	LYS	ENGINEERED	UNP O00330
M	176	LEU	THR	ENGINEERED	UNP O00330
M	178	HIS	-	EXPRESSION TAG	UNP O00330
M	179	HIS	-	EXPRESSION TAG	UNP O00330
M	180	HIS	-	EXPRESSION TAG	UNP O00330
M	181	HIS	-	EXPRESSION TAG	UNP O00330
M	182	HIS	-	EXPRESSION TAG	UNP O00330
M	183	HIS	-	EXPRESSION TAG	UNP O00330
N	120	GLY	LYS	ENGINEERED	UNP O00330
N	176	LEU	THR	ENGINEERED	UNP O00330
N	178	HIS	-	EXPRESSION TAG	UNP O00330
N	179	HIS	-	EXPRESSION TAG	UNP O00330
N	180	HIS	-	EXPRESSION TAG	UNP O00330
N	181	HIS	-	EXPRESSION TAG	UNP O00330
N	182	HIS	-	EXPRESSION TAG	UNP O00330
N	183	HIS	-	EXPRESSION TAG	UNP O00330
O	120	GLY	LYS	ENGINEERED	UNP O00330
O	176	LEU	THR	ENGINEERED	UNP O00330
O	178	HIS	-	EXPRESSION TAG	UNP O00330
O	179	HIS	-	EXPRESSION TAG	UNP O00330

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Chain	Residue	Modelled	Actual	Comment	Reference
O	180	HIS	-	EXPRESSION TAG	UNP O00330
O	181	HIS	-	EXPRESSION TAG	UNP O00330
O	182	HIS	-	EXPRESSION TAG	UNP O00330
O	183	HIS	-	EXPRESSION TAG	UNP O00330

- Molecule 3 is SULFATE ION (three-letter code: SO₄) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		

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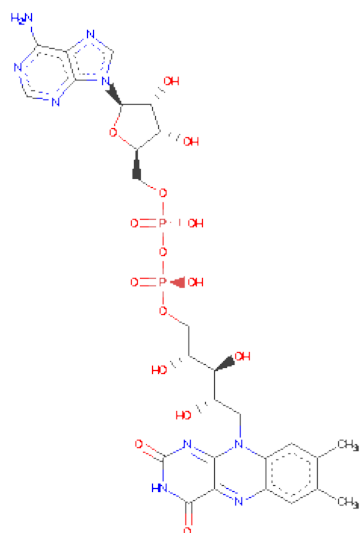
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	D	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	E	1	Total	O	S	0	0
			5	4	1		
3	E	1	Total	O	S	0	0
			5	4	1		
3	E	1	Total	O	S	0	0
			5	4	1		
3	E	1	Total	O	S	0	0
			5	4	1		
3	F	1	Total	O	S	0	0
			5	4	1		
3	F	1	Total	O	S	0	0
			5	4	1		
3	F	1	Total	O	S	0	0
			5	4	1		
3	F	1	Total	O	S	0	0
			5	4	1		
3	G	1	Total	O	S	0	0
			5	4	1		
3	G	1	Total	O	S	0	0
			5	4	1		
3	G	1	Total	O	S	0	0
			5	4	1		
3	G	1	Total	O	S	0	0
			5	4	1		
3	H	1	Total	O	S	0	0
			5	4	1		
3	H	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	H	1	Total	O	S	0	0
			5	4	1		
3	H	1	Total	O	S	0	0
			5	4	1		
3	I	1	Total	O	S	0	0
			5	4	1		
3	I	1	Total	O	S	0	0
			5	4	1		
3	I	1	Total	O	S	0	0
			5	4	1		
3	J	1	Total	O	S	0	0
			5	4	1		
3	J	1	Total	O	S	0	0
			5	4	1		
3	J	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula: $C_{27}H_{33}N_9O_{15}P_2$).



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

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	C	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
4	D	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
4	E	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
4	F	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
4	G	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
4	H	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
4	I	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
4	J	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	183	Total	O	0	0
			183	183		
5	B	101	Total	O	0	0
			101	101		
5	C	172	Total	O	0	0
			172	172		
5	D	243	Total	O	0	0
			243	243		
5	E	150	Total	O	0	0
			150	150		
5	F	217	Total	O	0	0
			217	217		
5	G	188	Total	O	0	0
			188	188		
5	H	315	Total	O	0	0
			315	315		
5	I	73	Total	O	0	0
			73	73		
5	J	132	Total	O	0	0
			132	132		
5	N	1	Total	O	0	0
			1	1		

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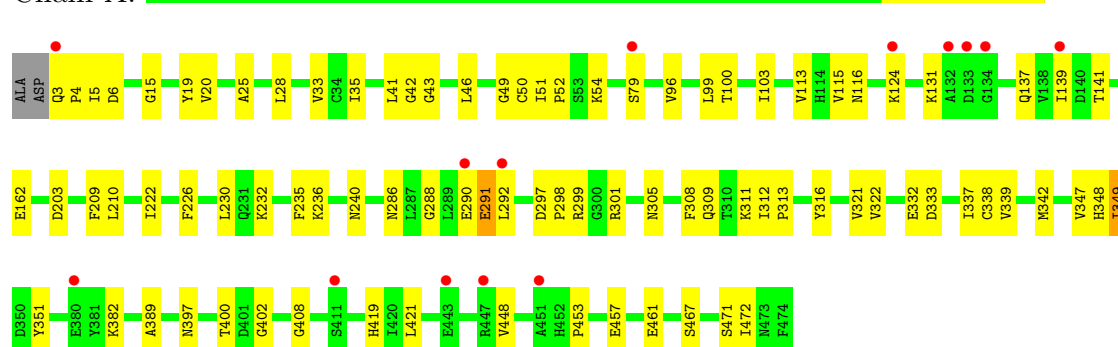
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	O	1	Total	O	0	0
			1	1		

3 Residue-property plots

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

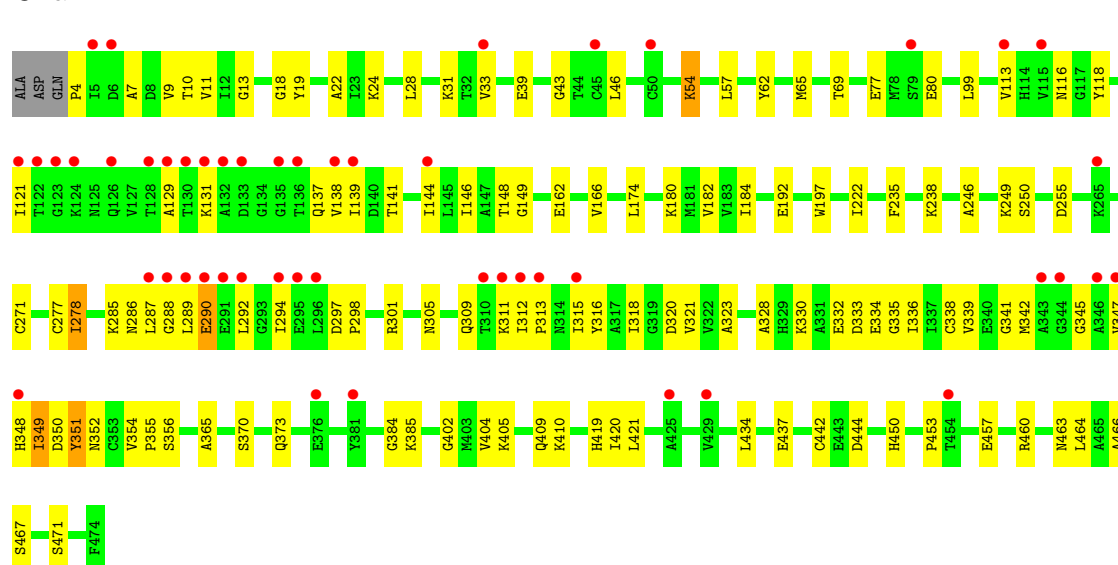
• Molecule 1: Dihydrolipoyl dehydrogenase

Chain A:



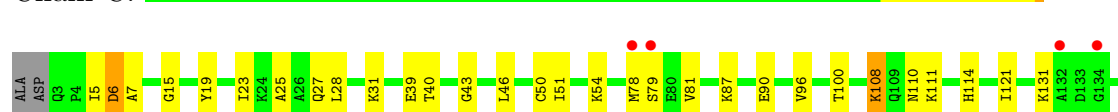
• Molecule 1: Dihydrolipoyl dehydrogenase

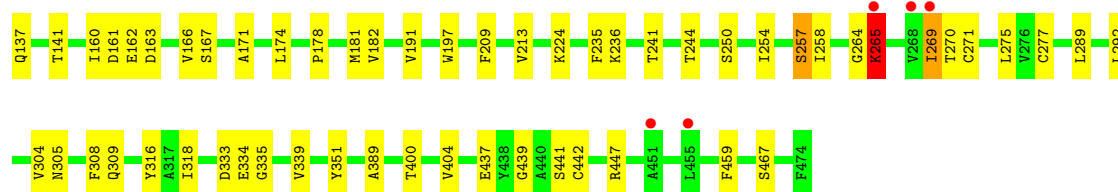
Chain B:



• Molecule 1: Dihydrolipoyl dehydrogenase

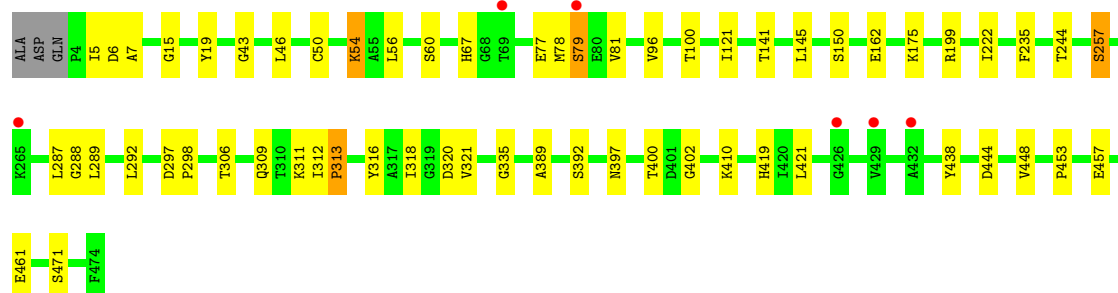
Chain C:





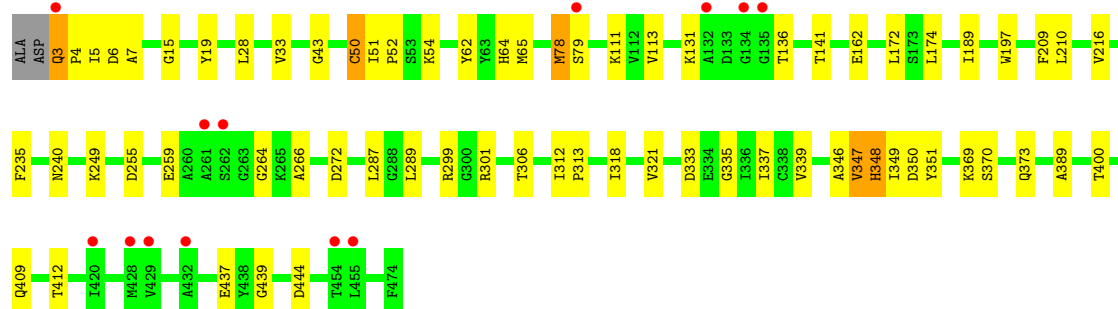
• Molecule 1: Dihydrolipoyl dehydrogenase

Chain D:



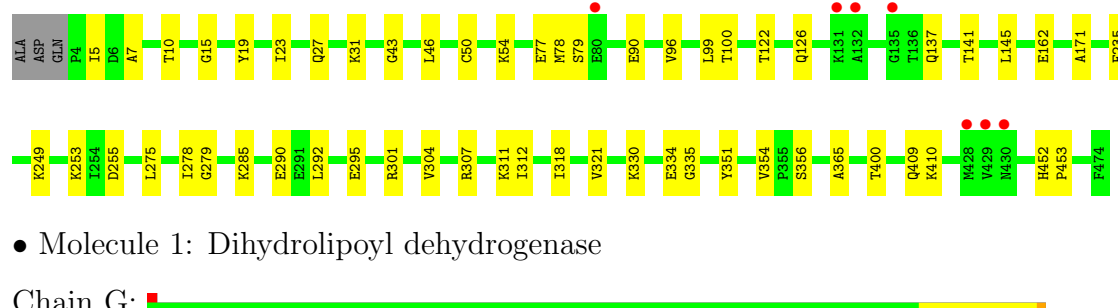
• Molecule 1: Dihydrolipoyl dehydrogenase

Chain E:



• Molecule 1: Dihydrolipoyl dehydrogenase

Chain F:



• Molecule 1: Dihydrolipoyl dehydrogenase

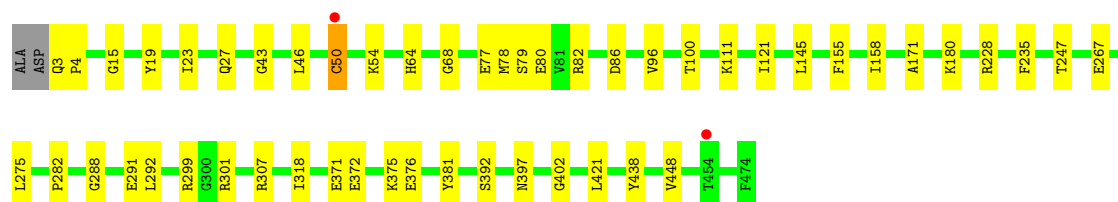
Chain G:





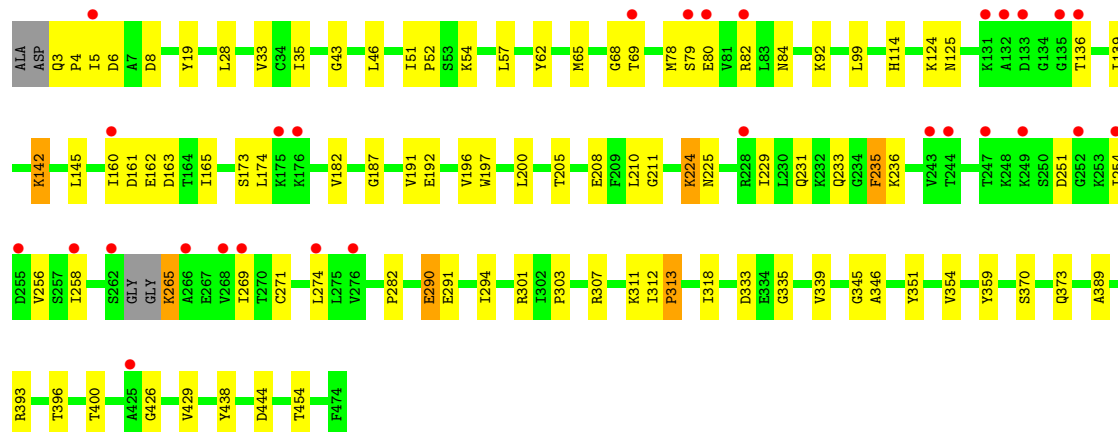
• Molecule 1: Dihydrolipoyl dehydrogenase

Chain H:



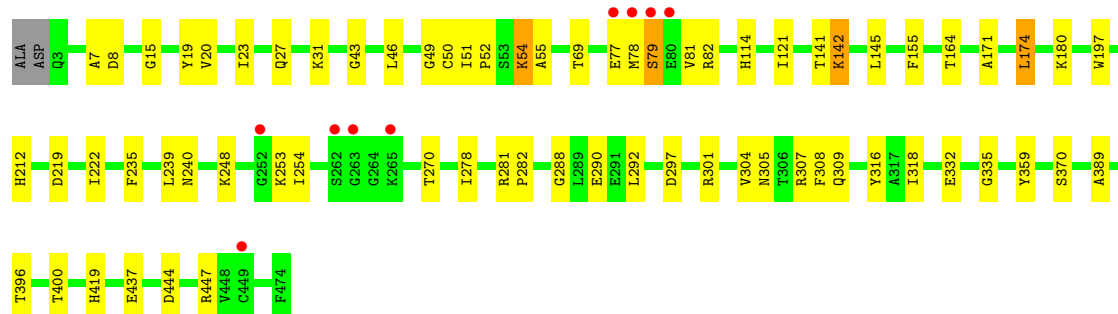
• Molecule 1: Dihydrolipoyl dehydrogenase

Chain I:



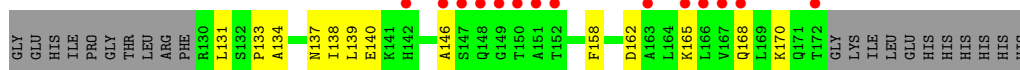
• Molecule 1: Dihydrolipoyl dehydrogenase

Chain J:



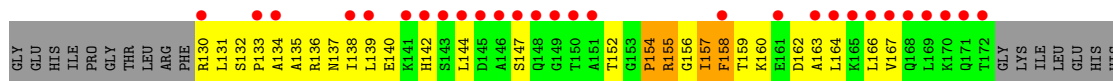
• Molecule 2: Pyruvate dehydrogenase protein X component

Chain K:



- Molecule 2: Pyruvate dehydrogenase protein X component

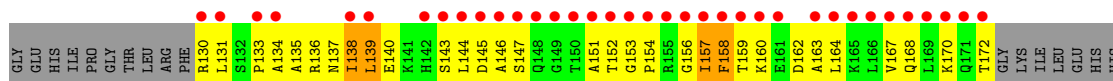
Chain L:



HIS
HIS
HIS

- Molecule 2: Pyruvate dehydrogenase protein X component

Chain M:



HIS
HIS
HIS

- Molecule 2: Pyruvate dehydrogenase protein X component

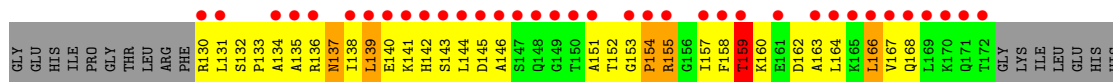
Chain N:



HIS
HIS
HIS

- Molecule 2: Pyruvate dehydrogenase protein X component

Chain O:



HIS
HIS
HIS

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	171.15Å 187.73Å 224.39Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	37.40 – 2.18 46.92 – 2.18	Depositor EDS
% Data completeness (in resolution range)	98.1 (37.40-2.18) 98.3 (46.92-2.18)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.79 (at 2.18Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.205 , 0.242 0.209 , 0.244	Depositor DCC
R_{free} test set	4517 reflections (1.23%)	DCC
Wilson B-factor (Å ²)	35.9	Xtriage
Anisotropy	0.311	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 36.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 367467 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	40176	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.16% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, 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	0.46	0/3560	0.70	1/4809 (0.0%)
1	B	0.42	0/3551	0.67	0/4796
1	C	0.52	0/3560	0.74	1/4809 (0.0%)
1	D	0.55	0/3551	0.75	1/4796 (0.0%)
1	E	0.51	0/3560	0.70	1/4809 (0.0%)
1	F	0.55	0/3551	0.74	3/4796 (0.1%)
1	G	0.52	0/3560	0.73	2/4809 (0.0%)
1	H	0.62	0/3560	0.77	1/4809 (0.0%)
1	I	0.46	0/3550	0.69	0/4796
1	J	0.50	0/3560	0.72	1/4809 (0.0%)
2	K	0.37	0/304	0.70	0/411
2	L	0.31	0/596	0.62	0/808
2	M	0.34	0/596	0.65	0/808
2	N	0.35	0/596	0.70	0/808
2	O	0.32	0/596	0.63	0/808
All	All	0.50	0/38251	0.72	11/51681 (0.0%)

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	50	CYS	CA-CB-SG	7.22	127.00	114.00
1	F	50	CYS	CA-CB-SG	6.26	125.27	114.00
1	J	50	CYS	CA-CB-SG	6.15	125.07	114.00
1	G	50	CYS	CA-CB-SG	6.04	124.86	114.00
1	E	50	CYS	CA-CB-SG	5.96	124.72	114.00
1	A	50	CYS	CA-CB-SG	5.91	124.64	114.00
1	D	50	CYS	CA-CB-SG	5.72	124.31	114.00
1	G	400	THR	N-CA-C	5.54	125.97	111.00
1	F	400	THR	N-CA-C	5.32	125.36	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	50	CYS	CA-CB-SG	5.18	123.32	114.00
1	F	31	LYS	N-CA-C	-5.02	97.45	111.00

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3505	0	3556	62	0
1	B	3496	0	3549	100	0
1	C	3505	0	3556	55	0
1	D	3496	0	3549	41	0
1	E	3505	0	3556	58	0
1	F	3496	0	3549	35	0
1	G	3505	0	3556	43	0
1	H	3505	0	3556	39	0
1	I	3496	0	3549	71	0
1	J	3505	0	3556	55	0
2	K	301	0	289	15	0
2	L	590	0	546	37	0
2	M	590	0	546	35	0
2	N	590	0	546	44	0
2	O	590	0	546	55	0
3	A	20	0	0	2	0
3	B	20	0	0	4	0
3	C	20	0	0	1	0
3	D	20	0	0	0	0
3	E	20	0	0	3	0
3	F	20	0	0	2	0
3	G	20	0	0	1	0
3	H	20	0	0	4	0
3	I	15	0	0	3	0
3	J	20	0	0	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	53	0	31	2	0
4	B	53	0	31	2	0
4	C	53	0	31	3	0
4	D	53	0	31	2	0
4	E	53	0	31	3	0
4	F	53	0	31	4	0
4	G	53	0	31	2	0
4	H	53	0	31	5	0
4	I	53	0	31	1	0
4	J	53	0	31	3	0
5	A	183	0	0	0	0
5	B	101	0	0	3	0
5	C	172	0	0	2	0
5	D	243	0	0	1	0
5	E	150	0	0	2	0
5	F	217	0	0	4	0
5	G	188	0	0	1	0
5	H	315	0	0	5	0
5	I	73	0	0	3	0
5	J	132	0	0	0	0
5	N	1	0	0	1	0
5	O	1	0	0	0	0
All	All	40176	0	38315	692	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 9.

All (692) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:I:290:GLU:HB3	3:I:2516:SO4:O4	1.54	1.07
1:A:461:GLU:HG2	1:A:471:SER:HB2	1.37	1.02
1:G:337:ILE:HG12	1:G:349:ILE:HD11	1.37	1.02
1:B:144:ILE:HB	1:B:315:ILE:HD13	1.39	1.01
2:O:158[B]:PHE:O	2:O:162[B]:ASP:HB2	1.71	0.90
1:A:308:PHE:HE2	1:A:337:ILE:HD11	1.33	0.89
2:N:142[B]:HIS:HB3	2:N:144[B]:LEU:HD13	1.54	0.89
2:O:157[B]:ILE:HG22	2:O:159[B]:THR:H	1.38	0.89
1:I:80:GLU:HG2	1:J:79:SER:HB3	1.58	0.85
1:B:370:SER:H	1:B:373:GLN:NE2	1.75	0.85
1:H:307:ARG:NH1	5:H:2731:HOH:O	2.08	0.85
2:O:154[B]:PRO:HG2	2:O:157[B]:ILE:HD13	1.59	0.84
1:E:78:MET:HG2	1:F:78:MET:HG2	1.59	0.84

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:67:HIS:HA	1:D:81:VAL:HG11	1.59	0.82
2:N:132[B]:SER:HB3	2:N:157[B]:ILE:HD11	1.63	0.81
2:O:159[B]:THR:HG23	2:O:162[B]:ASP:H	1.45	0.80
2:M:138[B]:ILE:HD11	2:M:163[B]:ALA:HB3	1.64	0.79
1:B:288:GLY:O	1:B:292:LEU:HD13	1.82	0.79
1:E:79:SER:HB2	1:F:79:SER:OG	1.83	0.79
1:F:307:ARG:NH1	5:F:2714:HOH:O	2.16	0.78
1:E:370:SER:H	1:E:373:GLN:NE2	1.82	0.78
1:B:11:VAL:HG11	1:B:22:ALA:HB2	1.64	0.78
2:N:130[B]:ARG:O	2:N:157[B]:ILE:HD12	1.84	0.78
1:A:308:PHE:CE2	1:A:337:ILE:HD11	2.19	0.77
1:I:79:SER:HB2	1:J:79:SER:OG	1.85	0.76
1:A:5:ILE:HG22	1:A:137:GLN:HE21	1.51	0.76
1:J:447:ARG:NE	2:O:137[A]:ASN:ND2	2.34	0.76
2:N:166[A]:LEU:HD13	2:N:166[A]:LEU:O	1.86	0.76
1:B:250:SER:HB3	1:J:304:VAL:HG23	1.67	0.76
1:H:438:TYR:O	2:N:130[B]:ARG:HD2	1.86	0.75
2:M:157[A]:ILE:HG22	2:M:159[A]:THR:H	1.50	0.75
2:K:162:ASP:O	2:K:165:LYS:HG2	1.87	0.75
2:N:159[B]:THR:OG1	2:N:161[B]:GLU:HG2	1.87	0.75
2:O:130[A]:ARG:NH2	2:O:155[A]:ARG:HD2	2.03	0.72
1:C:305:ASN:HD21	1:C:309:GLN:HE21	1.37	0.72
1:C:79:SER:HB2	1:D:79:SER:OG	1.89	0.72
1:B:10:THR:OG1	1:B:141:THR:HG21	1.90	0.72
1:E:216:VAL:HG23	5:E:2646:HOH:O	1.88	0.72
1:J:20:VAL:HG21	1:J:332:GLU:HG2	1.69	0.72
1:E:249:LYS:HD2	1:E:255:ASP:CG	2.10	0.72
2:N:140[B]:GLU:OE2	5:N:1791:HOH:O	2.08	0.71
2:O:159[B]:THR:HG22	2:O:162[B]:ASP:OD2	1.90	0.71
2:M:133[B]:PRO:O	2:M:134[B]:ALA:HB3	1.91	0.70
2:L:164[A]:LEU:O	2:L:167[A]:VAL:HG22	1.91	0.70
1:E:28:LEU:HD12	1:E:339:VAL:HG12	1.72	0.70
1:G:397:ASN:HB3	5:H:2648:HOH:O	1.92	0.70
2:M:164[B]:LEU:O	2:M:167[B]:VAL:HG12	1.92	0.70
2:K:162:ASP:HA	2:K:165:LYS:HD3	1.72	0.70
1:F:122:THR:OG1	1:F:126:GLN:HG2	1.91	0.70
1:G:308:PHE:HE1	1:G:337:ILE:HD11	1.55	0.70
2:O:138[B]:ILE:HD11	2:O:163[B]:ALA:HB3	1.73	0.70
1:I:258:ILE:HD13	1:I:269:ILE:HD11	1.73	0.69
1:H:301:ARG:NH1	3:H:2510:SO4:O1	2.25	0.69
1:I:35:ILE:HD11	1:I:139:ILE:HD12	1.74	0.69
1:A:297:ASP:HB2	1:A:298:PRO:HD2	1.74	0.69

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:O:157[B]:ILE:H	2:O:157[B]:ILE:HD12	1.58	0.69
2:L:164[B]:LEU:O	2:L:167[B]:VAL:HG22	1.91	0.69
2:O:133[B]:PRO:O	2:O:134[B]:ALA:HB3	1.93	0.68
2:O:140[A]:GLU:HG2	2:O:141[A]:LYS:H	1.57	0.68
1:G:409:GLN:HG2	1:G:412:THR:OG1	1.93	0.68
1:I:182:VAL:HG12	1:I:271:CYS:SG	2.33	0.68
1:H:288:GLY:O	1:H:292:LEU:HD13	1.94	0.68
1:I:62:TYR:CD1	1:I:65:MET:HE3	2.29	0.68
2:L:144[A]:LEU:HD21	2:L:166[A]:LEU:HG	1.75	0.67
2:M:131[A]:LEU:HD11	2:M:146[A]:ALA:HB2	1.75	0.67
2:O:135[A]:ALA:O	2:O:139[A]:LEU:HB2	1.95	0.67
1:E:51:ILE:HB	1:E:52:PRO:HD3	1.77	0.66
1:J:248:LYS:HD2	1:J:254:ILE:HG12	1.77	0.65
1:J:121:ILE:HG21	1:J:292:LEU:HD21	1.77	0.65
1:D:60:SER:OG	1:D:199:ARG:HD2	1.96	0.65
1:D:96:VAL:O	1:D:100:THR:HG23	1.95	0.65
1:E:370:SER:H	1:E:373:GLN:HE21	1.42	0.65
1:H:121:ILE:HG21	1:H:292:LEU:HD21	1.78	0.65
2:L:159[A]:THR:HG22	2:L:162[A]:ASP:OD2	1.96	0.64
2:N:139[A]:LEU:HD21	2:N:146[A]:ALA:HB2	1.79	0.64
2:N:158[A]:PHE:O	2:N:162[A]:ASP:HB2	1.98	0.64
1:I:444:ASP:OD2	2:O:133[B]:PRO:O	2.16	0.64
2:L:158[A]:PHE:HE1	2:L:163[A]:ALA:HB2	1.62	0.64
1:E:64:HIS:CE1	3:E:2499:SO4:O4	2.50	0.64
2:O:133[A]:PRO:O	2:O:134[A]:ALA:HB3	1.98	0.64
1:I:307:ARG:NH1	5:I:2581:HOH:O	2.31	0.64
1:B:62:TYR:HD1	1:B:65:MET:HE3	1.62	0.64
1:E:79:SER:HB2	1:F:79:SER:HG	1.61	0.64
1:D:297:ASP:HB2	1:D:298:PRO:HD2	1.80	0.63
1:I:318:ILE:HD13	1:I:335:GLY:HA2	1.80	0.63
2:O:160[B]:LYS:O	2:O:164[B]:LEU:HD23	1.98	0.63
1:I:265:LYS:O	1:I:265:LYS:HD2	1.99	0.63
1:G:121:ILE:HG21	1:G:292:LEU:HD11	1.81	0.63
1:I:69:THR:HG22	5:I:2522:HOH:O	1.99	0.62
2:O:130[A]:ARG:HH21	2:O:155[A]:ARG:HD2	1.63	0.62
1:I:211:GLY:N	3:I:2515:SO4:O2	2.31	0.62
2:L:133[A]:PRO:O	2:L:134[A]:ALA:HB3	1.99	0.62
1:E:249:LYS:HD2	1:E:255:ASP:OD1	2.00	0.62
1:C:5:ILE:HG23	1:C:137:GLN:NE2	2.14	0.62
2:K:133:PRO:O	2:K:134:ALA:HB3	1.99	0.62
1:A:382:LYS:HG2	1:A:408:GLY:O	1.99	0.62
1:A:288:GLY:O	1:A:292:LEU:HG	2.00	0.62

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:369:LYS:HA	1:E:373:GLN:NE2	2.15	0.61
1:I:84:ASN:HB2	1:J:77:GLU:OE1	1.99	0.61
1:A:25:ALA:HA	1:A:339:VAL:HG11	1.81	0.61
1:B:312:ILE:HD12	1:B:312:ILE:N	2.15	0.61
2:O:159[B]:THR:CG2	2:O:162[B]:ASP:H	2.13	0.61
2:M:158[A]:PHE:O	2:M:162[A]:ASP:HB2	1.99	0.61
2:L:130[B]:ARG:HB2	2:L:157[B]:ILE:HA	1.83	0.61
1:I:51:ILE:HG23	1:J:396:THR:OG1	2.00	0.61
2:M:143[A]:SER:O	2:M:144[A]:LEU:HB3	2.01	0.61
1:I:33:VAL:HG11	1:I:139:ILE:HD13	1.83	0.60
1:A:290:GLU:OE1	1:A:290:GLU:HA	2.01	0.60
1:A:305:ASN:HD21	1:A:309:GLN:HB2	1.65	0.60
1:I:182:VAL:HG13	1:I:274:LEU:HD12	1.83	0.60
1:I:161:ASP:OD1	1:I:163:ASP:HB3	2.01	0.60
2:N:145[A]:ASP:O	2:N:146[A]:ALA:HB3	2.01	0.60
1:J:15:GLY:HA3	4:J:480:FAD:H52A	1.84	0.60
1:C:277:CYS:HA	5:C:2512:HOH:O	2.01	0.60
2:N:138[A]:ILE:HD11	2:N:163[A]:ALA:HB3	1.83	0.60
1:J:290:GLU:HB3	3:J:2520:SO4:O1	2.00	0.60
1:E:62:TYR:HD1	1:E:65:MET:HE2	1.66	0.60
2:M:134[B]:ALA:HA	2:M:137[B]:ASN:HB3	1.83	0.60
1:B:402:GLY:HA3	1:B:421:LEU:O	2.02	0.60
1:B:182:VAL:HG23	1:B:271:CYS:HB3	1.84	0.60
2:L:156[B]:GLY:O	2:L:157[B]:ILE:HB	2.02	0.60
1:H:292:LEU:CD1	5:H:2814:HOH:O	2.49	0.60
1:C:5:ILE:HG23	1:C:137:GLN:HE22	1.66	0.60
1:B:43:GLY:HA2	4:B:480:FAD:O3B	2.02	0.60
1:H:3:GLN:HB2	1:H:4:PRO:HD3	1.84	0.59
2:N:164[B]:LEU:HA	2:N:167[B]:VAL:HG22	1.84	0.59
1:B:301:ARG:HH11	1:B:323:ALA:HA	1.67	0.59
1:B:9:VAL:HG22	1:B:342:MET:HE1	1.85	0.59
1:B:9:VAL:CG2	1:B:342:MET:HE1	2.32	0.59
1:B:384:GLY:HA3	1:B:466:ALA:HB2	1.85	0.59
1:G:308:PHE:CE1	1:G:337:ILE:HD11	2.36	0.59
1:E:369:LYS:HA	1:E:373:GLN:HE22	1.67	0.59
1:B:328:ALA:O	1:B:332:GLU:HG3	2.03	0.59
1:H:375:LYS:HE3	1:H:381:TYR:OH	2.03	0.59
1:I:57:LEU:HD22	1:I:359:TYR:O	2.03	0.58
1:I:307:ARG:NH2	1:I:345:GLY:O	2.36	0.58
2:M:152[B]:THR:HG22	2:M:153[B]:GLY:H	1.69	0.58
2:N:135[A]:ALA:O	2:N:139[A]:LEU:HB2	2.03	0.58
2:M:138[B]:ILE:HD11	2:M:163[B]:ALA:CB	2.32	0.58

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:L:144[B]:LEU:HD11	2:L:167[B]:VAL:HA	1.86	0.58
1:A:51:ILE:HB	1:A:52:PRO:HD3	1.84	0.58
1:I:389:ALA:HA	1:I:400:THR:HB	1.86	0.58
2:O:159[B]:THR:HG22	2:O:162[B]:ASP:CG	2.23	0.58
1:B:370:SER:H	1:B:373:GLN:HE21	1.51	0.58
2:M:167[B]:VAL:HG13	2:M:168[B]:GLN:N	2.19	0.58
1:E:43:GLY:HA2	4:E:480:FAD:O3B	2.03	0.58
1:D:309:GLN:HG2	1:D:316:TYR:CE2	2.39	0.58
1:H:372:GLU:O	1:H:376:GLU:HG3	2.03	0.58
1:B:116:ASN:O	1:B:131:LYS:HG2	2.03	0.58
1:B:144:ILE:HD12	1:B:315:ILE:HD11	1.86	0.58
1:J:8:ASP:HA	1:J:142:LYS:HD3	1.84	0.58
1:E:62:TYR:HD1	1:E:65:MET:CE	2.17	0.57
1:A:96:VAL:O	1:A:100:THR:HG23	2.04	0.57
2:N:134[B]:ALA:O	2:N:138[B]:ILE:HG23	2.03	0.57
2:O:135[B]:ALA:O	2:O:139[B]:LEU:HB2	2.05	0.57
1:E:259:GLU:HB2	1:E:264:GLY:O	2.03	0.57
2:N:131[B]:LEU:HD12	2:N:158[B]:PHE:HB3	1.86	0.57
2:N:146[A]:ALA:C	2:N:148[A]:GLN:H	2.08	0.57
1:I:51:ILE:HB	1:I:52:PRO:HD3	1.86	0.57
1:C:250:SER:HB3	2:K:131:LEU:HD11	1.87	0.57
1:E:349:ILE:O	1:E:350:ASP:HB3	2.04	0.57
1:I:62:TYR:HD1	1:I:65:MET:HE3	1.68	0.57
1:I:294:ILE:HD11	1:I:312:ILE:HD12	1.86	0.57
1:I:229:ILE:O	1:I:233:GLN:HG3	2.05	0.57
1:E:174:LEU:HD12	1:E:197:TRP:CE2	2.40	0.57
1:B:348:HIS:CD2	1:B:349:ILE:H	2.22	0.57
1:B:139:ILE:HD12	1:B:139:ILE:N	2.20	0.57
1:I:231:GLN:HA	1:I:235:PHE:O	2.05	0.57
1:E:209:PHE:O	1:E:240:ASN:HA	2.05	0.56
1:F:301:ARG:HD3	5:F:2713:HOH:O	2.04	0.56
2:O:157[B]:ILE:HD12	2:O:157[B]:ILE:N	2.19	0.56
1:A:5:ILE:HG22	1:A:137:GLN:NE2	2.19	0.56
1:B:289:LEU:HB3	1:B:294:ILE:HB	1.85	0.56
1:G:28:LEU:HD12	1:G:339:VAL:HG12	1.86	0.56
1:F:15:GLY:HA3	4:F:480:FAD:H52A	1.88	0.56
1:J:46:LEU:O	1:J:46:LEU:HD12	2.05	0.56
1:J:318:ILE:HD13	1:J:335:GLY:HA2	1.87	0.56
1:H:171:ALA:HB2	1:H:275:LEU:HD13	1.88	0.56
1:I:57:LEU:HD13	1:I:359:TYR:HB2	1.88	0.56
1:A:43:GLY:HA2	4:A:480:FAD:O3B	2.05	0.56
2:L:159[A]:THR:HG23	2:L:162[A]:ASP:H	1.70	0.55

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:O:134[B]:ALA:HA	2:O:137[B]:ASN:HB3	1.88	0.55
1:C:182:VAL:HG23	1:C:271:CYS:HB3	1.86	0.55
1:E:33:VAL:HG22	1:E:113:VAL:HB	1.87	0.55
2:L:135[B]:ALA:O	2:L:139[B]:LEU:HG	2.06	0.55
1:H:371:GLU:O	1:H:375:LYS:HG2	2.05	0.55
1:D:402:GLY:HA3	1:D:421:LEU:O	2.05	0.55
2:M:156[B]:GLY:O	2:M:157[B]:ILE:HG23	2.07	0.55
1:G:23:ILE:O	1:G:27:GLN:HG3	2.05	0.55
2:N:133[B]:PRO:O	2:N:134[B]:ALA:HB3	2.07	0.55
1:C:318:ILE:HD13	1:C:335:GLY:HA2	1.88	0.55
2:M:134[B]:ALA:O	2:M:138[B]:ILE:HG23	2.06	0.55
1:I:5:ILE:HD12	1:I:6:ASP:H	1.72	0.55
1:E:62:TYR:CD1	1:E:65:MET:HE2	2.41	0.55
1:G:307:ARG:HG2	1:G:347:VAL:HG21	1.88	0.55
1:D:318:ILE:HD13	1:D:335:GLY:HA2	1.89	0.55
1:B:137:GLN:HG3	1:B:139:ILE:HD11	1.89	0.55
1:G:265:LYS:O	1:G:266:ALA:HB2	2.07	0.55
1:D:444:ASP:OD2	2:L:133[A]:PRO:O	2.24	0.55
1:A:49:GLY:O	1:A:52:PRO:HD2	2.05	0.55
1:A:210:LEU:HB3	3:A:2484:SO4:O2	2.06	0.55
1:C:437:GLU:HG2	1:D:448:VAL:HG22	1.90	0.54
1:J:51:ILE:HB	1:J:52:PRO:HD3	1.89	0.54
2:O:138[B]:ILE:HD11	2:O:163[B]:ALA:CB	2.38	0.54
1:H:77:GLU:C	1:H:78:MET:HG3	2.28	0.54
1:D:56:LEU:O	1:D:60:SER:CB	2.55	0.54
1:D:392:SER:HB2	5:D:2701:HOH:O	2.07	0.54
1:J:49:GLY:O	1:J:52:PRO:HD2	2.06	0.54
1:B:305:ASN:OD1	1:B:309:GLN:HB2	2.07	0.54
1:B:312:ILE:HD12	1:B:312:ILE:H	1.73	0.54
1:C:162:GLU:HA	1:C:166:VAL:HG12	1.90	0.54
2:L:155[A]:ARG:HE	2:L:155[A]:ARG:N	2.06	0.54
1:B:121:ILE:HG21	1:B:292:LEU:HD21	1.87	0.54
2:L:158[A]:PHE:CE1	2:L:163[A]:ALA:HB2	2.42	0.54
1:H:3:GLN:CB	1:H:4:PRO:HD3	2.38	0.53
1:H:43:GLY:HA2	4:H:480:FAD:O3B	2.08	0.53
2:O:136[B]:ARG:HH11	2:O:136[B]:ARG:HG2	1.73	0.53
1:C:15:GLY:HA3	4:C:480:FAD:H52A	1.89	0.53
1:H:438:TYR:O	2:N:157[A]:ILE:HD11	2.07	0.53
1:C:305:ASN:HD21	1:C:309:GLN:NE2	2.05	0.53
2:L:131[A]:LEU:HD13	2:L:136[A]:ARG:HG3	1.91	0.53
1:B:349:ILE:O	1:B:349:ILE:HG23	2.08	0.53
1:B:138:VAL:C	1:B:139:ILE:HD12	2.28	0.53

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:108:LYS:HB2	1:C:108:LYS:NZ	2.23	0.53
1:G:15:GLY:HA3	4:G:480:FAD:H52A	1.90	0.53
1:D:444:ASP:OD2	2:L:134[A]:ALA:HB3	2.09	0.53
1:C:289:LEU:N	3:C:2493:SO4:O4	2.41	0.53
1:D:43:GLY:HA2	4:D:480:FAD:O3B	2.08	0.53
1:I:438:TYR:O	2:O:130[A]:ARG:HD2	2.09	0.53
1:B:370:SER:N	1:B:373:GLN:NE2	2.54	0.53
1:H:228:ARG:NH1	1:H:228:ARG:HB2	2.24	0.53
1:F:43:GLY:HA2	4:F:480:FAD:O3B	2.09	0.53
2:L:135[A]:ALA:HA	2:L:138[A]:ILE:HD11	1.91	0.53
1:G:43:GLY:HA2	4:G:480:FAD:O3B	2.07	0.53
1:G:145:LEU:HD11	1:G:318:ILE:HG12	1.89	0.53
2:L:159[A]:THR:HG22	2:L:162[A]:ASP:CG	2.29	0.52
1:I:43:GLY:HA2	4:I:480:FAD:O3B	2.10	0.52
1:J:309:GLN:HG2	1:J:316:TYR:CE2	2.45	0.52
1:C:79:SER:HB2	1:D:79:SER:CB	2.40	0.52
1:B:318:ILE:HD13	1:B:335:GLY:HA2	1.91	0.52
1:B:180:LYS:HE3	5:B:2575:HOH:O	2.09	0.52
1:B:149:GLY:HA2	1:B:320:ASP:HB2	1.91	0.52
1:F:23:ILE:O	1:F:27:GLN:HG3	2.10	0.52
1:C:23:ILE:O	1:C:27:GLN:HG3	2.09	0.52
1:I:52:PRO:HB3	1:I:92:LYS:HD3	1.90	0.52
1:A:46:LEU:HD11	1:A:99:LEU:HB2	1.90	0.52
2:L:134[A]:ALA:O	2:L:138[A]:ILE:HG12	2.10	0.52
1:C:121:ILE:HG21	1:C:292:LEU:HD11	1.91	0.52
1:J:23:ILE:O	1:J:27:GLN:HG3	2.09	0.52
1:J:121:ILE:CG2	1:J:292:LEU:HD21	2.38	0.52
1:J:69:THR:OG1	3:J:2518:SO4:O4	2.27	0.52
1:E:318:ILE:HD13	1:E:335:GLY:HA2	1.91	0.52
1:I:351:TYR:HA	1:I:354:VAL:HG23	1.92	0.52
1:A:5:ILE:HG13	1:A:6:ASP:N	2.24	0.52
1:I:311:LYS:HD2	5:I:2563:HOH:O	2.10	0.52
1:C:110:ASN:O	1:C:111:LYS:HB2	2.09	0.52
5:C:2664:HOH:O	1:D:397:ASN:HB3	2.09	0.52
1:H:46:LEU:O	1:H:46:LEU:HD12	2.10	0.52
1:B:464:LEU:HD23	1:B:471:SER:HA	1.91	0.52
1:J:7:ALA:O	1:J:141:THR:HA	2.09	0.52
1:H:282:PRO:HG2	1:H:301:ARG:HG3	1.92	0.51
1:D:7:ALA:O	1:D:141:THR:HA	2.10	0.51
1:E:28:LEU:HD12	1:E:339:VAL:CG1	2.41	0.51
1:B:370:SER:OG	1:B:373:GLN:HG3	2.10	0.51
1:E:28:LEU:CD1	1:E:339:VAL:HG12	2.41	0.51

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:F:7:ALA:O	1:F:141:THR:HA	2.11	0.51
1:H:121:ILE:CG2	1:H:292:LEU:HD21	2.39	0.51
1:B:62:TYR:CD1	1:B:65:MET:HE3	2.44	0.51
1:G:461:GLU:OE1	1:G:471:SER:HB2	2.10	0.51
2:O:131[A]:LEU:HD21	2:O:146[A]:ALA:HB2	1.93	0.51
2:N:142[A]:HIS:O	2:N:144[A]:LEU:HD22	2.10	0.51
2:K:134:ALA:HA	2:K:137:ASN:HB3	1.92	0.51
1:J:43:GLY:HA2	4:J:480:FAD:O3B	2.10	0.51
1:D:5:ILE:HG22	1:D:6:ASP:N	2.26	0.51
1:G:434:LEU:HD12	1:H:448:VAL:HG21	1.93	0.51
1:B:69:THR:HG23	3:B:2487:SO4:O2	2.10	0.51
2:O:131[B]:LEU:HD21	2:O:146[B]:ALA:HB2	1.94	0.51
1:J:15:GLY:CA	4:J:480:FAD:H52A	2.40	0.51
1:H:80:GLU:HG3	1:H:82:ARG:HH22	1.76	0.51
1:A:402:GLY:HA3	1:A:421:LEU:O	2.11	0.50
2:O:158[B]:PHE:HZ	2:O:166[B]:LEU:HD12	1.75	0.50
2:M:133[B]:PRO:O	2:M:134[B]:ALA:CB	2.55	0.50
1:C:258:ILE:HD13	1:C:269:ILE:HD13	1.92	0.50
2:N:130[A]:ARG:O	2:N:157[A]:ILE:HA	2.11	0.50
1:E:346:ALA:HB1	2:M:140[A]:GLU:CD	2.32	0.50
1:E:5:ILE:HD12	1:E:6:ASP:H	1.76	0.50
1:I:145:LEU:HD11	1:I:318:ILE:HG12	1.94	0.50
1:G:318:ILE:HD13	1:G:335:GLY:CA	2.42	0.50
1:H:80:GLU:HG3	1:H:82:ARG:NH2	2.25	0.50
1:C:447:ARG:CZ	2:L:137[B]:ASN:ND2	2.75	0.50
1:B:146:ILE:HG12	1:B:316:TYR:O	2.12	0.50
1:C:28:LEU:HD12	1:C:339:VAL:HG12	1.93	0.50
2:M:135[A]:ALA:O	2:M:139[A]:LEU:HB2	2.11	0.50
2:O:133[B]:PRO:O	2:O:134[B]:ALA:CB	2.59	0.50
1:C:162:GLU:HA	1:C:162:GLU:OE2	2.12	0.50
1:C:191:VAL:HG21	1:C:213:VAL:CG1	2.42	0.50
1:I:8:ASP:HA	1:I:142:LYS:HG3	1.93	0.50
1:A:124:LYS:HE2	1:A:312:ILE:HD12	1.94	0.50
1:F:5:ILE:HG12	1:F:137:GLN:NE2	2.26	0.50
1:F:318:ILE:HD13	1:F:335:GLY:CA	2.42	0.50
1:B:39:GLU:H	1:B:39:GLU:CD	2.15	0.50
1:D:438:TYR:O	2:L:130[B]:ARG:HD2	2.11	0.50
1:I:5:ILE:HD12	1:I:6:ASP:N	2.27	0.50
1:H:15:GLY:HA3	4:H:480:FAD:H52A	1.93	0.50
1:A:25:ALA:CA	1:A:339:VAL:HG11	2.42	0.49
1:J:389:ALA:HA	1:J:400:THR:HB	1.94	0.49
1:J:444:ASP:OD2	2:O:133[A]:PRO:O	2.30	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:J:437:GLU:HG3	2:O:133[B]:PRO:HG3	1.94	0.49
1:B:278:ILE:H	1:B:278:ILE:HD13	1.77	0.49
1:G:299:ARG:HB2	1:G:301:ARG:HD3	1.93	0.49
1:A:46:LEU:HD11	1:A:99:LEU:CB	2.42	0.49
1:C:178:PRO:HG3	1:C:181:MET:CE	2.42	0.49
2:M:156[B]:GLY:C	2:M:157[B]:ILE:HG13	2.32	0.49
2:M:158[B]:PHE:HE1	2:M:163[B]:ALA:HB2	1.76	0.49
2:L:139[A]:LEU:HG	2:L:163[A]:ALA:HB1	1.94	0.49
1:B:33:VAL:HG22	1:B:113:VAL:HB	1.95	0.49
1:G:402:GLY:HA3	1:G:421:LEU:O	2.12	0.49
1:F:96:VAL:O	1:F:100:THR:HG23	2.12	0.49
1:D:244:THR:OG1	1:D:257:SER:OG	2.28	0.49
1:J:447:ARG:CD	2:O:137[A]:ASN:ND2	2.75	0.49
1:C:178:PRO:HG3	1:C:181:MET:HE2	1.95	0.49
1:E:437:GLU:HG3	2:M:133[A]:PRO:HB3	1.95	0.49
1:G:264:GLY:O	1:G:265:LYS:O	2.31	0.49
1:B:341:GLY:HA2	1:B:345:GLY:O	2.13	0.49
1:B:285:LYS:HG2	1:B:286:ASN:HD22	1.78	0.49
1:I:205:THR:HA	1:I:236:LYS:O	2.13	0.49
1:B:311:LYS:O	1:B:313:PRO:HD3	2.12	0.49
1:A:305:ASN:ND2	1:A:309:GLN:HB2	2.26	0.49
1:B:404:VAL:HG22	1:B:420:ILE:HG23	1.95	0.49
1:E:439:GLY:HA3	2:M:154[B]:PRO:HG3	1.94	0.49
1:I:192:GLU:O	1:I:196:VAL:HG23	2.12	0.49
1:D:145:LEU:HD11	1:D:318:ILE:HG12	1.94	0.49
1:A:162:GLU:HA	1:A:162:GLU:OE1	2.12	0.49
1:E:272:ASP:OD2	5:E:2540:HOH:O	2.20	0.49
1:B:7:ALA:CB	1:B:31:LYS:HD2	2.42	0.48
1:J:288:GLY:O	1:J:292:LEU:HD13	2.12	0.48
1:H:64:HIS:CE1	3:H:2511:SO4:O1	2.66	0.48
1:F:318:ILE:HD13	1:F:335:GLY:HA2	1.94	0.48
2:N:154[B]:PRO:HD2	2:N:157[B]:ILE:CG2	2.43	0.48
1:D:389:ALA:HA	1:D:400:THR:HB	1.95	0.48
1:A:226:PHE:CZ	1:A:230:LEU:HD11	2.48	0.48
1:B:144:ILE:HD12	1:B:315:ILE:CD1	2.44	0.48
2:M:164[A]:LEU:O	2:M:167[A]:VAL:HG22	2.13	0.48
1:D:15:GLY:HA3	4:D:480:FAD:H52A	1.96	0.48
1:E:52:PRO:HB2	1:E:172:LEU:HD13	1.95	0.48
1:C:43:GLY:HA2	4:C:480:FAD:O3B	2.13	0.48
2:K:131:LEU:HD21	2:K:146:ALA:CB	2.43	0.48
1:F:301:ARG:NH1	3:F:2502:SO4:O3	2.46	0.48
1:B:162:GLU:OE1	1:B:162:GLU:HA	2.13	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:G:148:THR:OG1	1:G:321:VAL:HG11	2.14	0.48
1:A:290:GLU:C	1:A:292:LEU:H	2.16	0.48
1:A:337:ILE:HG13	1:A:338:CYS:N	2.29	0.48
1:C:39:GLU:HG2	1:C:40:THR:HG23	1.95	0.48
2:O:132[A]:SER:O	2:O:135[A]:ALA:HB3	2.14	0.48
1:B:7:ALA:HB2	1:B:31:LYS:HD2	1.95	0.48
1:A:46:LEU:HD21	1:A:100:THR:HA	1.96	0.48
1:G:174:LEU:HD12	1:G:197:TRP:CE2	2.49	0.48
1:B:249:LYS:HA	1:J:305:ASN:HA	1.96	0.48
1:F:330:LYS:HD3	1:F:334:GLU:CD	2.35	0.48
1:A:448:VAL:HG21	1:B:434:LEU:HD12	1.95	0.47
2:N:138[B]:ILE:HD11	2:N:163[B]:ALA:CB	2.44	0.47
2:N:164[B]:LEU:O	2:N:167[B]:VAL:HG22	2.14	0.47
1:J:7:ALA:HA	1:J:31:LYS:HD3	1.96	0.47
1:H:402:GLY:HA3	1:H:421:LEU:O	2.15	0.47
1:I:46:LEU:HD11	1:I:99:LEU:HB2	1.95	0.47
1:E:210:LEU:N	1:E:210:LEU:HD12	2.30	0.47
2:L:156[B]:GLY:C	2:L:157[B]:ILE:HD12	2.34	0.47
2:M:131[B]:LEU:HD13	2:M:136[B]:ARG:HA	1.96	0.47
1:D:461:GLU:OE1	1:D:471:SER:HB2	2.14	0.47
1:B:370:SER:CB	1:B:373:GLN:HE21	2.26	0.47
1:B:442:CYS:HB3	1:B:463:ASN:O	2.14	0.47
1:B:118:TYR:O	1:B:129:ALA:HA	2.15	0.47
1:H:96:VAL:O	1:H:100:THR:HG23	2.15	0.47
1:F:311:LYS:HE2	5:F:2545:HOH:O	2.14	0.47
1:B:121:ILE:CG2	1:B:292:LEU:HD21	2.44	0.47
2:O:140[A]:GLU:HG2	2:O:141[A]:LYS:N	2.26	0.47
1:D:56:LEU:O	1:D:60:SER:HB2	2.15	0.47
1:A:3:GLN:N	1:A:4:PRO:CD	2.78	0.47
1:C:389:ALA:HA	1:C:400:THR:HB	1.96	0.47
1:H:155:PHE:CD1	1:H:158:ILE:HD12	2.49	0.47
1:B:350:ASP:C	1:B:352:ASN:H	2.18	0.47
1:E:259:GLU:HB3	1:E:266:ALA:HA	1.97	0.47
1:G:318:ILE:HD13	1:G:335:GLY:HA2	1.97	0.47
1:B:13:GLY:O	1:B:18:GLY:HA3	2.15	0.47
1:E:349:ILE:O	1:E:350:ASP:CB	2.62	0.47
1:F:15:GLY:CA	4:F:480:FAD:H52A	2.44	0.47
2:O:134[B]:ALA:HA	2:O:137[B]:ASN:CB	2.45	0.47
1:B:442:CYS:CB	1:B:467:SER:HB2	2.45	0.47
1:I:165:ILE:HD11	1:I:254:ILE:HD12	1.97	0.46
1:E:64:HIS:NE2	3:E:2499:SO4:O4	2.48	0.46
1:A:46:LEU:HD12	1:A:51:ILE:HG13	1.97	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
5:G:2556:HOH:O	1:H:397:ASN:HB3	2.15	0.46
1:E:306:THR:O	1:E:347:VAL:HG11	2.15	0.46
1:B:28:LEU:HD12	1:B:339:VAL:HG12	1.97	0.46
1:A:35:ILE:HG12	1:A:115:VAL:HB	1.98	0.46
1:A:461:GLU:CG	1:A:471:SER:HB2	2.27	0.46
2:L:160[B]:LYS:O	2:L:164[B]:LEU:HG	2.15	0.46
1:I:258:ILE:CD1	1:I:269:ILE:HD11	2.44	0.46
1:A:15:GLY:HA3	4:A:480:FAD:H52A	1.97	0.46
1:E:210:LEU:CD1	1:E:210:LEU:N	2.78	0.46
1:A:286:ASN:HA	3:A:2485:SO4:O2	2.15	0.46
1:B:347:VAL:O	1:B:347:VAL:HG23	2.15	0.46
2:M:145[A]:ASP:C	2:M:147[A]:SER:H	2.18	0.46
1:D:287:LEU:HB3	1:D:289:LEU:HG	1.97	0.46
1:H:299:ARG:NH2	3:H:2510:SO4:O4	2.49	0.46
2:N:138[B]:ILE:CD1	2:N:164[B]:LEU:HD22	2.45	0.46
1:E:301:ARG:NH1	3:E:2498:SO4:O4	2.48	0.46
1:A:397:ASN:HB3	5:B:2533:HOH:O	2.14	0.46
2:K:133:PRO:O	2:K:134:ALA:CB	2.64	0.46
1:B:246:ALA:HA	1:B:255:ASP:O	2.16	0.46
1:B:444:ASP:OD1	2:K:134:ALA:HB3	2.14	0.46
1:B:442:CYS:HB2	1:B:467:SER:HB2	1.97	0.46
1:F:351:TYR:HA	1:F:354:VAL:HG13	1.98	0.46
1:A:299:ARG:CZ	1:A:301:ARG:HH12	2.28	0.46
1:C:171:ALA:HB2	1:C:275:LEU:HD13	1.98	0.46
1:I:370:SER:OG	1:I:373:GLN:HG3	2.15	0.46
2:K:139:LEU:CD2	2:K:146:ALA:HB2	2.46	0.46
1:J:212:HIS:ND1	3:J:2519:SO4:O3	2.48	0.46
1:J:155:PHE:CD2	1:J:278:ILE:HD13	2.50	0.46
2:M:167[B]:VAL:CG1	2:M:168[B]:GLN:N	2.79	0.46
1:I:258:ILE:HD13	1:I:269:ILE:CD1	2.45	0.46
1:J:164:THR:HG21	1:J:248:LYS:NZ	2.31	0.46
1:G:122:THR:OG1	1:G:126:GLN:HG2	2.16	0.46
1:E:62:TYR:CD1	1:E:65:MET:CE	2.98	0.45
1:H:15:GLY:N	4:H:480:FAD:H52A	2.31	0.45
1:E:321:VAL:HG22	1:E:321:VAL:O	2.16	0.45
1:I:290:GLU:HG3	1:I:291:GLU:N	2.31	0.45
2:O:157[A]:ILE:CG2	2:O:159[A]:THR:HG23	2.45	0.45
1:J:290:GLU:CB	3:J:2520:SO4:O1	2.63	0.45
1:F:171:ALA:HB2	1:F:275:LEU:HD13	1.99	0.45
1:G:160:ILE:HA	1:G:165:ILE:HG22	1.98	0.45
1:C:46:LEU:HD12	1:C:46:LEU:O	2.16	0.45
2:O:167[A]:VAL:HG23	2:O:168[A]:GLN:N	2.32	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:O:157[A]:ILE:HG22	2:O:159[A]:THR:HG23	1.99	0.45
2:O:160[A]:LYS:O	2:O:164[A]:LEU:HD13	2.15	0.45
1:G:160:ILE:HD11	1:G:276:VAL:HB	1.97	0.45
1:I:78:MET:HG2	1:J:81:VAL:HG22	1.98	0.45
1:J:180:LYS:HD3	1:J:270:THR:O	2.16	0.45
1:D:121:ILE:HG21	1:D:292:LEU:HD21	1.97	0.45
2:M:157[B]:ILE:O	2:M:158[B]:PHE:C	2.54	0.45
1:G:265:LYS:HD3	1:G:265:LYS:N	2.32	0.45
1:J:447:ARG:HD2	2:O:137[A]:ASN:ND2	2.31	0.45
1:C:439:GLY:O	2:L:130[A]:ARG:HD2	2.17	0.45
1:B:290:GLU:HB3	3:B:2489:SO4:O1	2.16	0.45
1:C:244:THR:OG1	1:C:257:SER:OG	2.34	0.45
1:E:444:ASP:OD2	2:M:133[B]:PRO:O	2.35	0.45
2:L:154[B]:PRO:C	2:L:156[B]:GLY:H	2.19	0.45
2:K:134:ALA:O	2:K:138:ILE:HG13	2.16	0.45
1:B:338:CYS:O	1:B:342:MET:HG3	2.17	0.45
1:E:189:ILE:HG23	4:E:480:FAD:HM73	1.98	0.45
1:B:250:SER:CB	1:J:304:VAL:HG23	2.43	0.45
1:A:309:GLN:HG2	1:A:316:TYR:CE2	2.52	0.45
1:J:155:PHE:HD2	1:J:278:ILE:HD13	1.81	0.45
1:I:426:GLY:O	1:I:429:VAL:HG12	2.17	0.45
1:A:139:ILE:HD12	1:A:139:ILE:N	2.31	0.45
1:B:444:ASP:OD1	2:K:133:PRO:O	2.35	0.45
2:K:131:LEU:HD21	2:K:146:ALA:HB2	1.99	0.45
1:I:124:LYS:HE2	1:I:125:ASN:OD1	2.17	0.45
1:H:180:LYS:HB3	1:H:180:LYS:HE2	1.85	0.45
1:A:222:ILE:HG13	1:A:419:HIS:HB3	1.98	0.45
1:B:277:CYS:HA	5:B:2547:HOH:O	2.17	0.45
1:C:309:GLN:HG2	1:C:316:TYR:CE2	2.52	0.44
1:E:7:ALA:O	1:E:141:THR:HA	2.17	0.44
1:G:383:VAL:HG11	1:G:385:LYS:HE3	1.99	0.44
1:B:57:LEU:HD11	1:B:192:GLU:HB3	1.98	0.44
2:N:145[A]:ASP:O	2:N:146[A]:ALA:CB	2.65	0.44
1:C:15:GLY:CA	4:C:480:FAD:H52A	2.47	0.44
1:F:290:GLU:HB3	3:F:2505:SO4:O4	2.17	0.44
1:C:161:ASP:OD1	1:C:163:ASP:HB3	2.18	0.44
2:O:158[B]:PHE:O	2:O:159[B]:THR:O	2.35	0.44
1:G:437:GLU:HG3	2:N:133[A]:PRO:HG3	2.00	0.44
2:L:140[A]:GLU:C	2:L:142[A]:HIS:H	2.21	0.44
1:A:41:LEU:HD23	1:A:103:ILE:HG22	1.99	0.44
1:A:389:ALA:HA	1:A:400:THR:HB	1.99	0.44
2:N:145[A]:ASP:C	2:N:147[A]:SER:H	2.21	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:79:SER:CB	1:D:79:SER:OG	2.62	0.44
2:K:131:LEU:HD23	2:K:158:PHE:HB3	1.98	0.44
1:B:54:LYS:HE3	1:B:54:LYS:N	2.33	0.44
1:I:79:SER:HB2	1:J:79:SER:CB	2.48	0.44
2:M:158[A]:PHE:O	2:M:159[A]:THR:O	2.35	0.44
1:E:370:SER:N	1:E:373:GLN:HE21	2.12	0.44
1:A:42:GLY:HA3	1:A:46:LEU:HD23	2.00	0.44
1:J:8:ASP:HA	1:J:142:LYS:CD	2.48	0.44
1:F:452:HIS:HA	1:F:453:PRO:HA	1.82	0.44
1:C:174:LEU:HD12	1:C:197:TRP:CE2	2.53	0.44
2:L:131[B]:LEU:HD23	2:L:136[B]:ARG:NH1	2.33	0.44
1:H:68:GLY:HA3	3:H:2511:SO4:O2	2.17	0.44
1:I:393:ARG:HD2	1:I:454:THR:HA	1.98	0.44
1:A:33:VAL:HG22	1:A:113:VAL:HB	1.98	0.44
2:N:144[B]:LEU:HD22	2:N:167[B]:VAL:HG12	1.99	0.44
1:B:62:TYR:CD1	1:B:65:MET:CE	3.00	0.44
1:B:148:THR:OG1	1:B:321:VAL:HG11	2.18	0.44
1:F:10:THR:HB	1:F:141:THR:HG21	2.00	0.44
1:F:249:LYS:HD3	1:F:255:ASP:OD2	2.18	0.44
1:E:78:MET:CE	1:F:78:MET:HG2	2.48	0.43
1:G:409:GLN:HG3	1:G:412:THR:H	1.81	0.43
1:B:4:PRO:HB3	1:B:138:VAL:HB	2.00	0.43
1:G:437:GLU:HG2	1:H:448:VAL:HG22	1.99	0.43
2:N:158[B]:PHE:HE1	2:N:163[B]:ALA:HA	1.82	0.43
2:M:160[A]:LYS:HE3	2:M:164[A]:LEU:HD11	2.00	0.43
1:E:15:GLY:HA3	4:E:480:FAD:H52A	2.00	0.43
1:C:87:LYS:HE3	1:C:90:GLU:OE2	2.18	0.43
1:F:292:LEU:HG	1:F:312:ILE:CD1	2.48	0.43
2:N:167[B]:VAL:HG23	2:N:168[B]:GLN:N	2.33	0.43
2:M:158[A]:PHE:O	2:M:162[A]:ASP:OD2	2.37	0.43
1:B:330:LYS:HE2	1:B:334:GLU:OE2	2.18	0.43
1:E:312:ILE:HA	1:E:313:PRO:HD2	1.82	0.43
1:I:173:SER:O	1:I:174:LEU:C	2.55	0.43
1:D:46:LEU:O	1:D:46:LEU:HD12	2.17	0.43
1:I:187:GLY:O	1:I:191:VAL:HG23	2.17	0.43
1:I:197:TRP:CE3	1:I:200:LEU:HD12	2.53	0.43
1:J:318:ILE:HD13	1:J:335:GLY:CA	2.49	0.43
1:C:81:VAL:HG22	1:D:78:MET:HG2	2.00	0.43
1:D:222:ILE:HG13	1:D:419:HIS:HB3	2.01	0.43
1:J:171:ALA:HA	1:J:174:LEU:HD13	2.01	0.43
2:N:164[A]:LEU:O	2:N:167[A]:VAL:HG22	2.19	0.43
2:L:144[B]:LEU:CD1	2:L:167[B]:VAL:HA	2.48	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:I:28:LEU:HD12	1:I:339:VAL:HG12	2.01	0.43
2:N:159[B]:THR:N	2:N:162[B]:ASP:HB2	2.33	0.43
1:C:121:ILE:CG2	1:C:292:LEU:HD11	2.48	0.43
1:A:301:ARG:HG2	1:A:322:VAL:C	2.39	0.43
1:G:444:ASP:OD1	2:N:133[B]:PRO:O	2.37	0.43
2:M:143[A]:SER:O	2:M:144[A]:LEU:CB	2.66	0.43
4:H:480:FAD:H51A	5:H:2517:HOH:O	2.18	0.43
1:D:288:GLY:O	1:D:292:LEU:HB2	2.19	0.43
1:C:304:VAL:HB	1:C:308:PHE:HA	2.01	0.43
1:B:46:LEU:HD11	1:B:99:LEU:HB2	2.00	0.43
1:G:209:PHE:O	1:G:240:ASN:HA	2.19	0.43
1:G:389:ALA:HA	1:G:400:THR:HB	2.00	0.43
2:O:131[B]:LEU:HD23	2:O:158[B]:PHE:CB	2.48	0.43
2:M:134[A]:ALA:O	2:M:138[A]:ILE:HG23	2.18	0.43
1:B:7:ALA:O	1:B:141:THR:HA	2.19	0.43
2:O:142[B]:HIS:C	2:O:144[B]:LEU:H	2.22	0.43
2:L:159[A]:THR:O	2:L:162[A]:ASP:HB2	2.18	0.43
1:B:287:LEU:HD23	1:B:289:LEU:HD21	2.01	0.43
1:E:3:GLN:N	1:E:4:PRO:CD	2.81	0.43
1:B:354:VAL:HA	1:B:355:PRO:HD3	1.83	0.43
1:A:203:ASP:OD1	1:A:236:LYS:NZ	2.42	0.43
2:O:141[B]:LYS:O	2:O:142[B]:HIS:HD2	2.02	0.43
1:I:182:VAL:HG13	1:I:274:LEU:CD1	2.47	0.43
1:B:162:GLU:HA	1:B:166:VAL:HG12	2.01	0.43
1:F:162:GLU:OE1	1:F:162:GLU:HA	2.19	0.43
1:C:209:PHE:HA	1:C:241:THR:O	2.19	0.43
1:I:346:ALA:HB1	2:O:140[A]:GLU:CD	2.40	0.42
1:I:254:ILE:HG23	1:I:271:CYS:SG	2.59	0.42
1:C:442:CYS:HB2	1:C:467:SER:HB2	2.01	0.42
1:G:288:GLY:N	3:G:2509:SO4:O4	2.48	0.42
1:I:3:GLN:N	1:I:4:PRO:CD	2.82	0.42
1:A:337:ILE:CD1	1:A:347:VAL:O	2.68	0.42
2:L:157[B]:ILE:O	2:L:158[B]:PHE:HB3	2.19	0.42
1:C:441:SER:CB	2:L:160[B]:LYS:HD3	2.49	0.42
1:C:404:VAL:HG11	1:C:459:PHE:HA	2.01	0.42
1:E:409:GLN:HB3	1:E:412:THR:OG1	2.19	0.42
1:A:116:ASN:O	1:A:131:LYS:HG2	2.20	0.42
2:O:151[A]:ALA:O	2:O:152[A]:THR:HB	2.20	0.42
2:N:146[A]:ALA:C	2:N:148[A]:GLN:N	2.73	0.42
1:B:301:ARG:NH1	3:B:2486:SO4:O4	2.53	0.42
1:D:54:LYS:N	1:D:54:LYS:HE3	2.34	0.42
2:K:168:GLN:C	2:K:170:LYS:H	2.21	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:J:281:ARG:HB2	1:J:282:PRO:HD2	2.00	0.42
2:L:139[A]:LEU:HD22	2:L:144[A]:LEU:O	2.20	0.42
1:H:145:LEU:HD11	1:H:318:ILE:HG12	2.01	0.42
1:D:162:GLU:OE1	1:D:162:GLU:HA	2.19	0.42
2:N:144[A]:LEU:HD22	2:N:144[A]:LEU:N	2.35	0.42
2:O:133[A]:PRO:O	2:O:135[A]:ALA:N	2.43	0.42
2:N:149[B]:GLY:HA3	2:N:166[B]:LEU:HD11	2.02	0.42
1:I:182:VAL:HG11	1:I:256:VAL:HG21	2.01	0.42
1:F:278:ILE:HD12	1:F:279:GLY:N	2.34	0.42
1:B:356:SER:O	1:B:365:ALA:HA	2.19	0.42
1:B:450:HIS:ND1	1:B:460:ARG:HD2	2.35	0.42
2:N:153[A]:GLY:HA3	2:N:159[A]:THR:CG2	2.49	0.42
2:L:131[B]:LEU:HA	2:L:157[B]:ILE:O	2.20	0.42
1:D:56:LEU:O	1:D:60:SER:HB3	2.19	0.42
1:H:15:GLY:CA	4:H:480:FAD:H52A	2.48	0.42
2:M:170[B]:LYS:C	2:M:172[B]:THR:H	2.23	0.42
1:D:453:PRO:HA	1:D:457:GLU:OE1	2.20	0.42
1:A:209:PHE:O	1:A:240:ASN:HA	2.19	0.42
1:I:162:GLU:HA	1:I:162:GLU:OE2	2.19	0.42
2:O:153[B]:GLY:HA3	2:O:154[B]:PRO:HD2	1.96	0.42
2:O:133[A]:PRO:O	2:O:134[A]:ALA:CB	2.64	0.42
2:L:131[B]:LEU:N	2:L:157[B]:ILE:O	2.52	0.42
1:A:339:VAL:HA	1:A:342:MET:HE3	2.02	0.42
1:A:299:ARG:NH1	1:A:301:ARG:HH12	2.17	0.42
1:D:312:ILE:HA	1:D:313:PRO:HD2	1.75	0.42
2:N:133[B]:PRO:O	2:N:135[B]:ALA:N	2.44	0.42
1:A:348:HIS:HB3	2:K:140:GLU:OE1	2.19	0.42
2:O:134[A]:ALA:O	2:O:138[A]:ILE:HG13	2.20	0.42
1:B:301:ARG:NH1	1:B:323:ALA:HA	2.32	0.42
1:G:299:ARG:HB2	1:G:301:ARG:CD	2.50	0.42
1:I:136:THR:O	1:I:136:THR:HG23	2.19	0.42
1:I:225:ASN:O	1:I:229:ILE:HG13	2.20	0.42
1:B:350:ASP:O	1:B:352:ASN:N	2.53	0.42
1:I:282:PRO:HG2	1:I:301:ARG:HG3	2.00	0.42
1:I:301:ARG:O	1:I:303:PRO:HD3	2.20	0.42
1:C:334:GLU:OE2	1:C:351:TYR:OH	2.31	0.42
1:J:239:LEU:O	1:J:240:ASN:C	2.58	0.42
1:B:238:LYS:HD3	1:B:238:LYS:HA	1.86	0.42
1:F:46:LEU:HD12	1:F:46:LEU:O	2.20	0.42
1:D:77:GLU:OE2	1:D:77:GLU:HA	2.21	0.41
1:G:178:PRO:HG3	1:G:181:MET:CE	2.50	0.41
1:J:297:ASP:OD2	1:J:297:ASP:C	2.58	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:265:LYS:HE3	1:C:265:LYS:HB3	1.89	0.41
1:I:396:THR:HG23	1:J:55:ALA:HB2	2.02	0.41
2:O:130[A]:ARG:NH2	2:O:157[A]:ILE:HD11	2.34	0.41
1:A:453:PRO:HA	1:A:457:GLU:OE2	2.20	0.41
1:C:7:ALA:O	1:C:141:THR:HA	2.20	0.41
1:I:438:TYR:O	2:O:154[B]:PRO:HG3	2.20	0.41
1:G:437:GLU:CG	2:N:133[A]:PRO:HG3	2.50	0.41
1:A:339:VAL:HG22	1:A:342:MET:HE1	2.02	0.41
1:H:228:ARG:HB2	1:H:228:ARG:CZ	2.51	0.41
1:C:160:ILE:HD13	1:C:167:SER:HB3	2.01	0.41
1:J:219:ASP:HA	1:J:370:SER:HB2	2.01	0.41
1:F:145:LEU:HD11	1:F:318:ILE:HG12	2.03	0.41
1:J:307:ARG:O	1:J:308:PHE:HB2	2.20	0.41
1:D:150:SER:HB3	1:D:320:ASP:HB3	2.02	0.41
1:J:54:LYS:HE2	1:J:359:TYR:HB2	2.02	0.41
2:L:131[A]:LEU:HB3	2:L:132[A]:SER:H	1.67	0.41
1:B:286:ASN:HA	3:B:2489:SO4:O2	2.21	0.41
1:I:208:GLU:HG3	1:I:210:LEU:H	1.86	0.41
1:E:370:SER:N	1:E:373:GLN:NE2	2.61	0.41
1:B:9:VAL:HG21	1:B:342:MET:HE1	2.01	0.41
1:B:351:TYR:O	1:B:354:VAL:HG12	2.21	0.41
1:H:23:ILE:O	1:H:27:GLN:HG3	2.21	0.41
1:B:174:LEU:HD12	1:B:197:TRP:CE2	2.55	0.41
1:E:162:GLU:OE2	1:E:162:GLU:HA	2.19	0.41
1:E:347:VAL:O	1:E:348:HIS:O	2.39	0.41
1:J:174:LEU:HB2	1:J:197:TRP:CZ2	2.56	0.41
2:M:157[A]:ILE:HG22	2:M:158[A]:PHE:N	2.36	0.41
2:O:133[A]:PRO:C	2:O:135[A]:ALA:H	2.22	0.41
2:O:139[A]:LEU:HD22	2:O:139[A]:LEU:HA	1.87	0.41
1:I:312:ILE:HA	1:I:313:PRO:HD3	1.77	0.41
1:B:335:GLY:O	1:B:339:VAL:HG23	2.21	0.41
1:C:87:LYS:HA	1:C:87:LYS:HD2	1.72	0.41
1:B:453:PRO:HA	1:B:457:GLU:OE1	2.21	0.41
1:E:333:ASP:O	1:E:337:ILE:HG12	2.21	0.41
2:O:152[A]:THR:HB	2:O:162[A]:ASP:OD2	2.20	0.41
1:A:290:GLU:O	1:A:292:LEU:N	2.54	0.41
1:A:25:ALA:HA	1:A:339:VAL:CG1	2.49	0.41
1:E:5:ILE:HG13	1:E:6:ASP:N	2.36	0.41
1:D:77:GLU:C	1:D:78:MET:HG3	2.41	0.41
1:A:349:ILE:HD13	1:A:351:TYR:CZ	2.55	0.41
1:E:389:ALA:HA	1:E:400:THR:HB	2.03	0.41
1:J:222:ILE:HG13	1:J:419:HIS:HB3	2.03	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:F:356:SER:O	1:F:365:ALA:HA	2.21	0.41
1:C:254:ILE:O	1:C:270:THR:HA	2.21	0.41
1:B:297:ASP:HB2	1:B:298:PRO:CD	2.51	0.41
1:B:222:ILE:HG13	1:B:419:HIS:HB3	2.03	0.41
2:M:151[B]:ALA:HB1	2:M:158[B]:PHE:H	1.85	0.41
2:N:149[B]:GLY:HA2	2:N:166[B]:LEU:HD21	2.03	0.41
1:I:211:GLY:CA	3:I:2515:SO4:O2	2.69	0.41
1:B:309:GLN:HG2	1:B:316:TYR:CE2	2.56	0.41
1:H:392:SER:HB2	5:H:2653:HOH:O	2.20	0.41
1:G:57:LEU:HD13	1:G:359:TYR:HB2	2.03	0.41
1:I:224:LYS:HE2	1:I:224:LYS:HA	2.03	0.41
1:A:308:PHE:HE2	1:A:337:ILE:CD1	2.16	0.40
2:N:146[B]:ALA:O	2:N:148[B]:GLN:N	2.54	0.40
1:D:60:SER:OG	1:D:199:ARG:CD	2.68	0.40
1:E:349:ILE:HG23	1:E:351:TYR:CE1	2.56	0.40
1:C:25:ALA:HA	1:C:339:VAL:HG11	2.02	0.40
1:F:253:LYS:HE3	5:F:2587:HOH:O	2.21	0.40
2:N:154[B]:PRO:HD2	2:N:157[B]:ILE:HG21	2.02	0.40
1:B:184:ILE:HG22	1:B:278:ILE:CG2	2.51	0.40
1:B:285:LYS:CG	1:B:286:ASN:HD22	2.34	0.40
1:G:385:LYS:HE2	1:G:405:LYS:HD2	2.04	0.40
1:A:20:VAL:HG21	1:A:332:GLU:HG2	2.02	0.40
1:C:96:VAL:O	1:C:100:THR:HG23	2.20	0.40
1:G:453:PRO:HA	1:G:457:GLU:OE2	2.21	0.40
1:A:311:LYS:O	1:A:313:PRO:HD3	2.21	0.40
1:H:438:TYR:CE1	2:N:132[B]:SER:HB2	2.57	0.40
1:B:148:THR:O	4:B:480:FAD:H8A	2.22	0.40
1:B:24:LYS:O	1:B:28:LEU:HG	2.20	0.40
1:E:287:LEU:HD23	1:E:289:LEU:HD21	2.03	0.40
1:I:160:ILE:O	1:I:160:ILE:HG13	2.21	0.40
1:A:337:ILE:HD12	1:A:347:VAL:O	2.22	0.40
2:N:144[B]:LEU:HD21	2:N:167[B]:VAL:HA	2.03	0.40
1:E:78:MET:HE3	1:F:78:MET:CE	2.51	0.40
1:J:447:ARG:HD2	2:O:137[A]:ASN:HD22	1.86	0.40
1:F:15:GLY:N	4:F:480:FAD:H52A	2.36	0.40
1:B:24:LYS:HA	1:B:24:LYS:HD2	1.84	0.40
1:J:78:MET:HG2	1:J:81:VAL:HG22	2.04	0.40
2:L:159[B]:THR:O	2:L:162[B]:ASP:HB2	2.21	0.40
1:J:145:LEU:HD11	1:J:318:ILE:HG12	2.03	0.40
1:G:299:ARG:HG3	1:G:299:ARG:HH11	1.87	0.40
2:M:145[A]:ASP:O	2:M:147[A]:SER:N	2.54	0.40
1:C:46:LEU:HD13	1:C:51:ILE:HG13	2.03	0.40

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:F:46:LEU:HD11	1:F:99:LEU:HB2	2.04	0.40
1:I:208:GLU:HG3	1:I:210:LEU:N	2.37	0.40
1:G:256:VAL:HB	1:G:269:ILE:HG12	2.03	0.40
1:A:472:ILE:CD1	1:B:336:ILE:HD11	2.51	0.40
1:C:6:ASP:O	1:C:31:LYS:HE3	2.21	0.40
1:G:114:HIS:CD2	1:G:114:HIS:C	2.93	0.40
1:B:385:LYS:HE2	1:B:405:LYS:CD	2.51	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	470/474 (99%)	445 (95%)	23 (5%)	2 (0%)	43	44
1	B	469/474 (99%)	440 (94%)	27 (6%)	2 (0%)	43	44
1	C	470/474 (99%)	455 (97%)	13 (3%)	2 (0%)	43	44
1	D	469/474 (99%)	456 (97%)	11 (2%)	2 (0%)	43	44
1	E	470/474 (99%)	451 (96%)	17 (4%)	2 (0%)	43	44
1	F	469/474 (99%)	452 (96%)	17 (4%)	0	100	100
1	G	470/474 (99%)	456 (97%)	11 (2%)	3 (1%)	33	31
1	H	470/474 (99%)	459 (98%)	10 (2%)	1 (0%)	56	60
1	I	466/474 (98%)	438 (94%)	26 (6%)	2 (0%)	43	44
1	J	470/474 (99%)	452 (96%)	16 (3%)	2 (0%)	43	44
2	K	41/64 (64%)	35 (85%)	6 (15%)	0	100	100
2	L	82/64 (128%)	68 (83%)	6 (7%)	8 (10%)	1	0
2	M	82/64 (128%)	64 (78%)	14 (17%)	4 (5%)	3	1
2	N	82/64 (128%)	60 (73%)	16 (20%)	6 (7%)	2	0
2	O	82/64 (128%)	68 (83%)	8 (10%)	6 (7%)	2	0
All	All	5062/5060 (100%)	4799 (95%)	221 (4%)	42 (1%)	33	23

All (42) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	349	ILE
1	D	79	SER
1	E	348	HIS
1	G	265	LYS
1	H	79	SER
1	J	253	LYS
2	L	154[A]	PRO
2	L	154[B]	PRO
2	M	157[A]	ILE
2	M	157[B]	ILE
2	M	158[A]	PHE
2	M	158[B]	PHE
2	N	148[A]	GLN
2	N	148[B]	GLN
2	O	159[A]	THR
2	O	159[B]	THR
1	C	265	LYS
1	G	266	ALA
2	L	147[A]	SER
2	L	147[B]	SER
2	O	154[A]	PRO
2	O	154[B]	PRO
1	A	79	SER
1	A	291	GLU
1	B	351	TYR
2	N	170[A]	LYS
2	N	170[B]	LYS
2	O	143[A]	SER
2	O	143[B]	SER
1	J	79	SER
1	E	347	VAL
1	I	68	GLY
2	L	158[A]	PHE
2	L	158[B]	PHE
1	D	313	PRO
2	N	154[A]	PRO
2	N	154[B]	PRO
1	C	264	GLY
1	G	313	PRO
2	L	157[A]	ILE
2	L	157[B]	ILE
1	I	313	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	372/373 (100%)	361 (97%)	11 (3%)	53	62
1	B	371/373 (100%)	360 (97%)	11 (3%)	53	62
1	C	372/373 (100%)	358 (96%)	14 (4%)	44	51
1	D	371/373 (100%)	362 (98%)	9 (2%)	61	72
1	E	372/373 (100%)	362 (97%)	10 (3%)	57	66
1	F	371/373 (100%)	360 (97%)	11 (3%)	53	62
1	G	372/373 (100%)	362 (97%)	10 (3%)	57	66
1	H	372/373 (100%)	363 (98%)	9 (2%)	61	72
1	I	372/373 (100%)	361 (97%)	11 (3%)	53	62
1	J	372/373 (100%)	364 (98%)	8 (2%)	64	75
2	K	28/53 (53%)	28 (100%)	0	100	100
2	L	52/53 (98%)	48 (92%)	4 (8%)	18	17
2	M	52/53 (98%)	46 (88%)	6 (12%)	8	6
2	N	52/53 (98%)	50 (96%)	2 (4%)	44	51
2	O	52/53 (98%)	40 (77%)	12 (23%)	1	0
All	All	3953/3995 (99%)	3825 (97%)	128 (3%)	53	60

All (128) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	TYR
1	A	28	LEU
1	A	54	LYS
1	A	141	THR
1	A	232	LYS
1	A	235	PHE
1	A	291	GLU
1	A	321	VAL
1	A	333	ASP
1	A	349	ILE
1	A	467	SER

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Mol	Chain	Res	Type
1	B	19	TYR
1	B	54	LYS
1	B	77	GLU
1	B	80	GLU
1	B	235	PHE
1	B	278	ILE
1	B	290	GLU
1	B	333	ASP
1	B	409	GLN
1	B	410	LYS
1	B	437	GLU
1	C	6	ASP
1	C	19	TYR
1	C	54	LYS
1	C	78	MET
1	C	108	LYS
1	C	114	HIS
1	C	131	LYS
1	C	224	LYS
1	C	235	PHE
1	C	236	LYS
1	C	257	SER
1	C	265	LYS
1	C	269	ILE
1	C	333	ASP
1	D	19	TYR
1	D	54	LYS
1	D	175	LYS
1	D	235	PHE
1	D	257	SER
1	D	306	THR
1	D	311	LYS
1	D	321	VAL
1	D	410	LYS
1	E	3	GLN
1	E	19	TYR
1	E	50	CYS
1	E	54	LYS
1	E	78	MET
1	E	111	LYS
1	E	131	LYS
1	E	136	THR

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Mol	Chain	Res	Type
1	E	235	PHE
1	E	299	ARG
1	F	19	TYR
1	F	54	LYS
1	F	77	GLU
1	F	90	GLU
1	F	235	PHE
1	F	285	LYS
1	F	295	GLU
1	F	304	VAL
1	F	321	VAL
1	F	409	GLN
1	F	410	LYS
1	G	3	GLN
1	G	19	TYR
1	G	54	LYS
1	G	78	MET
1	G	175	LYS
1	G	224	LYS
1	G	235	PHE
1	G	269	ILE
1	G	301	ARG
1	G	333	ASP
1	H	19	TYR
1	H	50	CYS
1	H	54	LYS
1	H	86	ASP
1	H	111	LYS
1	H	235	PHE
1	H	247	THR
1	H	267	GLU
1	H	291	GLU
1	I	19	TYR
1	I	54	LYS
1	I	82	ARG
1	I	114	HIS
1	I	142	LYS
1	I	224	LYS
1	I	235	PHE
1	I	251	ASP
1	I	265	LYS
1	I	290	GLU

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Mol	Chain	Res	Type
1	I	333	ASP
1	J	19	TYR
1	J	54	LYS
1	J	82	ARG
1	J	114	HIS
1	J	142	LYS
1	J	174	LEU
1	J	235	PHE
1	J	301	ARG
2	L	152[A]	THR
2	L	152[B]	THR
2	L	155[A]	ARG
2	L	155[B]	ARG
2	M	130[A]	ARG
2	M	130[B]	ARG
2	M	138[A]	ILE
2	M	138[B]	ILE
2	M	139[A]	LEU
2	M	139[B]	LEU
2	N	155[A]	ARG
2	N	155[B]	ARG
2	O	137[A]	ASN
2	O	137[B]	ASN
2	O	139[A]	LEU
2	O	139[B]	LEU
2	O	145[A]	ASP
2	O	145[B]	ASP
2	O	155[A]	ARG
2	O	155[B]	ARG
2	O	159[A]	THR
2	O	159[B]	THR
2	O	166[A]	LEU
2	O	166[B]	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	3	GLN
1	A	137	GLN
1	A	227	GLN
1	A	231	GLN
1	A	286	ASN

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Mol	Chain	Res	Type
1	B	125	ASN
1	B	231	GLN
1	B	286	ASN
1	B	348	HIS
1	B	373	GLN
1	C	67	HIS
1	C	137	GLN
1	C	225	ASN
1	C	286	ASN
1	C	309	GLN
1	D	67	HIS
1	E	373	GLN
1	F	137	GLN
1	G	114	HIS

5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

49 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	SO4	A	2482	-	4,4,4	1.00	0	6,6,6	0.06	0
3	SO4	A	2483	-	4,4,4	0.98	0	6,6,6	0.08	0
3	SO4	A	2484	-	4,4,4	0.98	0	6,6,6	0.07	0
3	SO4	A	2485	-	4,4,4	1.01	0	6,6,6	0.07	0
4	FAD	A	480	-	58,58,58	1.25	7 (12%)	85,89,89	1.44	9 (10%)
3	SO4	B	2486	-	4,4,4	1.02	0	6,6,6	0.08	0
3	SO4	B	2487	-	4,4,4	1.03	0	6,6,6	0.08	0
3	SO4	B	2488	-	4,4,4	1.00	0	6,6,6	0.08	0
3	SO4	B	2489	-	4,4,4	1.02	0	6,6,6	0.09	0
4	FAD	B	480	-	58,58,58	1.23	8 (13%)	85,89,89	1.38	9 (10%)
3	SO4	C	2490	-	4,4,4	1.01	0	6,6,6	0.09	0
3	SO4	C	2491	-	4,4,4	1.01	0	6,6,6	0.11	0
3	SO4	C	2492	-	4,4,4	1.02	0	6,6,6	0.06	0
3	SO4	C	2493	-	4,4,4	1.03	0	6,6,6	0.07	0
4	FAD	C	480	-	58,58,58	1.15	6 (10%)	85,89,89	1.43	8 (9%)
3	SO4	D	2494	-	4,4,4	1.02	0	6,6,6	0.09	0
3	SO4	D	2495	-	4,4,4	1.00	0	6,6,6	0.10	0
3	SO4	D	2496	-	4,4,4	1.01	0	6,6,6	0.07	0
3	SO4	D	2497	-	4,4,4	1.03	0	6,6,6	0.09	0
4	FAD	D	480	-	58,58,58	1.26	6 (10%)	85,89,89	1.38	9 (10%)
3	SO4	E	2498	-	4,4,4	1.02	0	6,6,6	0.08	0
3	SO4	E	2499	-	4,4,4	1.05	0	6,6,6	0.07	0
3	SO4	E	2500	-	4,4,4	1.02	0	6,6,6	0.10	0
3	SO4	E	2501	-	4,4,4	1.01	0	6,6,6	0.08	0
4	FAD	E	480	-	58,58,58	1.27	7 (12%)	85,89,89	1.33	9 (10%)
3	SO4	F	2502	-	4,4,4	1.01	0	6,6,6	0.08	0
3	SO4	F	2503	-	4,4,4	0.99	0	6,6,6	0.08	0
3	SO4	F	2504	-	4,4,4	1.00	0	6,6,6	0.11	0
3	SO4	F	2505	-	4,4,4	1.03	0	6,6,6	0.08	0
4	FAD	F	480	-	58,58,58	1.34	6 (10%)	85,89,89	1.39	8 (9%)
3	SO4	G	2506	-	4,4,4	1.00	0	6,6,6	0.06	0
3	SO4	G	2507	-	4,4,4	0.99	0	6,6,6	0.08	0
3	SO4	G	2508	-	4,4,4	1.00	0	6,6,6	0.07	0
3	SO4	G	2509	-	4,4,4	1.02	0	6,6,6	0.07	0
4	FAD	G	480	-	58,58,58	1.31	8 (13%)	85,89,89	1.37	8 (9%)
3	SO4	H	2510	-	4,4,4	1.04	0	6,6,6	0.10	0
3	SO4	H	2511	-	4,4,4	1.03	0	6,6,6	0.10	0
3	SO4	H	2512	-	4,4,4	0.98	0	6,6,6	0.11	0
3	SO4	H	2513	-	4,4,4	1.04	0	6,6,6	0.06	0
4	FAD	H	480	-	58,58,58	1.14	6 (10%)	85,89,89	1.45	9 (10%)
3	SO4	I	2514	-	4,4,4	1.00	0	6,6,6	0.07	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	SO4	I	2515	-	4,4,4	1.02	0	6,6,6	0.08	0
3	SO4	I	2516	-	4,4,4	1.02	0	6,6,6	0.05	0
4	FAD	I	480	-	58,58,58	1.26	8 (13%)	85,89,89	1.30	8 (9%)
3	SO4	J	2517	-	4,4,4	1.02	0	6,6,6	0.10	0
3	SO4	J	2518	-	4,4,4	1.01	0	6,6,6	0.07	0
3	SO4	J	2519	-	4,4,4	1.02	0	6,6,6	0.07	0
3	SO4	J	2520	-	4,4,4	1.01	0	6,6,6	0.09	0
4	FAD	J	480	-	58,58,58	1.18	5 (8%)	85,89,89	1.32	8 (9%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	SO4	A	2482	-	-	0/0/0/0	0/0/0/0
3	SO4	A	2483	-	-	0/0/0/0	0/0/0/0
3	SO4	A	2484	-	-	0/0/0/0	0/0/0/0
3	SO4	A	2485	-	-	0/0/0/0	0/0/0/0
4	FAD	A	480	-	-	0/34/50/50	0/1/6/6
3	SO4	B	2486	-	-	0/0/0/0	0/0/0/0
3	SO4	B	2487	-	-	0/0/0/0	0/0/0/0
3	SO4	B	2488	-	-	0/0/0/0	0/0/0/0
3	SO4	B	2489	-	-	0/0/0/0	0/0/0/0
4	FAD	B	480	-	-	0/34/50/50	0/1/6/6
3	SO4	C	2490	-	-	0/0/0/0	0/0/0/0
3	SO4	C	2491	-	-	0/0/0/0	0/0/0/0
3	SO4	C	2492	-	-	0/0/0/0	0/0/0/0
3	SO4	C	2493	-	-	0/0/0/0	0/0/0/0
4	FAD	C	480	-	-	0/34/50/50	0/1/6/6
3	SO4	D	2494	-	-	0/0/0/0	0/0/0/0
3	SO4	D	2495	-	-	0/0/0/0	0/0/0/0
3	SO4	D	2496	-	-	0/0/0/0	0/0/0/0
3	SO4	D	2497	-	-	0/0/0/0	0/0/0/0
4	FAD	D	480	-	-	0/34/50/50	0/1/6/6
3	SO4	E	2498	-	-	0/0/0/0	0/0/0/0
3	SO4	E	2499	-	-	0/0/0/0	0/0/0/0
3	SO4	E	2500	-	-	0/0/0/0	0/0/0/0
3	SO4	E	2501	-	-	0/0/0/0	0/0/0/0
4	FAD	E	480	-	-	0/34/50/50	0/1/6/6
3	SO4	F	2502	-	-	0/0/0/0	0/0/0/0
3	SO4	F	2503	-	-	0/0/0/0	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	SO4	F	2504	-	-	0/0/0/0	0/0/0/0
3	SO4	F	2505	-	-	0/0/0/0	0/0/0/0
4	FAD	F	480	-	-	0/34/50/50	0/1/6/6
3	SO4	G	2506	-	-	0/0/0/0	0/0/0/0
3	SO4	G	2507	-	-	0/0/0/0	0/0/0/0
3	SO4	G	2508	-	-	0/0/0/0	0/0/0/0
3	SO4	G	2509	-	-	0/0/0/0	0/0/0/0
4	FAD	G	480	-	-	0/34/50/50	0/1/6/6
3	SO4	H	2510	-	-	0/0/0/0	0/0/0/0
3	SO4	H	2511	-	-	0/0/0/0	0/0/0/0
3	SO4	H	2512	-	-	0/0/0/0	0/0/0/0
3	SO4	H	2513	-	-	0/0/0/0	0/0/0/0
4	FAD	H	480	-	-	0/34/50/50	0/1/6/6
3	SO4	I	2514	-	-	0/0/0/0	0/0/0/0
3	SO4	I	2515	-	-	0/0/0/0	0/0/0/0
3	SO4	I	2516	-	-	0/0/0/0	0/0/0/0
4	FAD	I	480	-	-	0/34/50/50	0/1/6/6
3	SO4	J	2517	-	-	0/0/0/0	0/0/0/0
3	SO4	J	2518	-	-	0/0/0/0	0/0/0/0
3	SO4	J	2519	-	-	0/0/0/0	0/0/0/0
3	SO4	J	2520	-	-	0/0/0/0	0/0/0/0
4	FAD	J	480	-	-	0/34/50/50	0/1/6/6

All (67) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	F	480	FAD	PA-O3P	4.13	1.67	1.59
4	G	480	FAD	C5X-N5	4.02	1.41	1.35
4	I	480	FAD	C9A-N10	3.55	1.44	1.38
4	J	480	FAD	C5X-N5	3.53	1.40	1.35
4	B	480	FAD	C9A-N10	3.50	1.44	1.38
4	F	480	FAD	C9A-N10	3.44	1.43	1.38
4	G	480	FAD	C9A-N10	3.26	1.43	1.38
4	E	480	FAD	C9A-N10	3.21	1.43	1.38
4	A	480	FAD	C5X-N5	3.06	1.40	1.35
4	D	480	FAD	C5X-N5	3.06	1.40	1.35
4	A	480	FAD	C9A-N10	3.05	1.43	1.38
4	I	480	FAD	C5X-N5	3.02	1.39	1.35
4	J	480	FAD	C5B-C4B	2.94	1.61	1.51
4	C	480	FAD	C5X-N5	2.93	1.39	1.35
4	B	480	FAD	C5X-N5	2.90	1.39	1.35
4	D	480	FAD	C5B-C4B	2.89	1.61	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	J	480	FAD	C1'-N10	2.84	1.51	1.48
4	H	480	FAD	O4B-C1B	2.77	1.45	1.41
4	C	480	FAD	C9A-N10	2.73	1.42	1.38
4	D	480	FAD	C9A-N10	2.73	1.42	1.38
4	H	480	FAD	C1'-C2'	-2.73	1.49	1.51
4	J	480	FAD	C9A-N10	2.69	1.42	1.38
4	F	480	FAD	C5X-N5	2.64	1.39	1.35
4	H	480	FAD	C2A-N3A	2.59	1.37	1.32
4	E	480	FAD	C2A-N1A	2.52	1.38	1.33
4	E	480	FAD	C1'-C2'	-2.50	1.49	1.51
4	E	480	FAD	C2A-N3A	2.44	1.37	1.32
4	A	480	FAD	C2A-N3A	2.42	1.37	1.32
4	B	480	FAD	C2A-N1A	2.39	1.38	1.33
4	I	480	FAD	C1'-N10	2.36	1.50	1.48
4	G	480	FAD	C2A-N3A	2.36	1.36	1.32
4	D	480	FAD	C2A-N3A	2.36	1.36	1.32
4	G	480	FAD	C4X-N5	2.34	1.41	1.36
4	D	480	FAD	C2A-N1A	2.33	1.38	1.33
4	G	480	FAD	C5B-C4B	2.32	1.59	1.51
4	F	480	FAD	C5B-C4B	2.29	1.59	1.51
4	H	480	FAD	C5X-N5	2.28	1.38	1.35
4	F	480	FAD	C2A-N1A	2.28	1.38	1.33
4	E	480	FAD	C5B-C4B	2.24	1.58	1.51
4	G	480	FAD	C1'-C2'	-2.24	1.49	1.51
4	A	480	FAD	C2-N3	2.23	1.41	1.37
4	C	480	FAD	C5B-C4B	2.23	1.58	1.51
4	I	480	FAD	C2A-N3A	2.23	1.36	1.32
4	H	480	FAD	C5B-C4B	2.22	1.58	1.51
4	C	480	FAD	C2A-N3A	2.22	1.36	1.32
4	B	480	FAD	C4X-N5	2.20	1.40	1.36
4	G	480	FAD	C2A-N1A	2.17	1.38	1.33
4	D	480	FAD	C1'-N10	2.17	1.50	1.48
4	F	480	FAD	C10-N1	2.16	1.39	1.35
4	E	480	FAD	C5X-N5	2.16	1.38	1.35
4	I	480	FAD	C2A-N1A	2.14	1.38	1.33
4	B	480	FAD	C2-N3	2.13	1.41	1.37
4	A	480	FAD	C1'-N10	2.12	1.50	1.48
4	A	480	FAD	C2A-N1A	2.12	1.38	1.33
4	J	480	FAD	C2A-N1A	2.11	1.38	1.33
4	I	480	FAD	C2-N3	2.10	1.41	1.37
4	I	480	FAD	C5B-C4B	2.09	1.58	1.51
4	B	480	FAD	C2A-N3A	2.09	1.36	1.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	480	FAD	PA-O3P	2.08	1.63	1.59
4	C	480	FAD	C2A-N1A	2.07	1.38	1.33
4	I	480	FAD	C4X-N5	2.06	1.40	1.36
4	B	480	FAD	C4X-C10	2.03	1.44	1.40
4	H	480	FAD	P-O2P	-2.03	1.46	1.55
4	C	480	FAD	C2-N3	2.02	1.41	1.37
4	E	480	FAD	C1'-N10	2.01	1.50	1.48
4	G	480	FAD	C6-C5X	-2.01	1.39	1.41
4	A	480	FAD	C6-C7	2.01	1.43	1.37

All (85) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	480	FAD	C2-N1-C10	5.79	120.82	114.98
4	H	480	FAD	C2-N1-C10	5.23	120.25	114.98
4	B	480	FAD	C4B-O4B-C1B	-5.19	104.12	109.75
4	G	480	FAD	C4B-O4B-C1B	-5.16	104.14	109.75
4	A	480	FAD	C4B-O4B-C1B	-5.15	104.15	109.75
4	A	480	FAD	C2-N1-C10	4.92	119.94	114.98
4	G	480	FAD	C2-N1-C10	4.83	119.85	114.98
4	I	480	FAD	C2-N1-C10	4.78	119.79	114.98
4	J	480	FAD	C2-N1-C10	4.77	119.78	114.98
4	B	480	FAD	C2-N1-C10	4.74	119.75	114.98
4	E	480	FAD	C2-N1-C10	4.71	119.73	114.98
4	C	480	FAD	N3A-C2A-N1A	-4.67	124.80	128.71
4	D	480	FAD	C2-N1-C10	4.67	119.69	114.98
4	D	480	FAD	N3A-C2A-N1A	-4.59	124.87	128.71
4	E	480	FAD	N3A-C2A-N1A	-4.56	124.90	128.71
4	F	480	FAD	N3A-C2A-N1A	-4.38	125.05	128.71
4	H	480	FAD	C4B-O4B-C1B	-4.37	105.00	109.75
4	F	480	FAD	C2-N1-C10	4.35	119.36	114.98
4	H	480	FAD	N3A-C2A-N1A	-4.30	125.11	128.71
4	G	480	FAD	N3A-C2A-N1A	-4.30	125.11	128.71
4	J	480	FAD	C4B-O4B-C1B	-4.27	105.12	109.75
4	C	480	FAD	C4B-O4B-C1B	-4.26	105.12	109.75
4	F	480	FAD	C4B-O4B-C1B	-4.18	105.21	109.75
4	I	480	FAD	N3A-C2A-N1A	-4.16	125.23	128.71
4	C	480	FAD	C4X-N5-C5X	4.14	121.34	116.69
4	H	480	FAD	C4X-N5-C5X	4.14	121.34	116.69
4	A	480	FAD	N3A-C2A-N1A	-4.08	125.30	128.71
4	B	480	FAD	O4B-C1B-C2B	-4.07	100.54	106.77
4	A	480	FAD	C4X-N5-C5X	3.99	121.18	116.69

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	F	480	FAD	C4X-N5-C5X	3.97	121.16	116.69
4	J	480	FAD	N3A-C2A-N1A	-3.97	125.39	128.71
4	E	480	FAD	C4X-N5-C5X	3.92	121.09	116.69
4	B	480	FAD	C4X-N5-C5X	3.88	121.06	116.69
4	B	480	FAD	N3A-C2A-N1A	-3.86	125.48	128.71
4	I	480	FAD	O4B-C1B-C2B	-3.85	100.87	106.77
4	G	480	FAD	O4B-C1B-C2B	-3.78	100.98	106.77
4	D	480	FAD	C4B-O4B-C1B	-3.73	105.70	109.75
4	D	480	FAD	C4X-N5-C5X	3.73	120.88	116.69
4	E	480	FAD	O4B-C1B-C2B	-3.67	101.14	106.77
4	A	480	FAD	O4B-C1B-C2B	-3.61	101.24	106.77
4	I	480	FAD	C4X-N5-C5X	3.55	120.68	116.69
4	F	480	FAD	C4X-C10-N10	-3.55	118.74	120.51
4	F	480	FAD	O4B-C1B-C2B	-3.52	101.38	106.77
4	J	480	FAD	C4X-N5-C5X	3.43	120.54	116.69
4	G	480	FAD	C4X-N5-C5X	3.40	120.51	116.69
4	H	480	FAD	O4B-C1B-C2B	-3.39	101.58	106.77
4	D	480	FAD	O4B-C1B-C2B	-3.34	101.65	106.77
4	A	480	FAD	C4X-C10-N10	-3.30	118.86	120.51
4	H	480	FAD	C4X-C10-N10	-3.23	118.90	120.51
4	E	480	FAD	C4B-O4B-C1B	-3.15	106.33	109.75
4	E	480	FAD	C4X-C10-N10	-3.12	118.95	120.51
4	C	480	FAD	O4B-C1B-C2B	-3.08	102.05	106.77
4	D	480	FAD	P-O3P-PA	2.99	140.44	131.68
4	I	480	FAD	C1'-N10-C9A	2.97	121.76	118.87
4	J	480	FAD	O4B-C1B-C2B	-2.93	102.27	106.77
4	C	480	FAD	C1'-N10-C9A	2.90	121.69	118.87
4	B	480	FAD	C1'-N10-C9A	2.70	121.50	118.87
4	D	480	FAD	C4X-C10-N10	-2.60	119.21	120.51
4	A	480	FAD	P-O3P-PA	2.57	139.23	131.68
4	A	480	FAD	C5X-C9A-N10	2.57	119.33	116.80
4	H	480	FAD	C5X-C9A-N10	2.56	119.32	116.80
4	D	480	FAD	C1'-N10-C9A	2.47	121.27	118.87
4	D	480	FAD	C5X-C9A-N10	2.43	119.20	116.80
4	B	480	FAD	P-O3P-PA	2.43	138.80	131.68
4	B	480	FAD	C4X-C10-N10	-2.40	119.31	120.51
4	C	480	FAD	P-O3P-PA	2.40	138.71	131.68
4	I	480	FAD	C4B-O4B-C1B	-2.38	107.17	109.75
4	E	480	FAD	C5X-C9A-N10	2.30	119.06	116.80
4	B	480	FAD	C5X-C9A-N10	2.30	119.06	116.80
4	J	480	FAD	C1'-N10-C9A	2.26	121.07	118.87
4	J	480	FAD	C5X-C9A-N10	2.24	119.01	116.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	E	480	FAD	C1'-N10-C9A	2.21	121.02	118.87
4	G	480	FAD	C1'-N10-C9A	2.20	121.02	118.87
4	G	480	FAD	N3A-C4A-N9A	2.19	129.39	125.43
4	J	480	FAD	C4X-C10-N10	-2.14	119.44	120.51
4	G	480	FAD	C5X-C9A-N10	2.11	118.88	116.80
4	F	480	FAD	C5X-C9A-N10	2.10	118.87	116.80
4	F	480	FAD	C1'-N10-C9A	2.09	120.91	118.87
4	C	480	FAD	C5X-C9A-N10	2.08	118.86	116.80
4	I	480	FAD	C5X-C9A-N10	2.05	118.82	116.80
4	E	480	FAD	N3A-C4A-N9A	2.04	129.11	125.43
4	A	480	FAD	C1'-N10-C9A	2.03	120.85	118.87
4	I	480	FAD	P-O3P-PA	2.02	137.62	131.68
4	H	480	FAD	O2B-C2B-C3B	2.02	118.39	111.83
4	H	480	FAD	P-O3P-PA	2.00	137.56	131.68

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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	472/474 (99%)	0.00	14 (2%) 48 50	26, 44, 68, 86	0
1	B	471/474 (99%)	0.53	49 (10%) 7 7	28, 51, 84, 93	0
1	C	472/474 (99%)	-0.09	9 (1%) 64 65	23, 36, 61, 79	0
1	D	471/474 (99%)	-0.14	6 (1%) 74 75	21, 31, 53, 74	0
1	E	472/474 (99%)	0.03	13 (2%) 50 52	24, 37, 60, 81	0
1	F	471/474 (99%)	-0.13	7 (1%) 70 72	20, 32, 54, 73	0
1	G	472/474 (99%)	-0.13	7 (1%) 70 72	21, 36, 60, 79	0
1	H	472/474 (99%)	-0.25	2 (0%) 90 92	16, 27, 48, 66	0
1	I	470/474 (99%)	0.33	29 (6%) 20 21	27, 49, 76, 86	0
1	J	472/474 (99%)	-0.10	9 (1%) 64 65	25, 38, 60, 82	0
2	K	43/64 (67%)	1.65	14 (32%) 1 1	54, 68, 81, 92	0
2	L	43/64 (67%)	3.83	28 (65%) 0 0	35, 47, 58, 60	0
2	M	43/64 (67%)	4.09	36 (83%) 0 0	33, 49, 56, 58	0
2	N	43/64 (67%)	4.61	39 (90%) 0 0	38, 51, 60, 61	0
2	O	43/64 (67%)	4.94	36 (83%) 0 0	36, 50, 58, 59	0
All	All	4930/5060 (97%)	0.17	298 (6%) 21 22	16, 38, 69, 93	0

All (298) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	N	172[A]	THR	12.9
2	O	167[A]	VAL	12.8
2	O	164[A]	LEU	12.4
2	O	144[A]	LEU	12.4
2	L	144[A]	LEU	11.9
2	O	166[A]	LEU	11.2
2	L	166[A]	LEU	10.0

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Mol	Chain	Res	Type	RSRZ
2	N	171[A]	GLN	10.0
2	L	167[A]	VAL	9.4
2	N	166[A]	LEU	9.2
2	N	167[A]	VAL	8.9
2	O	149[A]	GLY	8.6
2	L	146[A]	ALA	8.4
2	M	144[A]	LEU	8.4
2	O	139[A]	LEU	8.4
2	M	146[A]	ALA	7.6
2	L	171[A]	GLN	7.4
2	O	169[A]	LEU	7.4
2	O	150[A]	THR	7.2
2	L	170[A]	LYS	7.2
2	M	157[A]	ILE	7.1
2	N	163[A]	ALA	7.1
1	B	294	ILE	7.1
2	N	169[A]	LEU	7.0
2	O	145[A]	ASP	6.9
1	B	312	ILE	6.7
2	L	172[A]	THR	6.7
2	N	150[A]	THR	6.6
2	N	157[A]	ILE	6.6
2	N	145[A]	ASP	6.6
2	M	164[A]	LEU	6.5
2	N	158[A]	PHE	6.4
2	M	167[A]	VAL	6.4
2	M	149[A]	GLY	6.3
2	N	138[A]	ILE	6.3
2	L	169[A]	LEU	6.2
2	O	138[A]	ILE	6.2
1	G	265	LYS	6.2
2	K	172	THR	6.0
2	N	139[A]	LEU	6.0
2	M	145[A]	ASP	5.9
2	N	146[A]	ALA	5.9
2	O	158[A]	PHE	5.9
2	M	172[A]	THR	5.9
2	M	166[A]	LEU	5.9
2	O	148[A]	GLN	5.8
2	O	143[A]	SER	5.8
2	M	150[A]	THR	5.8
2	O	171[A]	GLN	5.8

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Mol	Chain	Res	Type	RSRZ
2	L	149[A]	GLY	5.8
2	M	148[A]	GLN	5.7
2	O	165[A]	LYS	5.6
2	M	169[A]	LEU	5.6
2	O	142[A]	HIS	5.5
2	M	156[A]	GLY	5.5
1	B	289	LEU	5.5
2	M	158[A]	PHE	5.4
2	O	163[A]	ALA	5.4
2	M	142[A]	HIS	5.4
1	B	139	ILE	5.3
2	O	168[A]	GLN	5.3
2	L	139[A]	LEU	5.3
2	L	145[A]	ASP	5.3
2	M	139[A]	LEU	5.3
2	M	171[A]	GLN	5.2
2	L	163[A]	ALA	5.2
2	N	151[A]	ALA	5.2
2	M	153[A]	GLY	5.1
2	O	170[A]	LYS	5.1
2	N	147[A]	SER	5.1
2	O	172[A]	THR	5.1
2	M	143[A]	SER	5.1
2	O	147[A]	SER	5.1
1	B	121	ILE	5.0
2	N	144[A]	LEU	4.9
2	N	149[A]	GLY	4.9
2	L	148[A]	GLN	4.9
1	I	254	ILE	4.9
2	M	147[A]	SER	4.8
2	M	138[A]	ILE	4.8
2	L	142[A]	HIS	4.8
2	L	164[A]	LEU	4.7
2	L	147[A]	SER	4.7
1	I	132	ALA	4.6
2	N	155[A]	ARG	4.6
2	K	149	GLY	4.6
2	N	164[A]	LEU	4.5
2	M	163[A]	ALA	4.5
2	O	159[A]	THR	4.5
1	B	313	PRO	4.4
2	L	151[A]	ALA	4.4

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Mol	Chain	Res	Type	RSRZ
2	K	152	THR	4.4
1	B	79	SER	4.3
1	C	79	SER	4.3
2	M	154[A]	PRO	4.3
1	B	290	GLU	4.2
2	O	146[A]	ALA	4.2
2	O	151[A]	ALA	4.2
2	M	151[A]	ALA	4.2
1	E	135	GLY	4.1
1	B	311	LYS	4.1
2	N	130[A]	ARG	4.1
2	N	152[A]	THR	4.1
1	B	123	GLY	4.1
1	B	315	ILE	4.1
2	L	165[A]	LYS	4.1
2	L	168[A]	GLN	4.1
1	B	129	ALA	4.0
1	I	79	SER	3.9
2	M	168[A]	GLN	3.9
2	N	162[A]	ASP	3.9
2	N	148[A]	GLN	3.9
2	O	154[A]	PRO	3.9
1	B	5	ILE	3.9
2	O	157[A]	ILE	3.9
1	B	287	LEU	3.9
2	O	161[A]	GLU	3.8
1	B	132	ALA	3.8
2	K	146	ALA	3.8
2	N	161[A]	GLU	3.8
1	B	292	LEU	3.7
2	N	170[A]	LYS	3.7
1	B	348	HIS	3.7
2	N	154[A]	PRO	3.7
2	K	150	THR	3.7
1	I	269	ILE	3.7
1	I	160	ILE	3.7
2	L	138[A]	ILE	3.6
1	B	113	VAL	3.6
1	I	136	THR	3.6
1	I	268	VAL	3.5
1	F	135	GLY	3.5
1	B	138	VAL	3.5

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Mol	Chain	Res	Type	RSRZ
2	M	159[A]	THR	3.4
2	O	134[A]	ALA	3.4
2	N	133[A]	PRO	3.4
2	O	140[A]	GLU	3.4
1	B	265	LYS	3.3
1	I	258	ILE	3.3
1	I	262	SER	3.3
2	N	143[A]	SER	3.3
1	B	288	GLY	3.3
2	N	136[A]	ARG	3.3
2	O	153[A]	GLY	3.3
1	B	295	GLU	3.3
1	B	133	ASP	3.3
2	N	131[A]	LEU	3.2
1	I	176	LYS	3.2
1	I	80	GLU	3.2
1	B	122	THR	3.2
2	L	150[A]	THR	3.2
1	I	266	ALA	3.2
1	C	265	LYS	3.2
1	J	80	GLU	3.1
1	A	451	ALA	3.1
2	L	143[A]	SER	3.1
2	N	168[A]	GLN	3.1
2	K	166	LEU	3.1
2	M	134[A]	ALA	3.1
1	J	265	LYS	3.1
2	M	170[A]	LYS	3.1
1	C	132	ALA	3.1
1	A	79	SER	3.1
1	E	262	SER	3.1
1	B	310	THR	3.1
1	D	429	VAL	3.1
1	B	381	TYR	3.0
1	E	261	ALA	3.0
1	A	132	ALA	3.0
1	I	249	LYS	3.0
2	K	148	GLN	3.0
1	B	115	VAL	3.0
1	A	3	GLN	2.9
2	N	134[A]	ALA	2.9
2	N	135[A]	ALA	2.9

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Mol	Chain	Res	Type	RSRZ
1	E	454	THR	2.9
1	E	132	ALA	2.9
2	N	165[A]	LYS	2.9
1	E	79	SER	2.9
2	N	153[A]	GLY	2.8
1	J	79	SER	2.8
2	M	155[A]	ARG	2.8
1	B	425	ALA	2.8
2	O	135[A]	ALA	2.8
1	A	133	ASP	2.8
1	B	344	GLY	2.8
1	E	455	LEU	2.8
1	B	144	ILE	2.8
1	D	265	LYS	2.8
2	L	158[A]	PHE	2.8
1	B	124	LYS	2.8
2	L	141[A]	LYS	2.7
2	K	167	VAL	2.7
2	K	142	HIS	2.7
1	B	376	GLU	2.7
1	G	3	GLN	2.7
1	A	290	GLU	2.7
1	B	135	GLY	2.7
1	I	131	LYS	2.6
2	L	161[A]	GLU	2.6
2	O	141[A]	LYS	2.6
1	I	82	ARG	2.6
1	B	291	GLU	2.6
1	J	78	MET	2.6
1	G	79	SER	2.6
2	O	136[A]	ARG	2.6
1	I	276	VAL	2.6
2	L	130[A]	ARG	2.6
1	A	124	LYS	2.6
2	K	147	SER	2.5
1	I	133	ASP	2.5
2	N	159[A]	THR	2.5
1	A	139	ILE	2.5
1	J	262	SER	2.5
1	A	447	ARG	2.5
2	L	133[A]	PRO	2.5
1	A	134	GLY	2.5

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Mol	Chain	Res	Type	RSRZ
1	J	263	GLY	2.5
1	D	79	SER	2.5
1	B	296	LEU	2.5
1	E	428	MET	2.5
1	I	135	GLY	2.4
1	B	346	ALA	2.4
1	B	131	LYS	2.4
1	B	130	THR	2.4
1	E	429	VAL	2.4
2	K	168	GLN	2.4
1	B	343	ALA	2.4
1	E	432	ALA	2.4
2	O	131[A]	LEU	2.4
1	I	244	THR	2.4
1	I	228	ARG	2.4
1	J	449	CYS	2.3
1	F	429	VAL	2.3
2	K	163	ALA	2.3
1	G	5	ILE	2.3
1	G	454	THR	2.3
1	C	268	VAL	2.3
1	J	252	GLY	2.3
2	K	151	ALA	2.3
1	A	380	GLU	2.3
1	I	243	VAL	2.3
1	C	134	GLY	2.3
1	F	131	LYS	2.2
2	N	160[A]	LYS	2.2
2	O	130[A]	ARG	2.2
2	N	156[A]	GLY	2.2
1	I	274	LEU	2.2
1	B	33	VAL	2.2
1	B	347	VAL	2.2
1	F	132	ALA	2.2
1	A	292	LEU	2.2
1	C	455	LEU	2.2
2	N	142[A]	HIS	2.2
1	D	426	GLY	2.2
1	G	132	ALA	2.2
1	F	428	MET	2.2
2	M	133[A]	PRO	2.2
2	K	165	LYS	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	136	THR	2.2
1	H	50	CYS	2.2
1	H	454	THR	2.2
2	M	152[A]	THR	2.2
1	G	124	LYS	2.2
1	I	252	GLY	2.2
2	O	155[A]	ARG	2.2
1	A	411	SER	2.2
1	E	3	GLN	2.2
1	C	451	ALA	2.2
1	B	126	GLN	2.1
2	M	130[A]	ARG	2.1
1	I	69	THR	2.1
1	B	429	VAL	2.1
2	M	161[A]	GLU	2.1
1	E	420	ILE	2.1
2	L	134[A]	ALA	2.1
1	I	247	THR	2.1
2	M	165[A]	LYS	2.1
1	E	134	GLY	2.1
1	F	80	GLU	2.1
1	D	69	THR	2.1
1	B	128	THR	2.1
1	I	175	LYS	2.1
1	C	269	ILE	2.1
1	C	78	MET	2.1
1	I	255	ASP	2.1
1	A	443	GLU	2.0
1	I	425	ALA	2.0
1	J	77	GLU	2.0
2	M	160[A]	LYS	2.0
1	B	6	ASP	2.0
1	B	454	THR	2.0
1	D	432	ALA	2.0
1	F	430	ASN	2.0
2	M	131[A]	LEU	2.0
1	B	50	CYS	2.0
1	I	5	ILE	2.0
1	B	45	CYS	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	SO4	F	2504	5/5	0.23	12.97	76,76,80,80	0
3	SO4	J	2519	5/5	0.22	12.53	86,86,88,89	0
3	SO4	H	2512	5/5	0.13	12.08	58,60,64,69	0
3	SO4	A	2483	5/5	0.22	9.24	88,92,93,95	0
3	SO4	I	2515	5/5	0.20	7.33	94,94,97,97	0
3	SO4	H	2511	5/5	0.21	5.48	75,79,83,83	0
3	SO4	E	2499	5/5	0.26	5.41	85,87,89,90	0
3	SO4	C	2490	5/5	0.15	4.06	73,77,81,82	0
3	SO4	G	2508	5/5	0.17	3.87	84,85,88,89	0
3	SO4	B	2488	5/5	0.26	3.74	92,93,94,97	0
3	SO4	F	2503	5/5	0.24	3.36	80,81,83,84	0
3	SO4	C	2491	5/5	0.17	3.18	86,89,92,92	0
3	SO4	E	2500	5/5	0.17	2.39	83,83,85,87	0
3	SO4	B	2487	5/5	0.20	2.33	89,91,92,92	0
3	SO4	D	2496	5/5	0.12	2.28	74,74,76,79	0
3	SO4	J	2518	5/5	0.22	1.88	93,95,96,96	0
3	SO4	H	2510	5/5	0.16	1.72	93,93,95,96	0
3	SO4	A	2484	5/5	0.10	1.51	77,78,79,81	0
3	SO4	G	2507	5/5	0.14	1.50	87,87,89,91	0
3	SO4	D	2495	5/5	0.21	1.07	83,86,87,87	0
3	SO4	I	2514	5/5	0.15	0.93	72,74,78,80	0
3	SO4	C	2492	5/5	0.12	0.80	86,88,88,88	0
4	FAD	J	480	53/53	0.13	0.75	23,30,37,39	0
4	FAD	C	480	53/53	0.13	0.73	25,32,39,43	0
3	SO4	E	2498	5/5	0.12	0.61	77,81,81,85	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
4	FAD	H	480	53/53	0.15	0.55	14,21,28,38	0
4	FAD	I	480	53/53	0.16	0.46	31,44,56,57	0
3	SO4	C	2493	5/5	0.15	0.46	80,81,84,86	0
3	SO4	A	2482	5/5	0.13	0.40	74,79,83,83	0
3	SO4	J	2517	5/5	0.13	0.39	65,68,69,70	0
4	FAD	D	480	53/53	0.13	0.38	21,27,32,36	0
3	SO4	A	2485	5/5	0.23	0.28	91,93,94,98	0
4	FAD	E	480	53/53	0.12	0.24	26,35,41,43	0
4	FAD	F	480	53/53	0.13	0.23	21,29,38,39	0
4	FAD	G	480	53/53	0.12	0.18	21,31,46,51	0
4	FAD	A	480	53/53	0.13	0.15	28,39,57,58	0
3	SO4	B	2489	5/5	0.33	0.11	100,101,102,106	0
3	SO4	G	2506	5/5	0.11	0.01	79,80,81,81	0
4	FAD	B	480	53/53	0.15	-0.06	42,55,68,70	0
3	SO4	D	2497	5/5	0.11	-0.40	58,62,66,67	0
3	SO4	F	2502	5/5	0.10	-0.51	54,59,65,66	0
3	SO4	H	2513	5/5	0.09	-0.52	60,61,67,70	0
3	SO4	I	2516	5/5	0.15	-0.54	84,86,90,91	0
3	SO4	B	2486	5/5	0.11	-0.56	87,90,90,92	0
3	SO4	J	2520	5/5	0.10	-0.58	72,73,76,76	0
3	SO4	G	2509	5/5	0.12	-0.76	88,88,91,94	0
3	SO4	D	2494	5/5	0.12	-0.78	61,64,65,68	0
3	SO4	F	2505	5/5	0.08	-0.97	64,65,71,72	0
3	SO4	E	2501	5/5	0.08	-1.08	79,81,81,84	0

6.5 Other polymers ⓘ

There are no such residues in this entry.