



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 03:29 AM UTC

PDB ID : 1DI0 / pdb_00001di0
Title : CRYSTAL STRUCTURE OF LUMAZINE SYNTHASE FROM BRUCELLA ABORTUS
Authors : Braden, B.C.; Velikovsky, C.A.; Cauwerhff, A.A.; Polikarpov, I.; Goldbaum, F.A.
Deposited on : 1999-11-28
Resolution : 2.70 Å(reported)

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

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A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

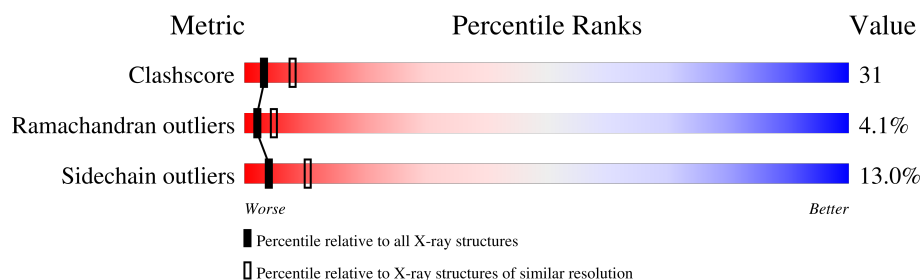
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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))
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)

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

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	158	
1	B	158	
1	C	158	
1	D	158	
1	E	158	

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
2	PO4	A	190	-	-	X	-
2	PO4	A	192	-	-	X	-
2	PO4	B	195	-	-	X	-
2	PO4	C	196	-	-	X	-
2	PO4	C	197	-	-	X	-
2	PO4	D	198	-	-	X	-
2	PO4	D	199	-	-	X	-
2	PO4	E	202	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5818 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 LUMAZINE SYNTHASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	148	Total	C	N	O	S	0	0	0
			1147	738	201	206	2			
1	B	148	Total	C	N	O	S	0	0	0
			1147	738	201	206	2			
1	C	146	Total	C	N	O	S	0	0	0
			1135	731	199	203	2			
1	D	147	Total	C	N	O	S	0	0	0
			1142	735	200	205	2			
1	E	146	Total	C	N	O	S	0	0	0
			1135	731	199	203	2			

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	88	ASP	ARG	conflict	UNP P61711
B	88	ASP	ARG	conflict	UNP P61711
C	88	ASP	ARG	conflict	UNP P61711
D	88	ASP	ARG	conflict	UNP P61711
E	88	ASP	ARG	conflict	UNP P61711

- Molecule 2 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	P	0	0
			5	4	1		
2	A	1	Total	O	P	0	0
			5	4	1		
2	A	1	Total	O	P	0	0
			5	4	1		
2	B	1	Total	O	P	0	0
			5	4	1		
2	B	1	Total	O	P	0	0
			5	4	1		
2	B	1	Total	O	P	0	0
			5	4	1		
2	C	1	Total	O	P	0	0
			5	4	1		
2	C	1	Total	O	P	0	0
			5	4	1		
2	D	1	Total	O	P	0	0
			5	4	1		
2	D	1	Total	O	P	0	0
			5	4	1		
2	E	1	Total	O	P	0	0
			5	4	1		
2	E	1	Total	O	P	0	0
			5	4	1		
2	E	1	Total	O	P	0	0
			5	4	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	10	Total 10	O 10	0	0
3	B	8	Total 8	O 8	0	0
3	C	12	Total 12	O 12	0	0
3	D	7	Total 7	O 7	0	0
3	E	10	Total 10	O 10	0	0

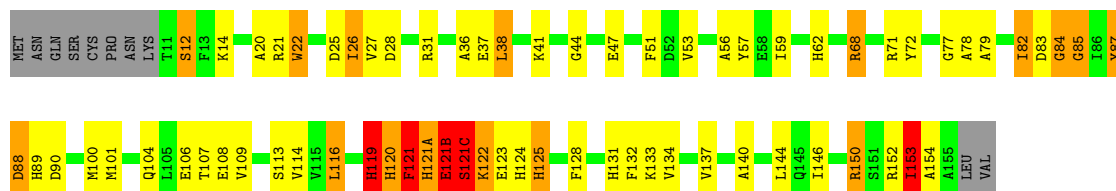
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. 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. 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.

Note EDS was not executed.

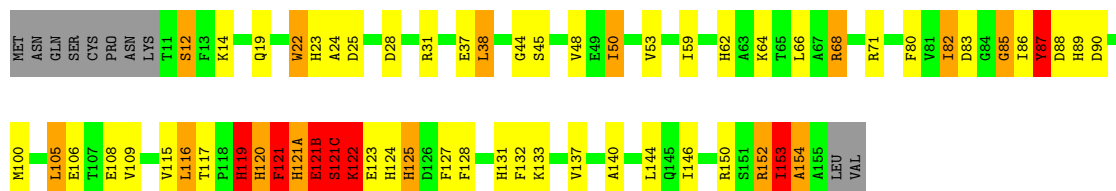
• Molecule 1: LUMAZINE SYNTHASE

Chain A: 

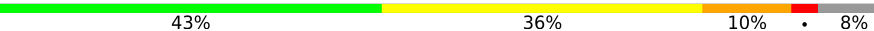


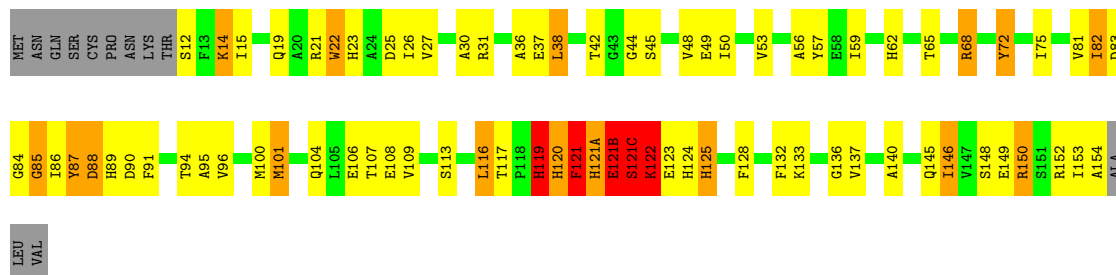
• Molecule 1: LUMAZINE SYNTHASE

Chain B: 

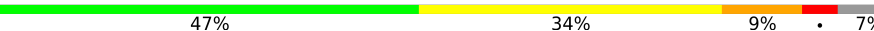


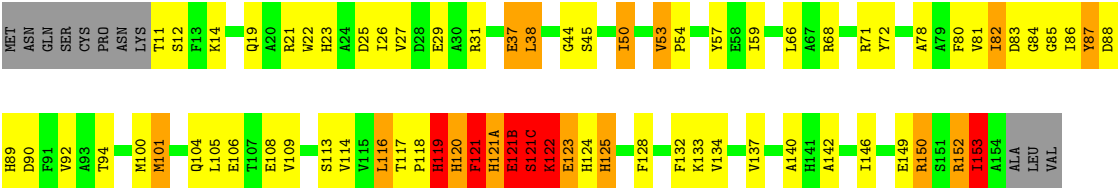
• Molecule 1: LUMAZINE SYNTHASE

Chain C: 

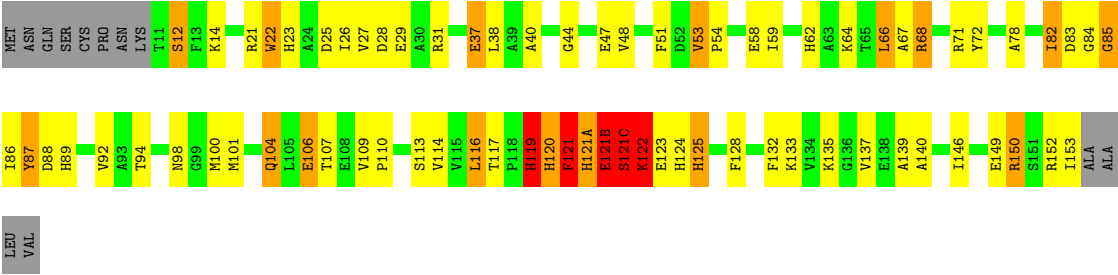


• Molecule 1: LUMAZINE SYNTHASE

Chain D: 



● Molecule 1: LUMAZINE SYNTHASE



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	142.25Å 142.25Å 242.30Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	150.00 – 2.70	Depositor
% Data completeness (in resolution range)	90.0 (150.00-2.70)	Depositor
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS	Depositor
R, R_{free}	0.180 , 0.230	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5818	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PO4

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	1.73	11/1176 (0.9%)	1.60	20/1596 (1.3%)
1	B	1.71	9/1176 (0.8%)	1.58	16/1596 (1.0%)
1	C	1.86	19/1164 (1.6%)	1.57	18/1579 (1.1%)
1	D	1.84	17/1171 (1.5%)	1.59	16/1589 (1.0%)
1	E	1.73	17/1164 (1.5%)	1.54	18/1579 (1.1%)
All	All	1.77	73/5851 (1.2%)	1.58	88/7939 (1.1%)

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
1	B	0	1
1	C	0	1
All	All	0	2

All (73) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	100	MET	SD-CE	-12.83	1.47	1.79
1	D	100	MET	SD-CE	-12.74	1.47	1.79
1	C	100	MET	SD-CE	-12.55	1.48	1.79
1	C	59	ILE	CA-CB	-10.93	1.48	1.54
1	C	36	ALA	CA-CB	-8.86	1.39	1.53
1	E	59	ILE	C-O	-8.56	1.13	1.24
1	C	48	VAL	CA-CB	-8.37	1.44	1.54
1	E	100	MET	SD-CE	-7.89	1.59	1.79
1	C	146	ILE	CA-CB	7.78	1.64	1.54
1	D	50	ILE	C-O	-7.64	1.13	1.24
1	C	15	ILE	C-O	-7.05	1.16	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	153	ILE	CA-C	-6.93	1.46	1.53
1	E	12	SER	CA-C	6.92	1.60	1.52
1	C	30	ALA	CA-CB	-6.75	1.42	1.53
1	C	101	MET	CG-SD	-6.66	1.64	1.80
1	D	106	GLU	C-O	-6.61	1.15	1.24
1	D	59	ILE	CA-CB	-6.42	1.50	1.54
1	A	134	VAL	CA-CB	-6.42	1.47	1.54
1	A	20	ALA	CA-CB	-6.41	1.42	1.53
1	E	106	GLU	C-O	-6.33	1.16	1.24
1	B	100	MET	SD-CE	-6.28	1.63	1.79
1	E	67	ALA	CA-CB	-6.18	1.43	1.53
1	D	142	ALA	CA-CB	-6.15	1.43	1.53
1	E	66	LEU	C-O	-6.12	1.17	1.24
1	C	42	THR	CA-CB	6.07	1.62	1.54
1	A	121(B)	GLU	CA-C	6.00	1.60	1.52
1	D	94	THR	CA-CB	-5.96	1.43	1.53
1	E	140	ALA	CA-CB	-5.90	1.44	1.53
1	B	109	VAL	CA-CB	-5.89	1.47	1.54
1	A	26	ILE	CA-CB	-5.88	1.47	1.54
1	C	95	ALA	CA-CB	-5.82	1.44	1.53
1	B	105	LEU	N-CA	-5.81	1.38	1.46
1	B	59	ILE	C-O	-5.80	1.19	1.24
1	A	72	TYR	CA-C	-5.71	1.45	1.52
1	C	136	GLY	C-O	-5.68	1.17	1.23
1	D	66	LEU	C-O	-5.68	1.17	1.24
1	A	36	ALA	CA-CB	-5.68	1.44	1.53
1	E	139	ALA	CA-CB	-5.63	1.44	1.53
1	D	80	PHE	C-O	-5.57	1.17	1.24
1	A	84	GLY	C-O	5.53	1.30	1.24
1	C	106	GLU	CA-C	-5.46	1.45	1.52
1	C	91	PHE	C-O	-5.44	1.17	1.24
1	B	108	GLU	CA-CB	-5.43	1.45	1.53
1	C	65	THR	N-CA	-5.42	1.40	1.46
1	C	148	SER	N-CA	-5.39	1.39	1.46
1	E	48	VAL	CA-CB	-5.37	1.47	1.54
1	B	115	VAL	CA-C	-5.37	1.46	1.52
1	D	37	GLU	N-CA	-5.36	1.39	1.46
1	C	75	ILE	CA-CB	-5.35	1.47	1.54
1	E	71	ARG	C-O	-5.34	1.17	1.24
1	E	37	GLU	C-O	-5.34	1.17	1.24
1	A	79	ALA	C-O	-5.30	1.17	1.23
1	D	80	PHE	CA-C	-5.29	1.46	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	47	GLU	CA-C	5.29	1.59	1.52
1	E	104	GLN	N-CA	-5.29	1.39	1.46
1	D	134	VAL	C-O	5.28	1.30	1.24
1	A	59	ILE	C-O	-5.26	1.17	1.24
1	B	50	ILE	C-O	-5.25	1.18	1.24
1	D	101	MET	CG-SD	-5.23	1.67	1.80
1	E	107	THR	CB-CG2	-5.22	1.35	1.52
1	B	117	THR	C-O	-5.21	1.20	1.24
1	E	109	VAL	CA-CB	-5.19	1.50	1.54
1	E	40	ALA	CA-CB	-5.14	1.45	1.53
1	E	117	THR	C-O	-5.14	1.21	1.25
1	E	72	TYR	CA-C	-5.11	1.46	1.52
1	D	53	VAL	CA-CB	5.11	1.60	1.54
1	C	145	GLN	CA-C	5.11	1.59	1.52
1	D	153	ILE	CG1-CD1	5.10	1.71	1.51
1	C	75	ILE	C-O	-5.07	1.18	1.24
1	C	109	VAL	CA-CB	-5.07	1.48	1.54
1	D	121(B)	GLU	CA-C	5.06	1.59	1.52
1	D	68	ARG	NE-CZ	5.05	1.38	1.33
1	B	59	ILE	CA-CB	-5.04	1.51	1.54

All (88) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	119	HIS	N-CA-C	12.01	125.52	111.71
1	D	119	HIS	N-CA-C	11.08	124.45	111.71
1	C	119	HIS	N-CA-C	10.48	123.77	111.71
1	A	119	HIS	N-CA-C	10.18	123.42	111.71
1	E	119	HIS	N-CA-C	9.67	122.83	111.71
1	E	109	VAL	N-CA-C	-9.00	96.91	107.61
1	A	121	PHE	N-CA-C	-8.77	99.22	110.53
1	D	153	ILE	N-CA-C	8.63	120.22	106.32
1	A	150	ARG	NE-CZ-NH1	-8.09	113.41	121.50
1	B	121	PHE	N-CA-C	-7.99	100.58	110.41
1	D	121	PHE	N-CA-C	-7.57	101.10	110.41
1	B	85	GLY	N-CA-C	-7.54	100.73	110.38
1	C	121	PHE	N-CA-C	-7.27	101.15	110.53
1	A	85	GLY	N-CA-C	-7.20	101.17	110.38
1	E	85	GLY	N-CA-C	-7.11	101.29	110.38
1	E	150	ARG	NE-CZ-NH1	-7.02	114.48	121.50
1	E	94	THR	N-CA-C	6.59	118.47	111.28
1	E	121	PHE	N-CA-C	-6.47	102.45	110.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	85	GLY	N-CA-C	-6.46	101.71	110.43
1	E	153	ILE	CB-CA-C	-6.42	98.76	111.60
1	B	71	ARG	CA-C-N	-6.42	112.74	122.81
1	B	71	ARG	C-N-CA	-6.42	112.74	122.81
1	A	113	SER	N-CA-C	6.37	119.12	108.73
1	E	121(B)	GLU	N-CA-C	6.30	124.23	110.80
1	D	150	ARG	NE-CZ-NH1	-6.30	115.20	121.50
1	D	71	ARG	N-CA-C	6.29	120.22	112.54
1	C	94	THR	N-CA-C	6.25	117.76	111.07
1	C	87	TYR	CA-C-N	6.17	129.59	120.95
1	C	87	TYR	C-N-CA	6.17	129.59	120.95
1	C	88	ASP	N-CA-C	-6.14	100.75	109.96
1	A	22	TRP	N-CA-C	-6.13	100.76	109.96
1	B	109	VAL	N-CA-C	-6.12	101.22	107.89
1	A	71	ARG	N-CA-C	6.06	119.93	112.54
1	A	88	ASP	N-CA-C	-6.05	100.88	109.96
1	B	153	ILE	CB-CA-C	-6.00	101.46	111.29
1	E	153	ILE	CA-C-O	-5.95	110.69	120.80
1	A	121(B)	GLU	N-CA-C	5.92	123.42	110.80
1	D	109	VAL	N-CA-C	-5.90	101.46	107.89
1	C	68	ARG	NE-CZ-NH2	-5.89	113.90	119.20
1	D	88	ASP	N-CA-C	-5.89	101.13	109.96
1	A	77	GLY	N-CA-C	-5.88	99.25	113.18
1	A	12	SER	N-CA-C	5.86	118.45	108.02
1	B	152	ARG	N-CA-C	-5.84	104.55	111.03
1	C	72	TYR	N-CA-C	5.83	118.97	109.59
1	E	113	SER	N-CA-C	5.81	118.21	108.73
1	D	153	ILE	O-C-N	5.81	129.64	122.32
1	A	71	ARG	CA-C-N	-5.79	114.25	122.94
1	A	71	ARG	C-N-CA	-5.79	114.25	122.94
1	C	113	SER	N-CA-C	5.72	118.57	109.24
1	B	22	TRP	N-CA-C	-5.71	101.59	110.10
1	C	59	ILE	N-CA-C	5.71	119.20	112.35
1	D	152	ARG	N-CA-C	-5.62	104.00	111.24
1	E	22	TRP	N-CA-C	-5.60	101.56	109.96
1	B	88	ASP	N-CA-C	-5.59	101.38	109.59
1	D	87	TYR	N-CA-C	5.57	118.71	109.46
1	E	71	ARG	CA-C-N	-5.57	114.58	122.94
1	E	71	ARG	C-N-CA	-5.57	114.58	122.94
1	E	98	ASN	N-CA-C	-5.56	105.30	111.36
1	D	108	GLU	N-CA-C	5.51	119.30	112.24
1	E	88	ASP	N-CA-C	-5.51	101.69	109.96

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	80	PHE	N-CA-C	-5.49	99.47	108.41
1	A	109	VAL	CA-C-O	5.48	123.11	119.38
1	D	113	SER	N-CA-C	5.46	118.14	109.24
1	C	22	TRP	N-CA-C	-5.43	102.01	110.10
1	A	87	TYR	CA-C-N	5.41	128.52	120.95
1	A	87	TYR	C-N-CA	5.41	128.52	120.95
1	C	121(B)	GLU	N-CA-C	5.38	122.27	110.80
1	E	92	VAL	N-CA-C	5.38	115.58	110.42
1	C	96	VAL	N-CA-C	5.34	115.55	110.53
1	A	41	LYS	CA-C-N	-5.33	113.67	121.99
1	A	41	LYS	C-N-CA	-5.33	113.67	121.99
1	B	87	TYR	CA-C-N	5.32	128.90	121.24
1	B	87	TYR	C-N-CA	5.32	128.90	121.24
1	D	72	TYR	N-CA-C	5.30	118.04	109.40
1	B	59	ILE	N-CA-C	5.24	118.63	112.35
1	A	68	ARG	N-CA-C	5.20	118.72	112.38
1	E	87	TYR	N-CA-C	5.18	117.69	109.24
1	C	106	GLU	CA-CB-CG	5.17	124.44	114.10
1	D	59	ILE	N-CA-C	5.17	118.55	112.35
1	D	92	VAL	N-CA-C	5.16	115.88	110.62
1	C	14	LYS	N-CA-C	5.15	118.46	110.17
1	B	48	VAL	N-CA-C	5.12	115.23	107.80
1	A	106	GLU	CA-CB-CG	5.09	124.28	114.10
1	B	12	SER	N-CA-C	5.04	116.61	108.34
1	E	48	VAL	N-CA-C	5.04	115.34	107.99
1	D	106	GLU	CA-CB-CG	5.03	124.17	114.10
1	C	150	ARG	NE-CZ-NH1	-5.03	116.47	121.50
1	C	150	ARG	NE-CZ-NH2	5.00	123.70	119.20

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	87	TYR	Sidechain
1	C	72	TYR	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	1147	0	1127	72	0
1	B	1147	0	1127	71	0
1	C	1135	0	1115	80	0
1	D	1142	0	1122	78	0
1	E	1135	0	1112	73	0
2	A	15	0	0	5	0
2	B	15	0	0	3	0
2	C	10	0	0	8	0
2	D	10	0	0	5	0
2	E	15	0	0	3	0
3	A	10	0	0	1	0
3	B	8	0	0	0	0
3	C	12	0	0	1	0
3	D	7	0	0	0	0
3	E	10	0	0	3	0
All	All	5818	0	5603	350	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 31.

All (350) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:152:ARG:NH2	1:D:153:ILE:HD13	1.61	1.14
1:D:152:ARG:HH21	1:D:153:ILE:CD1	1.64	1.11
1:D:153:ILE:HG12	1:E:51:PHE:HE2	1.09	1.05
1:B:121(B):GLU:O	1:B:121(C):SER:HB3	1.56	1.03
1:C:83:ASP:HB3	3:C:2027:HOH:O	1.59	1.02
1:D:152:ARG:HH21	1:D:153:ILE:HD13	0.88	1.02
1:B:85:GLY:HA3	2:B:195:PO4:O4	1.58	1.00
1:E:121(B):GLU:O	1:E:121(C):SER:HB3	1.57	1.00
1:D:121(B):GLU:O	1:D:121(C):SER:HB3	1.62	0.99
1:C:121(B):GLU:O	1:C:121(C):SER:HB3	1.60	0.98
1:D:121(C):SER:OG	1:D:123:GLU:HB2	1.63	0.98
1:A:85:GLY:HA3	2:A:192:PO4:O3	1.65	0.97
1:A:121(B):GLU:O	1:A:121(C):SER:HB3	1.59	0.96
1:D:153:ILE:HG12	1:E:51:PHE:CE2	2.02	0.94
1:A:37:GLU:HG3	1:A:137:VAL:HG13	1.53	0.90
1:C:121:PHE:HD2	1:C:128:PHE:HE1	1.21	0.89
1:C:121(C):SER:OG	1:C:123:GLU:HB2	1.71	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:153:ILE:HG22	1:D:153:ILE:O	1.70	0.88
1:D:153:ILE:O	1:D:153:ILE:CG2	2.21	0.87
1:C:121:PHE:CD2	1:C:128:PHE:CE1	2.64	0.86
1:B:121(C):SER:OG	1:B:123:GLU:HB2	1.76	0.85
1:A:153:ILE:HD13	1:A:153:ILE:O	1.78	0.83
1:D:22:TRP:CE3	1:D:23:HIS:HD2	1.97	0.83
1:D:12:SER:HA	1:D:45:SER:O	1.80	0.81
1:C:121:PHE:HD2	1:C:128:PHE:CE1	1.97	0.81
1:E:121:PHE:CD2	1:E:128:PHE:CE1	2.69	0.80
1:E:121:PHE:HD2	1:E:128:PHE:HE1	1.28	0.80
1:A:121(C):SER:OG	1:A:123:GLU:HB2	1.79	0.80
1:C:56:ALA:N	2:C:196:PO4:O2	2.15	0.79
1:B:37:GLU:HG3	1:B:137:VAL:HG13	1.65	0.79
1:D:37:GLU:HG3	1:D:137:VAL:HG13	1.64	0.78
1:E:37:GLU:HG3	1:E:137:VAL:HG13	1.65	0.78
1:E:121(C):SER:OG	1:E:123:GLU:HB2	1.83	0.78
1:A:56:ALA:HB3	2:A:190:PO4:O1	1.83	0.78
1:A:153:ILE:HD11	1:B:66:LEU:HD23	1.66	0.78
1:E:22:TRP:CE3	1:E:23:HIS:HD2	2.00	0.78
1:A:121:PHE:HD2	1:A:128:PHE:HE1	1.32	0.77
1:C:121:PHE:CD2	1:C:128:PHE:HE1	1.99	0.77
1:E:121:PHE:HD2	1:E:128:PHE:CE1	2.04	0.76
1:E:146:ILE:O	1:E:150:ARG:HG3	1.86	0.75
1:B:22:TRP:CE3	1:B:23:HIS:HD2	2.04	0.74
1:B:82:ILE:HD13	1:B:82:ILE:C	2.13	0.74
1:A:121:PHE:CD2	1:A:128:PHE:CE1	2.75	0.74
1:E:121:PHE:CD2	1:E:128:PHE:HE1	2.06	0.74
1:C:37:GLU:HG3	1:C:137:VAL:HG13	1.70	0.74
1:C:82:ILE:HG23	1:C:82:ILE:O	1.87	0.73
1:D:121:PHE:HD2	1:D:128:PHE:HE1	1.36	0.73
1:D:85:GLY:HA3	2:D:199:PO4:P	2.28	0.73
1:D:82:ILE:O	1:D:82:ILE:HG23	1.89	0.72
1:E:119:HIS:O	1:E:120:HIS:HB2	1.86	0.72
1:B:121:PHE:CD2	1:B:128:PHE:CE1	2.78	0.72
1:D:119:HIS:O	1:D:120:HIS:HB2	1.89	0.72
1:A:82:ILE:HG23	1:A:82:ILE:O	1.88	0.72
1:B:22:TRP:HH2	1:B:82:ILE:CG2	2.02	0.72
1:D:85:GLY:HA3	2:D:199:PO4:O1	1.89	0.72
1:C:119:HIS:O	1:C:120:HIS:HB2	1.89	0.71
1:A:121:PHE:CD2	1:A:128:PHE:HE1	2.10	0.70
1:E:116:LEU:HD23	1:E:132:PHE:CZ	2.27	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:82:ILE:HD13	1:A:82:ILE:C	2.16	0.70
1:C:12:SER:HA	1:C:45:SER:O	1.91	0.70
1:C:22:TRP:CE3	1:C:23:HIS:HD2	2.10	0.70
1:D:121:PHE:CD2	1:D:128:PHE:CE1	2.79	0.70
1:C:82:ILE:C	1:C:82:ILE:HD13	2.17	0.70
1:C:121(C):SER:O	1:C:123:GLU:N	2.25	0.70
1:A:146:ILE:O	1:A:150:ARG:HG3	1.92	0.69
1:D:25:ASP:OD1	1:D:125:HIS:HE1	1.75	0.69
1:B:119:HIS:O	1:B:120:HIS:HB2	1.92	0.69
1:A:121:PHE:HD2	1:A:128:PHE:CE1	2.09	0.69
1:A:22:TRP:HH2	1:A:82:ILE:CG2	2.06	0.69
1:B:22:TRP:CH2	1:B:82:ILE:HG22	2.28	0.69
1:B:121:PHE:HD2	1:B:128:PHE:HE1	1.41	0.69
1:D:57:TYR:N	2:D:198:PO4:O2	2.19	0.68
1:C:128:PHE:HD2	1:D:87:TYR:HH	1.40	0.68
1:A:38:LEU:HD23	1:A:140:ALA:HB1	1.76	0.68
1:A:153:ILE:HD11	1:B:66:LEU:CD2	2.24	0.68
1:C:22:TRP:HH2	1:C:82:ILE:CG2	2.06	0.67
1:B:82:ILE:HG23	1:B:82:ILE:O	1.95	0.67
1:E:83:ASP:OD1	3:E:2045:HOH:O	2.13	0.67
1:D:153:ILE:HG21	1:E:66:LEU:HD21	1.77	0.67
1:D:57:TYR:HB2	2:D:198:PO4:O2	1.95	0.67
1:D:121:PHE:HD2	1:D:128:PHE:CE1	2.13	0.67
1:D:121:PHE:CD2	1:D:128:PHE:HE1	2.14	0.67
1:B:152:ARG:O	1:B:153:ILE:C	2.37	0.66
1:C:85:GLY:HA3	2:C:197:PO4:P	2.35	0.66
1:A:85:GLY:HA3	2:A:192:PO4:P	2.34	0.66
1:C:25:ASP:OD1	1:C:125:HIS:HE1	1.77	0.66
1:B:89:HIS:HD2	2:B:194:PO4:O1	1.78	0.66
1:B:121(C):SER:O	1:B:123:GLU:N	2.29	0.65
1:E:116:LEU:HD23	1:E:132:PHE:CE1	2.32	0.65
1:D:82:ILE:HD13	1:D:82:ILE:C	2.22	0.65
1:D:22:TRP:HH2	1:D:82:ILE:CG2	2.10	0.65
1:C:22:TRP:CH2	1:C:82:ILE:HG22	2.32	0.65
1:E:82:ILE:C	1:E:82:ILE:HD13	2.22	0.65
1:E:22:TRP:HH2	1:E:82:ILE:CG2	2.09	0.65
1:D:146:ILE:O	1:D:150:ARG:HG3	1.97	0.64
1:B:85:GLY:HA3	2:B:195:PO4:P	2.38	0.64
1:B:121:PHE:HD2	1:B:128:PHE:CE1	2.14	0.64
1:A:22:TRP:CH2	1:A:82:ILE:HG22	2.33	0.64
1:A:25:ASP:OD1	1:A:125:HIS:HE1	1.80	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:121(C):SER:O	1:E:123:GLU:N	2.31	0.63
1:A:87:TYR:HH	1:E:128:PHE:HD2	1.46	0.63
1:C:116:LEU:HD23	1:C:132:PHE:CE1	2.34	0.63
1:E:22:TRP:CH2	1:E:82:ILE:HG22	2.33	0.63
1:D:121(C):SER:O	1:D:123:GLU:N	2.32	0.62
1:E:82:ILE:HG23	1:E:82:ILE:O	1.98	0.62
1:E:85:GLY:HA3	2:E:202:PO4:P	2.40	0.62
1:C:128:PHE:HD2	1:D:87:TYR:OH	1.83	0.62
1:B:38:LEU:HD23	1:B:140:ALA:HB1	1.82	0.61
1:E:22:TRP:CZ3	1:E:23:HIS:HD2	2.18	0.61
1:C:119:HIS:HB2	1:D:86:ILE:O	2.00	0.61
1:D:121(A):HIS:O	1:D:121(B):GLU:O	2.18	0.61
1:B:116:LEU:HD23	1:B:132:PHE:CZ	2.35	0.61
1:B:121(B):GLU:O	1:B:121(C):SER:CB	2.41	0.60
1:D:19:GLN:HE22	1:D:31:ARG:HE	1.48	0.60
1:B:12:SER:HA	1:B:45:SER:O	2.00	0.60
1:D:22:TRP:CZ3	1:D:23:HIS:HD2	2.18	0.60
1:D:22:TRP:CH2	1:D:82:ILE:HG22	2.37	0.60
1:B:121(A):HIS:O	1:B:121(B):GLU:O	2.20	0.60
1:A:121(C):SER:O	1:A:123:GLU:N	2.35	0.59
1:E:83:ASP:CG	3:E:2045:HOH:O	2.45	0.59
1:C:19:GLN:HE22	1:C:31:ARG:HE	1.50	0.59
1:C:22:TRP:HH2	1:C:82:ILE:HG21	1.67	0.59
1:C:85:GLY:HA3	2:C:197:PO4:O3	2.02	0.59
1:D:22:TRP:CZ3	1:D:23:HIS:CD2	2.91	0.59
1:E:25:ASP:OD1	1:E:125:HIS:HE1	1.85	0.59
1:C:107:THR:O	1:C:108:GLU:HB2	2.02	0.59
1:C:82:ILE:O	1:C:82:ILE:CG2	2.51	0.58
1:A:22:TRP:HH2	1:A:82:ILE:HG21	1.67	0.58
1:E:22:TRP:CZ3	1:E:23:HIS:CD2	2.91	0.58
1:C:86:ILE:N	2:C:197:PO4:O4	2.37	0.57
1:C:121(C):SER:C	1:C:123:GLU:N	2.62	0.57
1:D:26:ILE:O	1:D:27:VAL:C	2.47	0.57
1:A:82:ILE:O	1:A:82:ILE:CG2	2.53	0.57
1:A:119:HIS:O	1:A:120:HIS:HB2	2.05	0.57
1:B:22:TRP:HH2	1:B:82:ILE:HG21	1.68	0.57
1:E:64:LYS:NZ	1:E:106:GLU:OE1	2.34	0.57
1:C:128:PHE:CD2	1:D:87:TYR:OH	2.55	0.56
1:C:50:ILE:HD12	1:C:50:ILE:N	2.20	0.56
1:E:121(A):HIS:O	1:E:121(B):GLU:O	2.23	0.56
1:D:121(C):SER:C	1:D:123:GLU:N	2.64	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:122:LYS:N	1:B:122:LYS:HD2	2.20	0.55
1:B:116:LEU:HD23	1:B:132:PHE:CE1	2.42	0.55
1:D:22:TRP:HH2	1:D:82:ILE:HG21	1.72	0.55
1:B:82:ILE:C	1:B:82:ILE:CD1	2.78	0.55
1:A:87:TYR:OH	1:E:128:PHE:HD2	1.89	0.55
1:B:121:PHE:CD2	1:B:128:PHE:HE1	2.19	0.55
1:D:116:LEU:HD23	1:D:132:PHE:CZ	2.41	0.54
1:B:22:TRP:CZ3	1:B:23:HIS:HD2	2.25	0.54
1:B:146:ILE:O	1:B:150:ARG:HG3	2.06	0.54
1:C:122:LYS:HD2	1:C:122:LYS:N	2.22	0.54
1:B:25:ASP:OD1	1:B:125:HIS:HE1	1.89	0.54
1:B:64:LYS:NZ	1:B:106:GLU:OE1	2.39	0.54
1:A:82:ILE:C	1:A:82:ILE:CD1	2.81	0.54
1:D:82:ILE:O	1:D:82:ILE:CG2	2.54	0.54
1:C:82:ILE:C	1:C:82:ILE:CD1	2.80	0.54
1:B:152:ARG:O	1:B:153:ILE:O	2.26	0.53
1:C:121(A):HIS:O	1:C:121(B):GLU:O	2.26	0.53
1:E:22:TRP:HH2	1:E:82:ILE:HG21	1.72	0.53
1:A:116:LEU:HD23	1:A:132:PHE:CZ	2.42	0.53
1:C:121:PHE:HB2	1:C:128:PHE:CZ	2.43	0.53
1:A:150:ARG:HA	1:A:153:ILE:HG22	1.91	0.53
1:D:152:ARG:CZ	1:D:153:ILE:HD13	2.37	0.53
1:A:21:ARG:NE	1:A:21:ARG:HA	2.23	0.53
1:B:121(C):SER:C	1:B:123:GLU:N	2.65	0.53
1:B:50:ILE:N	1:B:50:ILE:HD12	2.23	0.53
1:A:121(B):GLU:O	1:A:121(C):SER:CB	2.43	0.52
1:D:121(C):SER:O	1:D:124:HIS:N	2.42	0.52
1:A:89:HIS:HD2	2:A:191:PO4:O2	1.92	0.52
1:E:121:PHE:HB2	1:E:128:PHE:CZ	2.44	0.52
1:A:121(C):SER:C	1:A:123:GLU:N	2.67	0.52
1:C:22:TRP:CH2	1:C:82:ILE:CG2	2.88	0.52
1:D:81:VAL:O	1:D:117:THR:HG23	2.10	0.52
1:E:121(C):SER:C	1:E:123:GLU:N	2.68	0.52
1:B:22:TRP:CH2	1:B:82:ILE:CG2	2.86	0.51
1:D:84:GLY:HA3	1:D:120:HIS:CE1	2.45	0.51
1:D:22:TRP:CH2	1:D:82:ILE:CG2	2.93	0.51
1:D:25:ASP:OD1	1:D:125:HIS:CE1	2.61	0.51
1:B:121:PHE:HB2	1:B:128:PHE:CZ	2.46	0.51
1:A:128:PHE:CD2	1:B:87:TYR:OH	2.64	0.51
1:E:121:PHE:CD1	1:E:121:PHE:C	2.88	0.51
1:D:22:TRP:HH2	1:D:82:ILE:HG22	1.76	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:81:VAL:O	1:C:117:THR:HG23	2.10	0.50
1:B:28:ASP:OD2	1:B:31:ARG:NH2	2.38	0.50
1:A:116:LEU:HD23	1:A:132:PHE:CE1	2.47	0.50
1:D:122:LYS:N	1:D:122:LYS:HD2	2.26	0.50
1:D:128:PHE:HD2	1:E:87:TYR:OH	1.93	0.50
1:A:121(A):HIS:O	1:A:121(B):GLU:O	2.29	0.50
1:E:82:ILE:HD13	1:E:83:ASP:C	2.37	0.50
1:A:150:ARG:O	1:A:153:ILE:HG22	2.11	0.49
1:D:101:MET:HA	1:D:104:GLN:HE21	1.77	0.49
1:D:116:LEU:HD23	1:D:132:PHE:CE1	2.47	0.49
1:E:28:ASP:OD2	1:E:31:ARG:NH2	2.44	0.49
1:B:82:ILE:CG2	1:B:82:ILE:O	2.60	0.49
1:D:128:PHE:CD2	1:E:87:TYR:OH	2.63	0.49
1:E:121(A):HIS:C	1:E:121(B):GLU:HG3	2.30	0.49
1:A:51:PHE:HB3	3:A:2036:HOH:O	2.12	0.49
1:A:25:ASP:OD1	1:A:125:HIS:CE1	2.64	0.49
1:A:128:PHE:HD2	1:B:87:TYR:OH	1.95	0.49
1:E:82:ILE:C	1:E:82:ILE:CD1	2.86	0.49
1:B:22:TRP:CZ3	1:B:23:HIS:CD2	3.00	0.49
1:A:153:ILE:HD13	1:A:153:ILE:C	2.38	0.49
1:D:149:GLU:OE2	1:D:152:ARG:NH1	2.45	0.49
1:E:82:ILE:CG2	1:E:82:ILE:O	2.61	0.49
1:C:119:HIS:N	1:D:86:ILE:O	2.27	0.49
1:D:19:GLN:NE2	1:D:31:ARG:HE	2.11	0.48
1:D:57:TYR:CB	2:D:198:PO4:O2	2.61	0.48
1:E:68:ARG:HG2	1:E:68:ARG:HH11	1.78	0.48
1:C:19:GLN:NE2	1:C:31:ARG:HE	2.10	0.48
1:C:88:ASP:OD1	1:C:90:ASP:OD2	2.32	0.48
1:A:22:TRP:CH2	1:A:82:ILE:CG2	2.89	0.48
1:C:121:PHE:HB2	1:C:128:PHE:CE1	2.48	0.48
1:C:152:ARG:O	1:C:154:ALA:N	2.46	0.48
1:A:28:ASP:OD2	1:A:31:ARG:NH2	2.38	0.48
1:E:68:ARG:HH11	1:E:68:ARG:CG	2.25	0.48
1:B:90:ASP:OD2	1:B:90:ASP:N	2.46	0.48
1:E:21:ARG:HA	1:E:21:ARG:NE	2.28	0.48
1:A:57:TYR:N	2:A:190:PO4:O1	2.43	0.48
1:C:116:LEU:HD23	1:C:132:PHE:CZ	2.49	0.48
1:D:82:ILE:C	1:D:82:ILE:CD1	2.86	0.48
1:E:135:LYS:NZ	3:E:2032:HOH:O	2.47	0.48
1:A:150:ARG:HA	1:A:153:ILE:CG2	2.44	0.48
1:B:121:PHE:HB2	1:B:128:PHE:CE1	2.48	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:85:GLY:HA3	2:E:202:PO4:O3	2.14	0.48
1:D:121(C):SER:C	1:D:123:GLU:H	2.22	0.47
1:A:152:ARG:O	1:A:154:ALA:N	2.47	0.47
1:D:21:ARG:HA	1:D:21:ARG:NE	2.29	0.47
1:A:121:PHE:CD1	1:A:121:PHE:C	2.90	0.47
1:B:87:TYR:CD1	1:B:87:TYR:N	2.83	0.47
1:E:101:MET:HA	1:E:104:GLN:HE21	1.79	0.47
1:D:152:ARG:HE	1:D:152:ARG:HB3	1.41	0.47
1:B:68:ARG:HG2	1:B:68:ARG:HH11	1.80	0.47
1:C:146:ILE:O	1:C:150:ARG:HG3	2.15	0.47
1:C:57:TYR:N	2:C:196:PO4:O2	2.43	0.46
1:C:22:TRP:CD2	1:C:23:HIS:HD2	2.33	0.46
1:D:29:GLU:HG3	1:D:133:LYS:HE3	1.97	0.46
1:A:84:GLY:HA3	1:A:120:HIS:CE1	2.51	0.46
1:A:121:PHE:HB2	1:A:128:PHE:CZ	2.50	0.46
1:C:121(C):SER:O	1:C:124:HIS:N	2.49	0.46
1:A:133:LYS:HB2	1:A:133:LYS:HE3	1.74	0.46
1:A:131:HIS:ND1	1:B:87:TYR:CZ	2.84	0.46
1:B:119:HIS:HB2	1:C:86:ILE:O	2.15	0.46
1:C:121(C):SER:C	1:C:123:GLU:H	2.23	0.46
1:E:22:TRP:CD2	1:E:23:HIS:HD2	2.34	0.46
1:A:152:ARG:C	1:A:154:ALA:H	2.22	0.46
1:B:19:GLN:HE22	1:B:31:ARG:HE	1.64	0.45
1:B:83:ASP:HA	1:B:89:HIS:HB2	1.98	0.45
1:E:26:ILE:O	1:E:27:VAL:C	2.57	0.45
1:E:83:ASP:HA	1:E:89:HIS:HB2	1.98	0.45
1:C:85:GLY:HA3	2:C:197:PO4:O1	2.17	0.45
1:C:121(C):SER:O	1:C:122:LYS:C	2.59	0.45
1:E:53:VAL:HG13	1:E:58:GLU:HB2	1.98	0.45
1:A:101:MET:HA	1:A:104:GLN:HE21	1.81	0.45
1:B:124:HIS:O	1:B:127:PHE:N	2.49	0.45
1:E:116:LEU:CD2	1:E:132:PHE:CZ	2.98	0.45
1:C:25:ASP:OD1	1:C:125:HIS:CE1	2.64	0.45
1:C:49:GLU:C	1:C:50:ILE:HD12	2.42	0.45
1:D:50:ILE:N	1:D:50:ILE:HD12	2.30	0.45
1:C:152:ARG:C	1:C:154:ALA:H	2.24	0.45
1:A:150:ARG:HH11	1:A:150:ARG:HD3	1.38	0.45
1:E:22:TRP:CH2	1:E:82:ILE:CG2	2.91	0.45
1:C:101:MET:HA	1:C:104:GLN:HE21	1.82	0.45
1:A:83:ASP:HA	1:A:89:HIS:HB2	1.98	0.45
1:C:38:LEU:HD23	1:C:140:ALA:HB1	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:122:LYS:HD2	1:E:122:LYS:N	2.32	0.45
1:A:87:TYR:OH	1:E:128:PHE:CD2	2.63	0.44
1:B:82:ILE:HD13	1:B:83:ASP:C	2.41	0.44
1:C:82:ILE:HD13	1:C:83:ASP:C	2.41	0.44
1:D:153:ILE:CG2	1:E:66:LEU:HD21	2.45	0.44
1:C:50:ILE:N	1:C:50:ILE:CD1	2.80	0.44
1:E:86:ILE:N	2:E:202:PO4:O3	2.47	0.44
1:B:133:LYS:HE3	1:B:133:LYS:HB2	1.81	0.44
1:B:121(C):SER:C	1:B:123:GLU:H	2.25	0.44
1:B:23:HIS:O	1:B:24:ALA:C	2.60	0.44
1:B:128:PHE:CD2	1:C:87:TYR:OH	2.70	0.44
1:C:22:TRP:CZ3	1:C:23:HIS:CD2	3.06	0.44
1:E:54:PRO:HD2	1:E:58:GLU:HG2	1.99	0.44
1:E:84:GLY:HA3	1:E:120:HIS:CE1	2.52	0.44
1:E:121(C):SER:C	1:E:123:GLU:H	2.26	0.44
1:B:19:GLN:NE2	1:B:31:ARG:HE	2.16	0.44
1:C:26:ILE:O	1:C:27:VAL:C	2.60	0.44
1:C:152:ARG:HE	1:C:152:ARG:HB3	1.34	0.44
1:D:29:GLU:HG3	1:D:133:LYS:CE	2.48	0.44
1:A:152:ARG:C	1:A:154:ALA:N	2.75	0.44
1:C:87:TYR:HD1	2:C:197:PO4:O4	2.00	0.44
1:D:83:ASP:HA	1:D:89:HIS:HB2	2.00	0.44
1:E:133:LYS:HE3	1:E:133:LYS:HB2	1.74	0.43
1:C:87:TYR:N	2:C:197:PO4:O4	2.41	0.43
1:C:146:ILE:HD12	1:C:146:ILE:HA	1.88	0.43
1:D:90:ASP:OD2	1:D:90:ASP:N	2.49	0.43
1:A:121:PHE:HB2	1:A:128:PHE:CE1	2.53	0.43
1:C:87:TYR:N	1:C:87:TYR:CD1	2.86	0.43
1:C:121:PHE:CD1	1:C:121:PHE:C	2.95	0.43
1:E:78:ALA:HA	1:E:114:VAL:O	2.18	0.43
1:E:121:PHE:HB2	1:E:128:PHE:CE1	2.53	0.43
1:A:144:LEU:HA	1:A:144:LEU:HD23	1.85	0.43
1:A:107:THR:O	1:A:108:GLU:HB2	2.19	0.43
1:E:121(B):GLU:O	1:E:121(C):SER:CB	2.42	0.43
1:A:26:ILE:O	1:A:27:VAL:C	2.61	0.43
1:A:78:ALA:HA	1:A:114:VAL:O	2.19	0.43
1:D:133:LYS:HE3	1:D:133:LYS:HB2	1.71	0.43
1:D:121:PHE:HB2	1:D:128:PHE:CZ	2.53	0.43
1:E:121(C):SER:O	1:E:124:HIS:N	2.50	0.43
1:B:22:TRP:CD2	1:B:23:HIS:HD2	2.36	0.43
1:B:50:ILE:N	1:B:50:ILE:CD1	2.81	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:121(C):SER:HG	1:D:123:GLU:HB2	1.79	0.42
1:A:87:TYR:CD1	1:A:87:TYR:N	2.87	0.42
1:C:83:ASP:HA	1:C:89:HIS:HB2	2.01	0.42
1:A:119:HIS:HB2	1:B:86:ILE:O	2.19	0.42
1:C:84:GLY:HA3	1:C:120:HIS:CE1	2.54	0.42
1:C:133:LYS:HE3	1:C:133:LYS:HB2	1.81	0.42
1:B:25:ASP:OD1	1:B:125:HIS:CE1	2.72	0.42
1:C:82:ILE:HD13	1:C:83:ASP:N	2.34	0.42
1:D:82:ILE:HD13	1:D:83:ASP:C	2.44	0.42
1:D:22:TRP:CE3	1:D:23:HIS:CD2	2.89	0.42
1:E:29:GLU:HG3	1:E:133:LYS:HE3	2.02	0.42
1:E:47:GLU:H	1:E:47:GLU:HG2	1.70	0.42
1:A:121(C):SER:C	1:A:123:GLU:H	2.26	0.42
1:B:68:ARG:HG2	1:B:68:ARG:NH1	2.35	0.42
1:B:121(C):SER:O	1:B:124:HIS:N	2.52	0.42
1:C:21:ARG:HA	1:C:21:ARG:NE	2.35	0.42
1:D:78:ALA:HA	1:D:114:VAL:O	2.20	0.42
1:E:22:TRP:CE3	1:E:23:HIS:CD2	2.92	0.42
1:C:22:TRP:CZ3	1:C:23:HIS:HD2	2.37	0.42
1:A:88:ASP:OD1	1:A:90:ASP:OD2	2.37	0.42
1:B:153:ILE:HB	1:B:154:ALA:H	1.51	0.42
1:C:19:GLN:O	1:C:19:GLN:HG3	2.20	0.42
1:A:121(C):SER:O	1:A:124:HIS:N	2.52	0.41
1:E:110:PRO:HG3	1:E:150:ARG:HD2	2.01	0.41
1:D:121:PHE:CD1	1:D:121:PHE:C	2.97	0.41
1:A:90:ASP:OD2	1:A:90:ASP:N	2.53	0.41
1:D:121:PHE:HB2	1:D:128:PHE:CE1	2.55	0.41
1:B:144:LEU:HA	1:B:144:LEU:HD23	1.83	0.41
1:B:131:HIS:ND1	1:C:87:TYR:CZ	2.89	0.41
1:C:121:PHE:CB	1:C:128:PHE:CZ	3.03	0.41
1:C:149:GLU:OE2	1:C:152:ARG:NH1	2.54	0.41
1:E:150:ARG:HH11	1:E:150:ARG:HD3	1.51	0.41
1:A:150:ARG:C	1:A:153:ILE:HG22	2.46	0.41
1:B:105:LEU:HD23	1:B:105:LEU:HA	1.84	0.41
1:D:38:LEU:HD23	1:D:140:ALA:HB1	2.02	0.41
1:A:152:ARG:HE	1:A:152:ARG:HB3	1.23	0.41
1:C:121(B):GLU:O	1:C:121(C):SER:CB	2.44	0.41
1:E:152:ARG:HE	1:E:152:ARG:HB3	1.33	0.41
1:B:124:HIS:O	1:B:127:PHE:HB3	2.21	0.41
1:A:21:ARG:HA	1:A:21:ARG:HE	1.86	0.40
1:B:128:PHE:HD2	1:C:87:TYR:OH	2.04	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:150:ARG:CA	1:A:153:ILE:HG22	2.52	0.40
1:D:105:LEU:HA	1:D:105:LEU:HD23	1.76	0.40
1:B:150:ARG:HH11	1:B:150:ARG:HD3	1.60	0.40
1:E:121(C):SER:O	1:E:122:LYS:C	2.64	0.40
1:E:149:GLU:OE2	1:E:152:ARG:NH1	2.54	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	146/158 (92%)	132 (90%)	8 (6%)	6 (4%)	2	5
1	B	146/158 (92%)	135 (92%)	4 (3%)	7 (5%)	2	3
1	C	144/158 (91%)	133 (92%)	5 (4%)	6 (4%)	2	4
1	D	145/158 (92%)	136 (94%)	3 (2%)	6 (4%)	2	5
1	E	144/158 (91%)	134 (93%)	5 (4%)	5 (4%)	3	6
All	All	725/790 (92%)	670 (92%)	25 (3%)	30 (4%)	2	5

All (30) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	120	HIS
1	A	121(B)	GLU
1	A	121(C)	SER
1	A	122	LYS
1	B	120	HIS
1	B	121(B)	GLU
1	B	121(C)	SER
1	B	122	LYS
1	B	153	ILE

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Mol	Chain	Res	Type
1	B	154	ALA
1	C	120	HIS
1	C	121(B)	GLU
1	C	121(C)	SER
1	C	122	LYS
1	D	120	HIS
1	D	121(B)	GLU
1	D	121(C)	SER
1	D	122	LYS
1	E	120	HIS
1	E	121(B)	GLU
1	E	121(C)	SER
1	E	122	LYS
1	A	153	ILE
1	C	153	ILE
1	A	44	GLY
1	C	44	GLY
1	B	44	GLY
1	D	44	GLY
1	E	44	GLY
1	D	54	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	116/126 (92%)	100 (86%)	16 (14%)	3	9
1	B	116/126 (92%)	102 (88%)	14 (12%)	5	12
1	C	115/126 (91%)	101 (88%)	14 (12%)	5	12
1	D	116/126 (92%)	100 (86%)	16 (14%)	3	9
1	E	115/126 (91%)	100 (87%)	15 (13%)	4	10
All	All	578/630 (92%)	503 (87%)	75 (13%)	4	10

All (75) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	SER
1	A	14	LYS
1	A	38	LEU
1	A	53	VAL
1	A	62	HIS
1	A	68	ARG
1	A	82	ILE
1	A	116	LEU
1	A	119	HIS
1	A	121	PHE
1	A	121(A)	HIS
1	A	121(B)	GLU
1	A	121(C)	SER
1	A	122	LYS
1	A	125	HIS
1	A	153	ILE
1	B	14	LYS
1	B	38	LEU
1	B	53	VAL
1	B	62	HIS
1	B	68	ARG
1	B	82	ILE
1	B	116	LEU
1	B	119	HIS
1	B	121	PHE
1	B	121(A)	HIS
1	B	121(B)	GLU
1	B	121(C)	SER
1	B	122	LYS
1	B	125	HIS
1	C	14	LYS
1	C	38	LEU
1	C	53	VAL
1	C	62	HIS
1	C	68	ARG
1	C	82	ILE
1	C	116	LEU
1	C	119	HIS
1	C	121	PHE
1	C	121(A)	HIS
1	C	121(B)	GLU
1	C	121(C)	SER
1	C	122	LYS

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Mol	Chain	Res	Type
1	C	125	HIS
1	D	11	THR
1	D	14	LYS
1	D	38	LEU
1	D	53	VAL
1	D	82	ILE
1	D	116	LEU
1	D	118	PRO
1	D	119	HIS
1	D	121	PHE
1	D	121(A)	HIS
1	D	121(B)	GLU
1	D	121(C)	SER
1	D	122	LYS
1	D	123	GLU
1	D	125	HIS
1	D	153	ILE
1	E	12	SER
1	E	14	LYS
1	E	38	LEU
1	E	53	VAL
1	E	62	HIS
1	E	68	ARG
1	E	82	ILE
1	E	116	LEU
1	E	119	HIS
1	E	121	PHE
1	E	121(A)	HIS
1	E	121(B)	GLU
1	E	121(C)	SER
1	E	122	LYS
1	E	125	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (36) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	19	GLN
1	A	23	HIS
1	A	89	HIS
1	A	98	ASN
1	A	104	GLN
1	A	119	HIS

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Mol	Chain	Res	Type
1	A	121(A)	HIS
1	A	125	HIS
1	A	141	HIS
1	B	19	GLN
1	B	23	HIS
1	B	89	HIS
1	B	98	ASN
1	B	119	HIS
1	B	121(A)	HIS
1	B	125	HIS
1	C	19	GLN
1	C	23	HIS
1	C	98	ASN
1	C	104	GLN
1	C	119	HIS
1	C	125	HIS
1	D	19	GLN
1	D	23	HIS
1	D	98	ASN
1	D	104	GLN
1	D	119	HIS
1	D	125	HIS
1	E	19	GLN
1	E	23	HIS
1	E	62	HIS
1	E	98	ASN
1	E	104	GLN
1	E	119	HIS
1	E	121(A)	HIS
1	E	125	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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

13 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	PO4	A	192	-	4,4,4	2.34	3 (75%)	6,6,6	0.68	0
2	PO4	A	190	-	4,4,4	1.47	0	6,6,6	0.53	0
2	PO4	C	197	-	4,4,4	1.84	1 (25%)	6,6,6	0.51	0
2	PO4	E	200	-	4,4,4	1.43	1 (25%)	6,6,6	0.39	0
2	PO4	A	191	-	4,4,4	1.52	0	6,6,6	0.60	0
2	PO4	C	196	-	4,4,4	1.38	1 (25%)	6,6,6	0.60	0
2	PO4	D	198	-	4,4,4	1.52	1 (25%)	6,6,6	0.49	0
2	PO4	D	199	-	4,4,4	2.18	2 (50%)	6,6,6	0.60	0
2	PO4	E	201	-	4,4,4	1.39	1 (25%)	6,6,6	0.45	0
2	PO4	B	193	-	4,4,4	1.27	0	6,6,6	0.48	0
2	PO4	E	202	-	4,4,4	1.74	1 (25%)	6,6,6	0.49	0
2	PO4	B	195	-	4,4,4	2.15	2 (50%)	6,6,6	0.49	0
2	PO4	B	194	-	4,4,4	1.38	1 (25%)	6,6,6	0.46	0

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	192	PO4	P-O2	-3.04	1.45	1.54
2	B	195	PO4	P-O2	-2.96	1.46	1.54
2	D	198	PO4	P-O2	-2.82	1.46	1.54
2	D	199	PO4	P-O3	-2.81	1.46	1.54
2	B	195	PO4	P-O4	-2.68	1.46	1.54
2	C	197	PO4	P-O4	-2.44	1.47	1.54
2	C	196	PO4	P-O2	-2.35	1.47	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	192	PO4	P-O3	-2.31	1.47	1.54
2	A	192	PO4	P-O4	-2.29	1.48	1.54
2	E	202	PO4	P-O4	-2.24	1.48	1.54
2	D	199	PO4	P-O4	-2.21	1.48	1.54
2	E	201	PO4	P-O4	-2.13	1.48	1.54
2	E	200	PO4	P-O2	-2.01	1.48	1.54
2	B	194	PO4	P-O2	-2.01	1.48	1.54

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

10 monomers are involved in 24 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	192	PO4	2	0
2	A	190	PO4	2	0
2	C	197	PO4	6	0
2	A	191	PO4	1	0
2	C	196	PO4	2	0
2	D	198	PO4	3	0
2	D	199	PO4	2	0
2	E	202	PO4	3	0
2	B	195	PO4	2	0
2	B	194	PO4	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.