



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:22 PM UTC

PDB ID : 4O6Z / pdb_00004o6z
Title : Crystal structure of serine hydroxymethyltransferase with covalently bound
PLP Schiff-base from Plasmodium falciparum
Authors : Chitnumsub, P.; Jaruwat, A.; Leartsakulpanich, U.
Deposited on : 2013-12-24
Resolution : 2.98 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

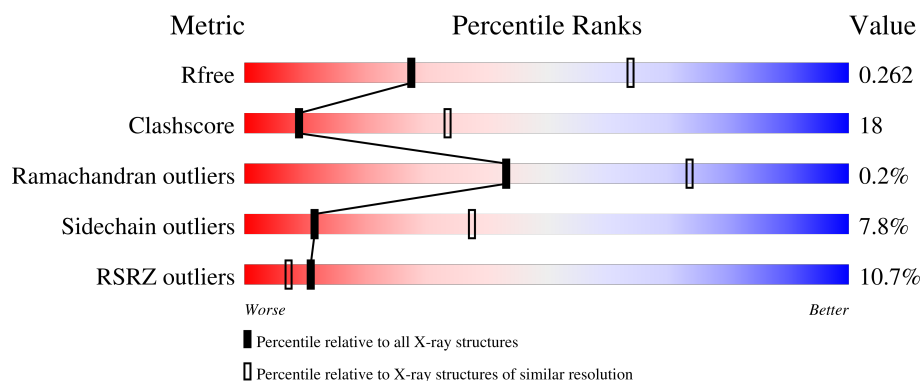
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.98 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3580 (3.00-2.96)
Clashscore	190562	3904 (3.00-2.96)
Ramachandran outliers	187476	3761 (3.00-2.96)
Sidechain outliers	187428	3764 (3.00-2.96)
RSRZ outliers	180081	3579 (3.00-2.96)

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

Mol	Chain	Length	Quality of chain
1	A	480	
1	B	480	
1	C	480	
1	D	480	

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
3	PO4	D	501	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 14254 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 Serine hydroxymethyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	448	Total	C	N	O	S	0	0	0
			3544	2251	603	672	18			
1	B	445	Total	C	N	O	S	0	0	0
			3516	2231	598	669	18			
1	C	441	Total	C	N	O	S	0	0	0
			3490	2215	594	664	17			
1	D	445	Total	C	N	O	S	0	0	0
			3516	2231	598	669	18			

There are 156 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-37	MET	-	expression tag	UNP Q8I566
A	-36	ARG	-	expression tag	UNP Q8I566
A	-35	GLY	-	expression tag	UNP Q8I566
A	-34	SER	-	expression tag	UNP Q8I566
A	-33	HIS	-	expression tag	UNP Q8I566
A	-32	HIS	-	expression tag	UNP Q8I566
A	-31	HIS	-	expression tag	UNP Q8I566
A	-30	HIS	-	expression tag	UNP Q8I566
A	-29	HIS	-	expression tag	UNP Q8I566
A	-28	HIS	-	expression tag	UNP Q8I566
A	-27	GLY	-	expression tag	UNP Q8I566
A	-26	MET	-	expression tag	UNP Q8I566
A	-25	ALA	-	expression tag	UNP Q8I566
A	-24	SER	-	expression tag	UNP Q8I566
A	-23	MET	-	expression tag	UNP Q8I566
A	-22	THR	-	expression tag	UNP Q8I566
A	-21	GLY	-	expression tag	UNP Q8I566
A	-20	GLY	-	expression tag	UNP Q8I566
A	-19	GLN	-	expression tag	UNP Q8I566
A	-18	GLN	-	expression tag	UNP Q8I566
A	-17	MET	-	expression tag	UNP Q8I566

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-16	GLY	-	expression tag	UNP Q8I566
A	-15	ARG	-	expression tag	UNP Q8I566
A	-14	ASP	-	expression tag	UNP Q8I566
A	-13	LEU	-	expression tag	UNP Q8I566
A	-12	TYR	-	expression tag	UNP Q8I566
A	-11	ASP	-	expression tag	UNP Q8I566
A	-10	ASP	-	expression tag	UNP Q8I566
A	-9	ASP	-	expression tag	UNP Q8I566
A	-8	ASP	-	expression tag	UNP Q8I566
A	-7	LYS	-	expression tag	UNP Q8I566
A	-6	ASP	-	expression tag	UNP Q8I566
A	-5	HIS	-	expression tag	UNP Q8I566
A	-4	PRO	-	expression tag	UNP Q8I566
A	-3	PHE	-	expression tag	UNP Q8I566
A	-2	THR	-	expression tag	UNP Q8I566
A	-1	PRO	-	expression tag	UNP Q8I566
A	0	GLY	-	expression tag	UNP Q8I566
A	292	GLU	PHE	engineered mutation	UNP Q8I566
B	-37	MET	-	expression tag	UNP Q8I566
B	-36	ARG	-	expression tag	UNP Q8I566
B	-35	GLY	-	expression tag	UNP Q8I566
B	-34	SER	-	expression tag	UNP Q8I566
B	-33	HIS	-	expression tag	UNP Q8I566
B	-32	HIS	-	expression tag	UNP Q8I566
B	-31	HIS	-	expression tag	UNP Q8I566
B	-30	HIS	-	expression tag	UNP Q8I566
B	-29	HIS	-	expression tag	UNP Q8I566
B	-28	HIS	-	expression tag	UNP Q8I566
B	-27	GLY	-	expression tag	UNP Q8I566
B	-26	MET	-	expression tag	UNP Q8I566
B	-25	ALA	-	expression tag	UNP Q8I566
B	-24	SER	-	expression tag	UNP Q8I566
B	-23	MET	-	expression tag	UNP Q8I566
B	-22	THR	-	expression tag	UNP Q8I566
B	-21	GLY	-	expression tag	UNP Q8I566
B	-20	GLY	-	expression tag	UNP Q8I566
B	-19	GLN	-	expression tag	UNP Q8I566
B	-18	GLN	-	expression tag	UNP Q8I566
B	-17	MET	-	expression tag	UNP Q8I566
B	-16	GLY	-	expression tag	UNP Q8I566
B	-15	ARG	-	expression tag	UNP Q8I566
B	-14	ASP	-	expression tag	UNP Q8I566

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-13	LEU	-	expression tag	UNP Q8I566
B	-12	TYR	-	expression tag	UNP Q8I566
B	-11	ASP	-	expression tag	UNP Q8I566
B	-10	ASP	-	expression tag	UNP Q8I566
B	-9	ASP	-	expression tag	UNP Q8I566
B	-8	ASP	-	expression tag	UNP Q8I566
B	-7	LYS	-	expression tag	UNP Q8I566
B	-6	ASP	-	expression tag	UNP Q8I566
B	-5	HIS	-	expression tag	UNP Q8I566
B	-4	PRO	-	expression tag	UNP Q8I566
B	-3	PHE	-	expression tag	UNP Q8I566
B	-2	THR	-	expression tag	UNP Q8I566
B	-1	PRO	-	expression tag	UNP Q8I566
B	0	GLY	-	expression tag	UNP Q8I566
B	292	GLU	PHE	engineered mutation	UNP Q8I566
C	-37	MET	-	expression tag	UNP Q8I566
C	-36	ARG	-	expression tag	UNP Q8I566
C	-35	GLY	-	expression tag	UNP Q8I566
C	-34	SER	-	expression tag	UNP Q8I566
C	-33	HIS	-	expression tag	UNP Q8I566
C	-32	HIS	-	expression tag	UNP Q8I566
C	-31	HIS	-	expression tag	UNP Q8I566
C	-30	HIS	-	expression tag	UNP Q8I566
C	-29	HIS	-	expression tag	UNP Q8I566
C	-28	HIS	-	expression tag	UNP Q8I566
C	-27	GLY	-	expression tag	UNP Q8I566
C	-26	MET	-	expression tag	UNP Q8I566
C	-25	ALA	-	expression tag	UNP Q8I566
C	-24	SER	-	expression tag	UNP Q8I566
C	-23	MET	-	expression tag	UNP Q8I566
C	-22	THR	-	expression tag	UNP Q8I566
C	-21	GLY	-	expression tag	UNP Q8I566
C	-20	GLY	-	expression tag	UNP Q8I566
C	-19	GLN	-	expression tag	UNP Q8I566
C	-18	GLN	-	expression tag	UNP Q8I566
C	-17	MET	-	expression tag	UNP Q8I566
C	-16	GLY	-	expression tag	UNP Q8I566
C	-15	ARG	-	expression tag	UNP Q8I566
C	-14	ASP	-	expression tag	UNP Q8I566
C	-13	LEU	-	expression tag	UNP Q8I566
C	-12	TYR	-	expression tag	UNP Q8I566
C	-11	ASP	-	expression tag	UNP Q8I566

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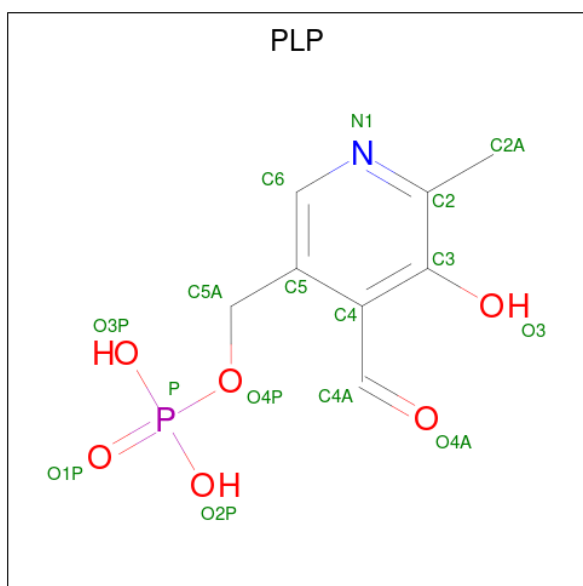
Chain	Residue	Modelled	Actual	Comment	Reference
C	-10	ASP	-	expression tag	UNP Q8I566
C	-9	ASP	-	expression tag	UNP Q8I566
C	-8	ASP	-	expression tag	UNP Q8I566
C	-7	LYS	-	expression tag	UNP Q8I566
C	-6	ASP	-	expression tag	UNP Q8I566
C	-5	HIS	-	expression tag	UNP Q8I566
C	-4	PRO	-	expression tag	UNP Q8I566
C	-3	PHE	-	expression tag	UNP Q8I566
C	-2	THR	-	expression tag	UNP Q8I566
C	-1	PRO	-	expression tag	UNP Q8I566
C	0	GLY	-	expression tag	UNP Q8I566
C	292	GLU	PHE	engineered mutation	UNP Q8I566
D	-37	MET	-	expression tag	UNP Q8I566
D	-36	ARG	-	expression tag	UNP Q8I566
D	-35	GLY	-	expression tag	UNP Q8I566
D	-34	SER	-	expression tag	UNP Q8I566
D	-33	HIS	-	expression tag	UNP Q8I566
D	-32	HIS	-	expression tag	UNP Q8I566
D	-31	HIS	-	expression tag	UNP Q8I566
D	-30	HIS	-	expression tag	UNP Q8I566
D	-29	HIS	-	expression tag	UNP Q8I566
D	-28	HIS	-	expression tag	UNP Q8I566
D	-27	GLY	-	expression tag	UNP Q8I566
D	-26	MET	-	expression tag	UNP Q8I566
D	-25	ALA	-	expression tag	UNP Q8I566
D	-24	SER	-	expression tag	UNP Q8I566
D	-23	MET	-	expression tag	UNP Q8I566
D	-22	THR	-	expression tag	UNP Q8I566
D	-21	GLY	-	expression tag	UNP Q8I566
D	-20	GLY	-	expression tag	UNP Q8I566
D	-19	GLN	-	expression tag	UNP Q8I566
D	-18	GLN	-	expression tag	UNP Q8I566
D	-17	MET	-	expression tag	UNP Q8I566
D	-16	GLY	-	expression tag	UNP Q8I566
D	-15	ARG	-	expression tag	UNP Q8I566
D	-14	ASP	-	expression tag	UNP Q8I566
D	-13	LEU	-	expression tag	UNP Q8I566
D	-12	TYR	-	expression tag	UNP Q8I566
D	-11	ASP	-	expression tag	UNP Q8I566
D	-10	ASP	-	expression tag	UNP Q8I566
D	-9	ASP	-	expression tag	UNP Q8I566
D	-8	ASP	-	expression tag	UNP Q8I566

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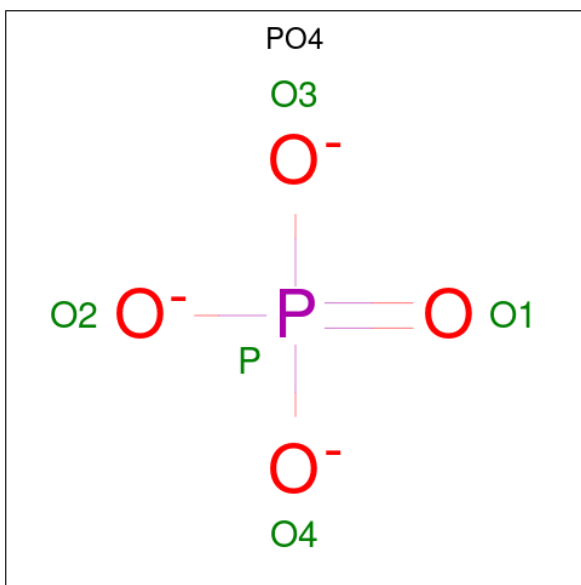
Chain	Residue	Modelled	Actual	Comment	Reference
D	-7	LYS	-	expression tag	UNP Q8I566
D	-6	ASP	-	expression tag	UNP Q8I566
D	-5	HIS	-	expression tag	UNP Q8I566
D	-4	PRO	-	expression tag	UNP Q8I566
D	-3	PHE	-	expression tag	UNP Q8I566
D	-2	THR	-	expression tag	UNP Q8I566
D	-1	PRO	-	expression tag	UNP Q8I566
D	0	GLY	-	expression tag	UNP Q8I566
D	292	GLU	PHE	engineered mutation	UNP Q8I566

- Molecule 2 is PYRIDOXAL-5'-PHOSPHATE (CCD ID: PLP) (formula: $C_8H_{10}NO_6P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			15	8	1	5	1		
2	B	1	Total	C	N	O	P	0	0
			15	8	1	5	1		
2	C	1	Total	C	N	O	P	0	0
			15	8	1	5	1		

- Molecule 3 is PHOSPHATE ION (CCD ID: PO4) (formula: O_4P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	D	1	Total	O	P	0	0
			5	4	1		

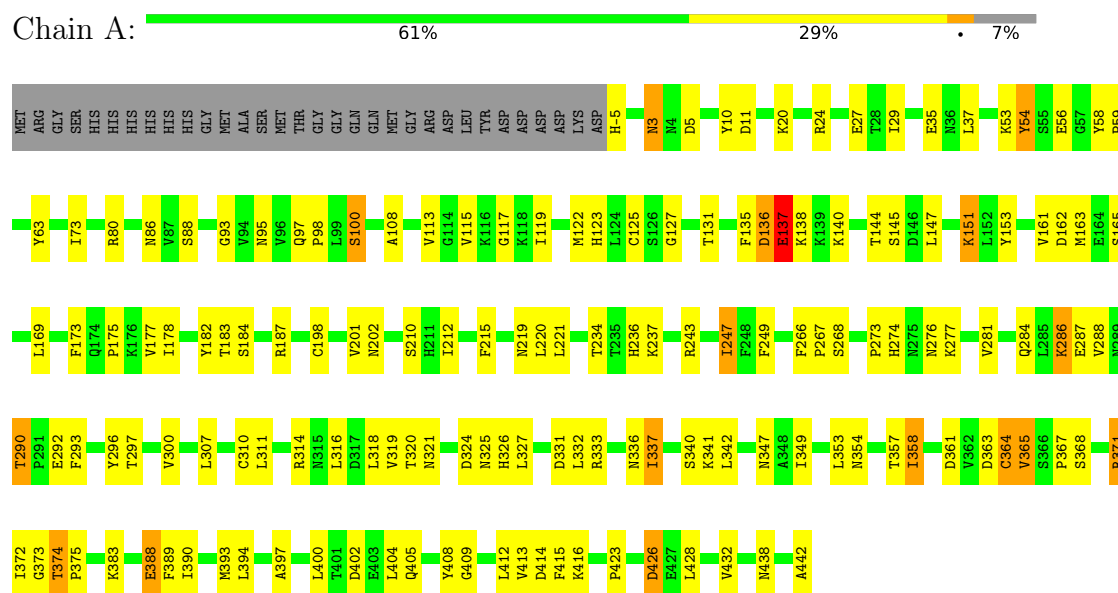
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	56	Total	O	0	0
			56	56		
4	B	55	Total	O	0	0
			55	55		
4	C	20	Total	O	0	0
			20	20		
4	D	7	Total	O	0	0
			7	7		

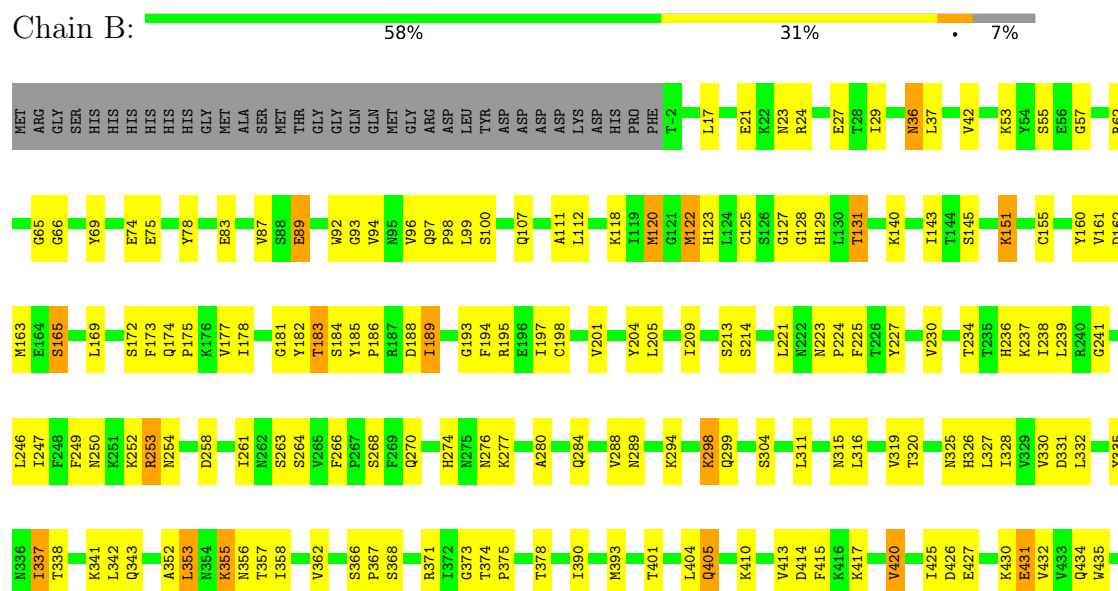
3 Residue-property plots

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

• Molecule 1: Serine hydroxymethyltransferase

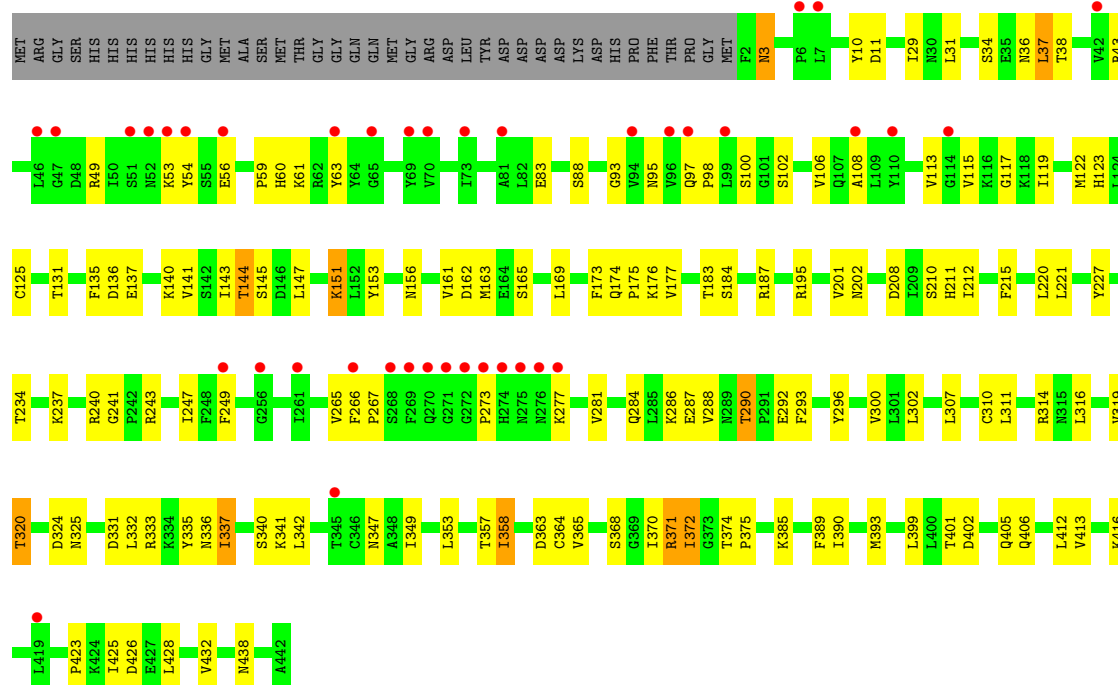


• Molecule 1: Serine hydroxymethyltransferase

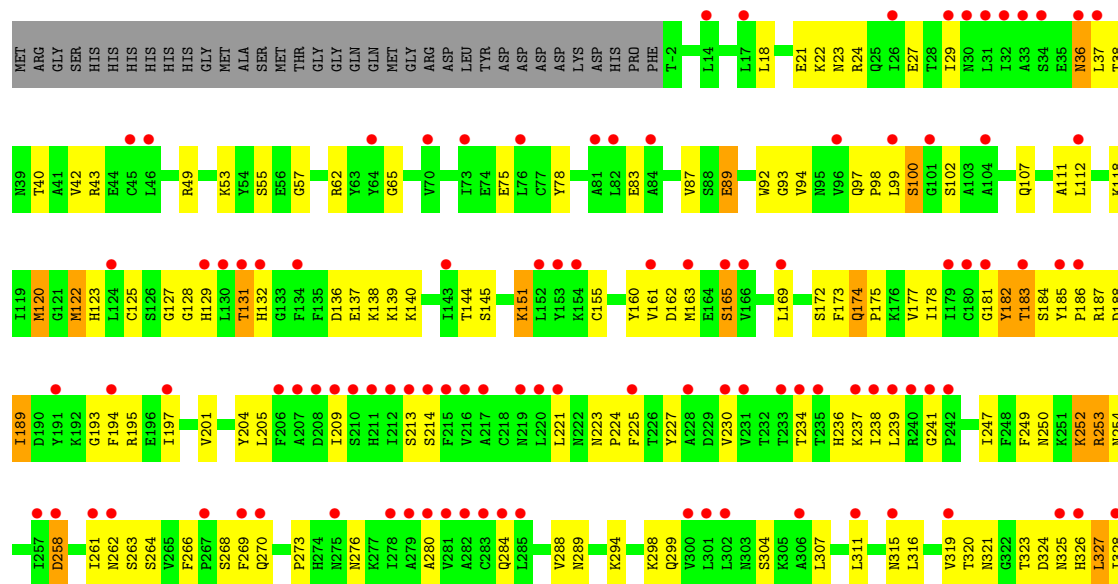


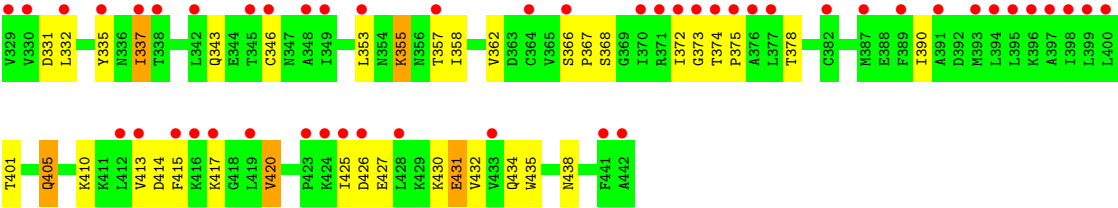


• Molecule 1: Serine hydroxymethyltransferase



• Molecule 1: Serine hydroxymethyltransferase





4 Data and refinement statistics

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, α , β , γ	254.95Å 254.95Å 61.40Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.85 – 2.98 29.85 – 2.98	Depositor EDS
% Data completeness (in resolution range)	99.7 (29.85-2.98) 99.8 (29.85-2.98)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	9.03 (at 3.00Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.224 , 0.270 0.216 , 0.262	Depositor DCC
R_{free} test set	4743 reflections (10.00%)	wwPDB-VP
Wilson B-factor (Å ²)	63.6	Xtriage
Anisotropy	0.253	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 56.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.021 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	14254	wwPDB-VP
Average B, all atoms (Å ²)	70.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.09% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PLP, 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	0.89	0/3612	1.07	7/4878 (0.1%)
1	B	0.90	0/3581	1.09	3/4835 (0.1%)
1	C	0.77	0/3554	1.00	2/4798 (0.0%)
1	D	0.68	0/3581	1.00	2/4835 (0.0%)
All	All	0.81	0/14328	1.04	14/19346 (0.1%)

There are no bond length outliers.

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	137	GLU	CB-CA-C	-6.80	108.74	116.63
1	B	294	LYS	N-CA-C	-5.82	104.53	111.69
1	A	325	ASN	N-CA-C	5.76	115.62	108.19
1	C	325	ASN	N-CA-C	5.75	115.60	108.19
1	A	247	ILE	CB-CA-C	-5.73	104.73	111.09
1	B	330	VAL	CB-CA-C	-5.60	103.97	110.91
1	A	5	ASP	CA-C-N	5.42	125.72	119.92
1	A	5	ASP	C-N-CA	5.42	125.72	119.92
1	B	353	LEU	N-CA-C	-5.34	100.02	108.14
1	D	294	LYS	N-CA-C	-5.32	105.14	111.69
1	D	182	TYR	N-CA-C	5.21	116.90	109.14
1	A	137	GLU	CB-CA-C	-5.12	110.69	116.63
1	A	364	CYS	N-CA-C	-5.06	106.93	113.01
1	A	409	GLY	N-CA-C	5.04	116.73	111.95

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3544	0	3558	138	0
1	B	3516	0	3537	141	0
1	C	3490	0	3508	122	0
1	D	3516	0	3539	143	0
2	A	15	0	6	1	0
2	B	15	0	7	1	0
2	C	15	0	6	3	0
3	D	5	0	0	2	0
4	A	56	0	0	6	0
4	B	55	0	0	2	0
4	C	20	0	0	3	0
4	D	7	0	0	1	0
All	All	14254	0	14161	520	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 18.

All (520) 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:337:ILE:HD11	1:D:401:THR:HB	1.29	1.14
1:A:122:MET:HE1	1:A:161:VAL:HG22	1.28	1.12
1:C:122:MET:HE1	1:C:161:VAL:HG22	1.29	1.08
1:B:337:ILE:HD11	1:B:401:THR:HB	1.22	1.07
1:A:151:LYS:HE3	1:A:173:PHE:CD1	1.90	1.07
1:C:290:THR:HG22	1:C:293:PHE:H	1.20	1.06
1:A:290:THR:HG22	1:A:293:PHE:H	1.17	1.05
1:A:268:SER:HB2	4:A:836:HOH:O	1.58	1.04
1:A:113:VAL:CG1	1:A:117:GLY:HA3	1.89	1.02
1:C:113:VAL:CG1	1:C:117:GLY:HA3	1.90	1.00
1:C:151:LYS:HE3	1:C:173:PHE:CD1	1.95	1.00
1:D:346:CYS:SG	1:D:353:LEU:HD21	2.06	0.95
1:D:250:ASN:ND2	1:D:253:ARG:HD2	1.82	0.94
1:A:219:ASN:HB2	1:C:399:LEU:CD2	1.98	0.93
1:A:184:SER:HA	1:A:327:LEU:HD21	1.50	0.92

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:337:ILE:HD11	1:B:401:THR:CB	2.00	0.92
1:D:89:GLU:HG2	1:D:89:GLU:O	1.69	0.91
1:B:250:ASN:ND2	1:B:253:ARG:HD2	1.87	0.89
1:A:151:LYS:HE3	1:A:173:PHE:HD1	1.34	0.88
1:C:311:LEU:HD22	1:C:316:LEU:HD12	1.56	0.87
1:D:337:ILE:HD11	1:D:401:THR:CB	2.04	0.87
1:B:123:HIS:HD2	1:B:125:CYS:H	1.21	0.86
1:C:151:LYS:HE3	1:C:173:PHE:HD1	1.39	0.85
1:A:287:GLU:O	1:A:290:THR:HB	1.78	0.84
1:C:287:GLU:O	1:C:290:THR:HB	1.79	0.82
1:C:358:ILE:HD11	1:C:368:SER:HB2	1.60	0.82
1:A:311:LEU:HD22	1:A:316:LEU:HD12	1.60	0.81
1:A:122:MET:CE	1:A:161:VAL:HG22	2.09	0.81
1:D:177:VAL:HG22	1:D:204:TYR:HB2	1.61	0.81
1:B:89:GLU:O	1:B:89:GLU:HG2	1.75	0.81
1:B:316:LEU:HD23	1:B:335:TYR:OH	1.80	0.81
1:A:113:VAL:HG13	1:A:117:GLY:HA3	1.62	0.81
1:B:177:VAL:HG22	1:B:204:TYR:HB2	1.63	0.80
1:A:374:THR:N	1:A:375:PRO:HD3	1.98	0.79
1:B:36:ASN:HD22	1:B:37:LEU:H	1.31	0.79
1:D:123:HIS:HD2	1:D:125:CYS:H	1.27	0.79
1:B:414:ASP:O	1:B:417:LYS:HB3	1.84	0.78
1:D:316:LEU:HD23	1:D:335:TYR:OH	1.84	0.78
1:B:123:HIS:CD2	1:B:125:CYS:H	2.02	0.77
1:D:417:LYS:O	1:D:420:VAL:HG23	1.84	0.77
1:C:122:MET:CE	1:C:161:VAL:HG22	2.12	0.77
1:A:219:ASN:HB2	1:C:399:LEU:HD22	1.65	0.77
1:D:36:ASN:HD22	1:D:37:LEU:H	1.30	0.76
1:A:358:ILE:HD11	1:A:368:SER:HB2	1.66	0.76
1:B:163:MET:HE1	1:B:194:PHE:CE2	2.20	0.75
1:B:358:ILE:CD1	1:B:368:SER:HB2	2.17	0.74
1:D:127:GLY:O	1:D:182:TYR:HB3	1.85	0.74
1:A:113:VAL:HG12	1:A:117:GLY:HA3	1.70	0.74
1:B:417:LYS:O	1:B:420:VAL:HG23	1.88	0.73
1:B:97:GLN:N	1:B:98:PRO:HD3	2.03	0.73
1:C:125:CYS:HB3	1:C:357:THR:HG21	1.68	0.73
1:B:332:LEU:HB3	1:B:337:ILE:HG22	1.70	0.73
1:C:113:VAL:HG13	1:C:117:GLY:HA3	1.69	0.73
1:B:89:GLU:O	1:B:89:GLU:CG	2.36	0.72
1:C:113:VAL:HG12	1:C:117:GLY:HA3	1.69	0.72
1:D:250:ASN:HD22	1:D:253:ARG:HD2	1.52	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:414:ASP:O	1:D:417:LYS:HB3	1.89	0.72
1:B:319:VAL:HG12	1:B:320:THR:HG23	1.72	0.71
1:D:97:GLN:N	1:D:98:PRO:HD3	2.06	0.71
1:A:140:LYS:HB3	1:A:145:SER:OG	1.91	0.71
1:D:89:GLU:O	1:D:89:GLU:CG	2.36	0.71
1:B:358:ILE:HD11	1:B:368:SER:HB2	1.73	0.71
1:D:311:LEU:HD22	1:D:316:LEU:CD1	2.21	0.71
1:C:119:ILE:HG22	1:C:177:VAL:CG1	2.21	0.70
1:A:122:MET:HE1	1:A:161:VAL:CG2	2.16	0.70
1:A:327:LEU:HD11	1:A:371:ARG:HD2	1.72	0.70
1:A:86:ASN:O	1:C:406:GLN:NE2	2.25	0.70
1:D:253:ARG:HG2	1:D:253:ARG:HH11	1.56	0.70
1:D:258:ASP:OD1	1:D:258:ASP:C	2.33	0.70
1:D:358:ILE:CD1	1:D:368:SER:HB2	2.22	0.70
1:B:127:GLY:O	1:B:182:TYR:HB3	1.91	0.69
1:A:423:PRO:HA	1:A:426:ASP:OD2	1.92	0.69
1:A:363:ASP:CG	1:A:365:VAL:HB	2.18	0.69
1:C:310:CYS:O	1:C:314:ARG:HD2	1.92	0.69
1:A:219:ASN:CB	1:C:399:LEU:CD2	2.70	0.69
1:C:3:ASN:HB3	1:C:10:TYR:CE2	2.27	0.69
1:D:319:VAL:HG12	1:D:320:THR:HG23	1.74	0.69
1:D:163:MET:HE1	1:D:194:PHE:CE2	2.27	0.69
1:D:205:LEU:HD23	1:D:227:TYR:O	1.93	0.69
1:B:87:VAL:HB	1:B:92:TRP:CD1	2.28	0.68
1:D:120:MET:HE2	1:D:178:ILE:HG12	1.72	0.68
1:A:371:ARG:HG3	1:A:371:ARG:HH11	1.58	0.68
1:B:258:ASP:OD1	1:B:258:ASP:C	2.35	0.68
1:D:93:GLY:HA3	1:D:249:PHE:CE1	2.30	0.67
1:D:123:HIS:CD2	1:D:125:CYS:H	2.10	0.67
1:C:119:ILE:HG22	1:C:177:VAL:HG12	1.77	0.67
1:C:371:ARG:HG3	1:C:371:ARG:HH11	1.59	0.67
1:C:108:ALA:HB2	1:C:247:ILE:HD13	1.76	0.67
1:B:120:MET:HE2	1:B:178:ILE:HG12	1.77	0.67
1:C:140:LYS:HB3	1:C:145:SER:OG	1.95	0.67
1:B:161:VAL:CG1	1:B:163:MET:HE2	2.25	0.67
1:C:290:THR:CG2	1:C:293:PHE:H	2.03	0.67
1:B:253:ARG:HH11	1:B:253:ARG:HG2	1.59	0.66
1:A:119:ILE:HG22	1:A:177:VAL:CG1	2.26	0.66
1:A:268:SER:O	1:B:143:ILE:HG22	1.95	0.66
1:D:161:VAL:CG1	1:D:163:MET:HE2	2.26	0.66
1:D:332:LEU:HB3	1:D:337:ILE:HG22	1.77	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:3:ASN:HB3	1:A:10:TYR:CE2	2.31	0.65
1:A:125:CYS:HB3	1:A:357:THR:HG21	1.78	0.65
1:B:205:LEU:HD23	1:B:227:TYR:O	1.96	0.65
1:D:87:VAL:HB	1:D:92:TRP:CD1	2.32	0.65
1:A:119:ILE:HG22	1:A:177:VAL:HG12	1.78	0.65
1:C:141:VAL:O	1:D:268:SER:HA	1.96	0.65
1:B:311:LEU:HD22	1:B:316:LEU:CD1	2.27	0.65
1:B:194:PHE:HB3	1:B:205:LEU:HD13	1.79	0.65
1:C:423:PRO:HA	1:C:426:ASP:OD2	1.96	0.65
1:D:223:ASN:OD1	1:D:225:PHE:N	2.27	0.65
1:C:363:ASP:C	1:C:365:VAL:H	2.05	0.64
1:D:326:HIS:O	1:D:373:GLY:HA2	1.97	0.64
1:C:3:ASN:CB	1:C:10:TYR:CE2	2.80	0.64
1:C:3:ASN:CB	1:C:10:TYR:HE2	2.11	0.64
1:A:412:LEU:O	1:A:416:LYS:HB2	1.97	0.63
1:A:332:LEU:HB2	1:A:368:SER:O	1.98	0.63
1:A:290:THR:CG2	1:A:293:PHE:H	2.04	0.63
1:D:426:ASP:O	1:D:430:LYS:HG3	1.98	0.63
1:C:371:ARG:HG3	4:C:810:HOH:O	1.97	0.63
1:B:223:ASN:OD1	1:B:225:PHE:N	2.28	0.63
1:A:202:ASN:ND2	1:A:202:ASN:O	2.32	0.63
1:B:93:GLY:HA3	1:B:249:PHE:CE1	2.34	0.63
1:D:311:LEU:HB3	1:D:316:LEU:HD12	1.81	0.63
1:C:201:VAL:HG12	1:C:201:VAL:O	1.97	0.62
1:C:122:MET:HE1	1:C:161:VAL:CG2	2.18	0.62
1:B:120:MET:CE	1:B:178:ILE:HG12	2.30	0.62
1:A:319:VAL:CG2	1:A:358:ILE:HG22	2.30	0.62
1:A:97:GLN:HG2	1:A:273:PRO:HB3	1.82	0.61
1:B:362:VAL:HG12	1:B:362:VAL:O	2.00	0.61
1:C:412:LEU:O	1:C:416:LYS:HB2	2.01	0.61
1:C:374:THR:N	1:C:375:PRO:HD3	2.15	0.61
1:A:290:THR:HG23	1:A:292:GLU:H	1.65	0.61
1:A:310:CYS:O	1:A:314:ARG:HD2	2.01	0.61
1:A:108:ALA:HB2	1:A:247:ILE:HD13	1.83	0.61
1:B:241:GLY:CA	1:B:284:GLN:HG2	2.31	0.61
1:B:122:MET:HE2	1:B:182:TYR:CD1	2.36	0.61
1:A:201:VAL:HG12	1:A:201:VAL:O	2.00	0.60
1:A:374:THR:N	1:A:375:PRO:CD	2.63	0.60
1:B:337:ILE:CD1	1:B:401:THR:HB	2.14	0.60
1:B:311:LEU:HB3	1:B:316:LEU:HD12	1.81	0.60
1:D:241:GLY:CA	1:D:284:GLN:HG2	2.31	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:211:HIS:HD2	2:C:701:PLP:O3	1.84	0.59
1:B:161:VAL:HG11	1:B:163:MET:HE2	1.84	0.59
1:D:236:HIS:O	1:D:237:LYS:HB2	2.02	0.59
1:B:250:ASN:HD22	1:B:253:ARG:HD2	1.65	0.59
1:D:183:THR:HG22	1:D:184:SER:HB3	1.83	0.59
1:A:113:VAL:CG1	1:A:117:GLY:CA	2.76	0.59
1:C:358:ILE:HD11	1:C:368:SER:CB	2.31	0.59
1:A:284:GLN:O	1:A:288:VAL:HG23	2.03	0.59
1:A:24:ARG:HH21	1:B:66:GLY:HA3	1.68	0.59
1:D:319:VAL:C	1:D:320:THR:HG23	2.28	0.59
1:B:122:MET:HE2	1:B:182:TYR:CE1	2.37	0.59
1:B:29:ILE:HD12	1:B:432:VAL:HG13	1.85	0.58
1:C:290:THR:HG23	1:C:292:GLU:H	1.68	0.58
1:B:182:TYR:CE2	1:B:189:ILE:HD13	2.37	0.58
1:D:435:TRP:O	1:D:438:ASN:HB3	2.04	0.58
1:B:36:ASN:ND2	1:B:37:LEU:H	2.02	0.58
1:A:220:LEU:O	1:A:221:LEU:HD23	2.04	0.58
1:B:75:GLU:HA	1:B:78:TYR:CD2	2.38	0.58
1:B:326:HIS:O	1:B:373:GLY:HA2	2.03	0.58
1:D:337:ILE:HD11	1:D:401:THR:CG2	2.34	0.58
1:D:129:HIS:ND1	1:D:131:THR:OG1	2.32	0.58
1:D:358:ILE:HD11	1:D:368:SER:HB2	1.85	0.57
1:D:362:VAL:HG12	1:D:362:VAL:O	2.02	0.57
1:B:97:GLN:N	1:B:98:PRO:CD	2.67	0.57
1:A:162:ASP:OD2	1:A:165:SER:N	2.37	0.57
1:C:374:THR:N	1:C:375:PRO:CD	2.67	0.57
1:B:29:ILE:CD1	1:B:432:VAL:HG13	2.34	0.57
1:C:363:ASP:C	1:C:365:VAL:N	2.61	0.57
1:D:311:LEU:HD22	1:D:316:LEU:HD12	1.87	0.57
1:B:371:ARG:HD3	4:B:823:HOH:O	2.04	0.57
1:D:161:VAL:HG11	1:D:163:MET:HE2	1.85	0.57
1:B:319:VAL:C	1:B:320:THR:HG23	2.30	0.57
1:D:75:GLU:HA	1:D:78:TYR:CD2	2.39	0.57
1:D:311:LEU:HD22	1:D:316:LEU:HD11	1.87	0.57
1:A:327:LEU:CD1	1:A:371:ARG:HD2	2.34	0.57
1:C:37:LEU:N	1:C:37:LEU:HD23	2.19	0.57
1:D:36:ASN:ND2	1:D:37:LEU:H	2.00	0.57
1:D:307:LEU:HD21	1:D:372:ILE:HD12	1.85	0.57
1:A:358:ILE:HD11	1:A:368:SER:CB	2.34	0.56
1:B:112:LEU:HD11	1:B:230:VAL:HG21	1.87	0.56
1:A:153:TYR:HB3	1:A:169:LEU:HD12	1.86	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:363:ASP:OD2	1:A:365:VAL:HB	2.03	0.56
1:D:122:MET:HE2	1:D:182:TYR:CD1	2.40	0.56
1:D:194:PHE:HB3	1:D:205:LEU:HD13	1.87	0.56
1:A:3:ASN:CB	1:A:10:TYR:CE2	2.89	0.56
1:B:21:GLU:OE2	1:B:24:ARG:NH1	2.39	0.56
1:C:202:ASN:O	1:C:202:ASN:ND2	2.38	0.56
1:B:357:THR:HG22	1:B:367:PRO:HB3	1.87	0.56
1:D:98:PRO:HA	1:D:270:GLN:HE22	1.69	0.56
1:D:343:GLN:OE1	1:D:355:LYS:HG2	2.06	0.56
1:A:162:ASP:HB3	1:A:165:SER:HB3	1.88	0.56
1:C:374:THR:H	1:C:375:PRO:HD3	1.71	0.56
1:C:113:VAL:CG1	1:C:117:GLY:CA	2.77	0.55
1:A:296:TYR:O	1:A:300:VAL:HG23	2.06	0.55
1:B:36:ASN:HD22	1:B:37:LEU:N	2.04	0.55
1:B:435:TRP:O	1:B:438:ASN:HB3	2.06	0.55
1:A:37:LEU:HD23	1:A:442:ALA:H	1.72	0.55
1:B:182:TYR:HE2	1:B:189:ILE:HD13	1.71	0.55
1:A:135:PHE:HB3	1:A:145:SER:HB2	1.88	0.55
1:B:263:SER:HA	1:B:266:PHE:O	2.06	0.55
1:C:162:ASP:OD2	1:C:165:SER:N	2.37	0.55
1:C:319:VAL:HG22	1:C:358:ILE:HG22	1.88	0.55
1:D:304:SER:OG	1:D:325:ASN:O	2.24	0.55
1:B:431:GLU:O	1:B:434:GLN:HG2	2.07	0.55
1:D:97:GLN:N	1:D:98:PRO:CD	2.70	0.55
1:A:319:VAL:HG22	1:A:358:ILE:HG22	1.88	0.55
1:B:57:GLY:O	1:B:62:ARG:NH1	2.40	0.55
1:D:337:ILE:CD1	1:D:401:THR:HB	2.19	0.55
1:B:123:HIS:HD2	1:B:125:CYS:N	2.00	0.55
1:C:153:TYR:HB3	1:C:169:LEU:HD12	1.87	0.55
1:B:140:LYS:HB3	1:B:145:SER:OG	2.07	0.55
1:C:284:GLN:O	1:C:288:VAL:HG23	2.07	0.55
1:B:284:GLN:O	1:B:288:VAL:HG23	2.07	0.54
1:A:97:GLN:N	1:A:98:PRO:CD	2.70	0.54
1:A:210:SER:HB3	1:A:234:THR:OG1	2.07	0.54
1:B:23:ASN:O	1:B:27:GLU:HG3	2.07	0.54
1:B:98:PRO:HA	1:B:270:GLN:HE22	1.71	0.54
1:C:135:PHE:CD1	1:C:135:PHE:C	2.85	0.54
1:B:311:LEU:HD22	1:B:316:LEU:HD11	1.89	0.54
1:C:296:TYR:O	1:C:300:VAL:HG23	2.06	0.54
1:B:326:HIS:HB2	1:B:375:PRO:HD3	1.90	0.54
1:C:49:ARG:NH1	1:D:22:LYS:HE2	2.22	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:337:ILE:HD11	1:B:401:THR:CG2	2.37	0.54
1:D:111:ALA:HB2	1:D:264:SER:HB2	1.89	0.54
1:D:434:GLN:HG3	1:D:435:TRP:N	2.22	0.54
1:D:284:GLN:O	1:D:288:VAL:HG23	2.07	0.54
1:C:122:MET:HE3	1:C:153:TYR:HE1	1.73	0.53
1:C:220:LEU:O	1:C:221:LEU:HD23	2.07	0.53
1:D:112:LEU:HD11	1:D:230:VAL:HG21	1.89	0.53
1:D:120:MET:CE	1:D:178:ILE:HG12	2.38	0.53
1:A:371:ARG:NE	4:A:804:HOH:O	2.41	0.53
1:B:111:ALA:HB2	1:B:264:SER:HB2	1.89	0.53
1:D:21:GLU:OE2	1:D:24:ARG:NH1	2.42	0.53
1:D:299:GLN:HB3	1:D:378:THR:HG23	1.90	0.53
1:D:57:GLY:O	1:D:62:ARG:NH1	2.42	0.53
1:D:263:SER:HA	1:D:266:PHE:O	2.08	0.53
1:A:135:PHE:C	1:A:135:PHE:CD1	2.87	0.53
1:B:107:GLN:HE22	1:B:270:GLN:NE2	2.06	0.53
1:D:151:LYS:HA	1:D:151:LYS:NZ	2.24	0.53
1:C:97:GLN:N	1:C:98:PRO:CD	2.72	0.52
1:C:135:PHE:HB3	1:C:145:SER:HB2	1.90	0.52
1:D:122:MET:HE2	1:D:182:TYR:CE1	2.44	0.52
1:D:362:VAL:O	1:D:362:VAL:CG1	2.57	0.52
1:A:122:MET:HE3	1:A:153:TYR:HE1	1.74	0.52
1:A:215:PHE:HB3	1:A:221:LEU:HG	1.91	0.52
1:A:337:ILE:HD11	1:A:402:ASP:CA	2.39	0.52
1:B:213:SER:OG	1:B:239:LEU:HB2	2.08	0.52
1:B:253:ARG:HH11	1:B:253:ARG:CG	2.21	0.52
1:B:343:GLN:OE1	1:B:355:LYS:HG2	2.09	0.52
1:D:29:ILE:HD12	1:D:432:VAL:HG13	1.92	0.52
1:B:122:MET:HG3	1:B:123:HIS:N	2.25	0.52
1:C:43:ARG:HH12	1:D:49:ARG:NH2	2.08	0.52
1:D:36:ASN:HD22	1:D:37:LEU:N	2.03	0.52
1:B:201:VAL:HG12	1:B:201:VAL:O	2.09	0.52
1:D:366:SER:N	1:D:367:PRO:HD3	2.25	0.52
1:D:188:ASP:OD2	1:D:221:LEU:HD22	2.10	0.52
1:B:223:ASN:OD1	1:B:225:PHE:HB2	2.10	0.51
1:C:337:ILE:HD11	1:C:402:ASP:CA	2.39	0.51
1:D:162:ASP:HB3	1:D:165:SER:HB3	1.91	0.51
1:D:29:ILE:CD1	1:D:432:VAL:HG13	2.40	0.51
1:D:366:SER:N	1:D:367:PRO:CD	2.72	0.51
1:C:332:LEU:HB2	1:C:368:SER:O	2.10	0.51
1:A:35:GLU:HA	4:A:819:HOH:O	2.09	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:213:SER:OG	1:D:239:LEU:HB2	2.10	0.51
1:C:277:LYS:O	1:C:281:VAL:HG23	2.10	0.51
1:D:100:SER:OG	3:D:501:PO4:O1	2.24	0.51
1:D:201:VAL:HG12	1:D:201:VAL:O	2.11	0.51
1:A:363:ASP:OD1	1:A:365:VAL:HB	2.11	0.51
1:A:234:THR:HB	1:A:236:HIS:CE1	2.46	0.51
1:A:243:ARG:O	1:A:277:LYS:HE3	2.10	0.51
1:A:331:ASP:OD1	1:A:333:ARG:HG2	2.11	0.51
1:C:389:PHE:CZ	1:C:393:MET:HE3	2.46	0.51
1:B:151:LYS:HA	1:B:151:LYS:NZ	2.25	0.51
1:D:323:THR:OG1	1:D:327:LEU:O	2.24	0.51
1:B:183:THR:HG22	1:B:184:SER:HB3	1.93	0.50
1:D:107:GLN:HE22	1:D:270:GLN:NE2	2.08	0.50
1:A:219:ASN:HB2	1:C:399:LEU:HD21	1.88	0.50
1:C:162:ASP:HA	4:C:816:HOH:O	2.12	0.50
1:D:42:VAL:HG13	1:D:280:ALA:HB1	1.93	0.50
1:B:311:LEU:HD22	1:B:316:LEU:HD12	1.92	0.50
1:B:426:ASP:O	1:B:430:LYS:HG3	2.11	0.50
1:A:123:HIS:CD2	1:A:125:CYS:H	2.28	0.50
1:A:336:ASN:O	1:A:337:ILE:HG13	2.10	0.50
1:B:366:SER:N	1:B:367:PRO:HD3	2.26	0.50
1:C:349:ILE:HG21	1:C:428:LEU:HB2	1.92	0.50
1:B:83:GLU:HG3	1:B:289:ASN:ND2	2.27	0.50
1:B:129:HIS:ND1	1:B:131:THR:OG1	2.39	0.50
1:D:36:ASN:C	1:D:37:LEU:HD12	2.37	0.50
1:B:188:ASP:OD2	1:B:221:LEU:HD22	2.12	0.50
1:B:21:GLU:OE1	1:B:21:GLU:HA	2.12	0.49
1:C:131:THR:O	1:C:144:THR:HG21	2.11	0.49
1:C:341:LYS:HE2	1:C:405:GLN:HG3	1.93	0.49
1:D:140:LYS:HB3	1:D:145:SER:OG	2.13	0.49
1:B:366:SER:N	1:B:367:PRO:CD	2.74	0.49
1:B:162:ASP:HB3	1:B:165:SER:HB3	1.93	0.49
1:D:374:THR:N	1:D:375:PRO:CD	2.74	0.49
1:A:80:ARG:NH2	1:A:286:LYS:HD3	2.27	0.49
1:A:274:HIS:HD2	1:B:277:LYS:HZ3	1.59	0.49
1:A:29:ILE:HD12	1:A:432:VAL:HG13	1.94	0.49
1:A:277:LYS:O	1:A:281:VAL:HG23	2.11	0.49
1:B:120:MET:SD	1:B:169:LEU:HD23	2.53	0.49
1:C:358:ILE:CD1	1:C:368:SER:HB2	2.36	0.49
1:D:83:GLU:HG3	1:D:289:ASN:ND2	2.28	0.49
1:D:155:CYS:HB2	1:D:160:TYR:O	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:337:ILE:HD11	1:A:402:ASP:CB	2.43	0.49
1:B:174:GLN:N	1:B:175:PRO:CD	2.76	0.49
1:A:389:PHE:CZ	1:A:393:MET:HE3	2.47	0.49
1:A:3:ASN:CB	1:A:10:TYR:HE2	2.26	0.48
1:A:311:LEU:HD22	1:A:316:LEU:CD1	2.37	0.48
1:C:36:ASN:C	1:C:37:LEU:HD23	2.38	0.48
1:C:97:GLN:HG2	1:C:273:PRO:HB3	1.95	0.48
1:D:326:HIS:HB2	1:D:375:PRO:HD3	1.94	0.48
1:A:187:ARG:HA	1:A:324:ASP:HB2	1.95	0.48
1:A:331:ASP:OD1	1:A:331:ASP:C	2.54	0.48
1:C:210:SER:HB3	1:C:234:THR:OG1	2.12	0.48
1:D:223:ASN:OD1	1:D:223:ASN:C	2.56	0.48
1:A:318:LEU:O	1:A:321:ASN:N	2.43	0.48
1:B:97:GLN:H	1:B:98:PRO:HD3	1.76	0.48
1:B:118:LYS:HB3	1:B:173:PHE:CE2	2.48	0.48
1:C:347:ASN:HD21	1:D:65:GLY:HA3	1.78	0.48
1:B:223:ASN:OD1	1:B:223:ASN:C	2.55	0.48
1:C:208:ASP:OD1	2:C:701:PLP:H2A2	2.14	0.48
1:D:405:GLN:HE21	1:D:410:LYS:HG3	1.78	0.48
1:C:161:VAL:CG1	1:C:163:MET:CE	2.92	0.48
1:D:97:GLN:HG2	1:D:273:PRO:HB3	1.96	0.48
1:B:332:LEU:HB3	1:B:337:ILE:CG2	2.41	0.48
1:C:208:ASP:OD1	2:C:701:PLP:C2A	2.61	0.48
1:D:21:GLU:HA	1:D:21:GLU:OE1	2.14	0.48
1:D:174:GLN:N	1:D:175:PRO:CD	2.77	0.48
1:A:123:HIS:HD2	1:A:125:CYS:H	1.62	0.47
1:A:173:PHE:CE2	1:A:175:PRO:HB3	2.49	0.47
1:C:336:ASN:O	1:C:337:ILE:HG13	2.14	0.47
1:D:99:LEU:H	1:D:270:GLN:HE22	1.61	0.47
1:A:365:VAL:O	1:A:367:PRO:HD3	2.14	0.47
1:D:23:ASN:O	1:D:27:GLU:HG3	2.13	0.47
1:A:131:THR:O	1:A:144:THR:HG21	2.14	0.47
1:B:36:ASN:C	1:B:37:LEU:HD12	2.40	0.47
1:D:151:LYS:HA	1:D:151:LYS:HZ3	1.80	0.47
1:A:73:ILE:HD11	1:B:17:LEU:HB3	1.96	0.47
1:A:93:GLY:HA3	1:A:249:PHE:CE1	2.50	0.47
1:A:375:PRO:HD2	4:A:848:HOH:O	2.14	0.47
1:C:195:ARG:HD3	1:C:227:TYR:O	2.15	0.47
1:C:29:ILE:HD12	1:C:432:VAL:HG13	1.96	0.47
1:C:123:HIS:CD2	1:C:125:CYS:H	2.33	0.47
1:C:319:VAL:HG12	1:C:320:THR:HG23	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:341:LYS:HE2	1:A:405:GLN:HG3	1.97	0.47
1:B:161:VAL:HG12	1:B:163:MET:HE2	1.96	0.47
1:B:331:ASP:OD1	1:B:331:ASP:C	2.57	0.47
1:A:63:TYR:N	1:A:63:TYR:CD1	2.82	0.47
1:D:221:LEU:HB3	4:D:607:HOH:O	2.15	0.47
1:B:304:SER:OG	1:B:325:ASN:O	2.27	0.46
1:C:135:PHE:CD1	1:C:135:PHE:O	2.68	0.46
1:B:94:VAL:HA	1:B:247:ILE:O	2.16	0.46
1:B:374:THR:N	1:B:375:PRO:CD	2.78	0.46
1:B:276:ASN:OD1	1:B:276:ASN:N	2.48	0.46
1:C:331:ASP:OD1	1:C:333:ARG:HG2	2.14	0.46
1:A:54:TYR:HA	4:A:829:HOH:O	2.15	0.46
1:A:219:ASN:CB	1:C:399:LEU:HD21	2.44	0.46
1:A:337:ILE:HD11	1:A:402:ASP:HB2	1.97	0.46
1:B:401:THR:HG23	1:B:415:PHE:CZ	2.50	0.46
1:C:162:ASP:HB3	1:C:165:SER:HB3	1.97	0.46
1:A:326:HIS:O	1:A:373:GLY:HA2	2.16	0.46
1:A:307:LEU:HD11	1:A:390:ILE:HG22	1.98	0.46
1:B:311:LEU:CB	1:B:316:LEU:HD12	2.45	0.46
1:D:401:THR:HG23	1:D:415:PHE:CZ	2.50	0.46
1:D:182:TYR:HE2	1:D:189:ILE:HD13	1.81	0.46
1:D:258:ASP:OD1	1:D:262:ASN:ND2	2.46	0.46
1:A:332:LEU:HD11	1:A:342:LEU:HD22	1.98	0.46
1:B:96:VAL:HG12	1:B:246:LEU:CD2	2.46	0.46
1:B:197:ILE:O	1:B:197:ILE:HG22	2.16	0.46
1:C:187:ARG:HA	1:C:324:ASP:HB2	1.97	0.46
1:B:332:LEU:HD11	1:B:342:LEU:HD23	1.98	0.45
1:A:219:ASN:HB3	1:C:335:TYR:HE2	1.80	0.45
1:B:393:MET:HE2	1:B:432:VAL:HG23	1.98	0.45
1:C:115:VAL:HG23	1:C:147:LEU:HD12	1.98	0.45
1:C:371:ARG:HH11	1:C:371:ARG:CG	2.28	0.45
1:A:63:TYR:N	1:A:63:TYR:HD1	2.14	0.45
1:A:219:ASN:HD22	1:C:399:LEU:HD23	1.82	0.45
1:B:87:VAL:HG11	1:B:225:PHE:CD1	2.51	0.45
1:B:99:LEU:H	1:B:270:GLN:HE22	1.64	0.45
1:B:193:GLY:O	1:B:197:ILE:HD12	2.16	0.45
1:C:319:VAL:CG2	1:C:358:ILE:HG22	2.45	0.45
1:D:182:TYR:CE2	1:D:189:ILE:HD13	2.51	0.45
1:A:202:ASN:O	1:A:202:ASN:CG	2.60	0.45
1:C:162:ASP:OD2	1:C:165:SER:CB	2.63	0.45
1:D:118:LYS:HB3	1:D:173:PHE:CE2	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:178:ILE:HD12	1:B:198:CYS:SG	2.57	0.45
1:C:243:ARG:O	1:C:277:LYS:HE3	2.16	0.45
1:A:100:SER:OG	1:B:274:HIS:HE1	2.00	0.45
1:B:42:VAL:HG13	1:B:280:ALA:HB1	1.98	0.45
1:C:401:THR:O	1:C:405:GLN:HB2	2.17	0.45
1:A:347:ASN:ND2	1:B:65:GLY:HA3	2.32	0.45
1:B:155:CYS:HB2	1:B:160:TYR:O	2.17	0.45
1:D:223:ASN:OD1	1:D:225:PHE:HB2	2.16	0.45
1:C:215:PHE:HB3	1:C:221:LEU:HG	1.99	0.45
1:D:173:PHE:HB3	1:D:175:PRO:HD3	1.99	0.45
1:D:237:LYS:HB3	1:D:238:ILE:H	1.65	0.45
1:A:371:ARG:HH11	1:A:371:ARG:CG	2.25	0.45
1:C:290:THR:HG23	1:C:292:GLU:N	2.32	0.45
1:D:311:LEU:CB	1:D:316:LEU:HD12	2.45	0.44
1:C:95:ASN:ND2	1:C:265:VAL:HG21	2.32	0.44
1:D:94:VAL:HA	1:D:247:ILE:O	2.16	0.44
1:D:185:TYR:HA	1:D:186:PRO:HD3	1.84	0.44
1:C:93:GLY:HA3	1:C:249:PHE:CE1	2.52	0.44
1:D:123:HIS:HD2	1:D:125:CYS:N	2.05	0.44
1:A:347:ASN:HD21	1:B:65:GLY:HA3	1.83	0.44
1:B:173:PHE:HB3	1:B:175:PRO:HD3	1.99	0.44
1:D:122:MET:HG3	1:D:123:HIS:N	2.33	0.44
1:A:161:VAL:CG1	1:A:163:MET:CE	2.96	0.44
1:C:102:SER:HA	1:C:131:THR:HG21	1.99	0.44
1:C:358:ILE:H	1:C:358:ILE:HG12	1.63	0.44
1:D:183:THR:HG22	1:D:184:SER:N	2.33	0.44
1:C:201:VAL:O	1:C:201:VAL:CG1	2.65	0.44
1:B:183:THR:HG22	1:B:184:SER:N	2.33	0.43
1:C:143:ILE:HG12	1:D:269:PHE:CE1	2.53	0.43
1:C:363:ASP:O	1:C:365:VAL:N	2.51	0.43
1:C:63:TYR:CD1	1:C:63:TYR:N	2.86	0.43
1:C:174:GLN:HG3	1:C:201:VAL:HG13	1.99	0.43
1:D:332:LEU:HB3	1:D:337:ILE:CG2	2.47	0.43
1:A:274:HIS:CD2	1:B:277:LYS:NZ	2.86	0.43
1:B:404:LEU:HA	1:B:404:LEU:HD23	1.77	0.43
1:C:173:PHE:CE2	1:C:175:PRO:HB3	2.53	0.43
1:D:276:ASN:N	1:D:276:ASN:OD1	2.50	0.43
1:B:338:THR:OG1	1:B:341:LYS:HG3	2.19	0.43
1:D:187:ARG:NH2	1:D:320:THR:O	2.46	0.43
1:A:394:LEU:O	1:A:397:ALA:HB3	2.17	0.43
1:D:252:LYS:HB2	1:D:252:LYS:HE2	1.70	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:209:ILE:HG21	1:B:224:PRO:HG3	2.01	0.43
1:D:38:THR:O	1:D:43:ARG:NH2	2.38	0.43
1:D:97:GLN:H	1:D:98:PRO:HD3	1.81	0.43
1:C:331:ASP:OD1	1:C:331:ASP:C	2.62	0.43
1:C:11:ASP:C	1:C:11:ASP:OD1	2.62	0.43
1:C:31:LEU:HB3	1:C:372:ILE:HG22	2.01	0.43
1:D:431:GLU:O	1:D:434:GLN:HG2	2.19	0.43
1:A:11:ASP:OD1	1:A:11:ASP:C	2.61	0.43
1:A:95:ASN:OD1	1:A:95:ASN:C	2.62	0.43
1:A:290:THR:HG23	1:A:292:GLU:N	2.32	0.43
1:D:136:ASP:OD1	1:D:137:GLU:N	2.49	0.43
1:A:115:VAL:HG23	1:A:147:LEU:HD12	2.01	0.42
1:A:136:ASP:O	1:A:136:ASP:OD2	2.37	0.42
1:A:404:LEU:HB2	1:A:415:PHE:HE1	1.84	0.42
1:B:214:SER:OG	1:B:239:LEU:HA	2.19	0.42
1:C:342:LEU:HD23	1:C:370:ILE:HD13	2.00	0.42
1:C:347:ASN:ND2	1:D:65:GLY:HA3	2.35	0.42
1:D:241:GLY:HA2	1:D:284:GLN:HG2	2.01	0.42
1:A:161:VAL:HG12	1:A:163:MET:CE	2.49	0.42
1:A:201:VAL:O	1:A:201:VAL:CG1	2.67	0.42
1:A:374:THR:H	1:A:375:PRO:HD3	1.79	0.42
1:B:298:LYS:HE2	1:B:298:LYS:HB3	1.76	0.42
1:B:362:VAL:O	1:B:362:VAL:CG1	2.65	0.42
1:C:102:SER:O	1:C:106:VAL:HG23	2.18	0.42
1:C:163:MET:N	4:C:816:HOH:O	2.52	0.42
1:D:140:LYS:HA	1:D:140:LYS:HD2	1.94	0.42
1:D:253:ARG:HH11	1:D:253:ARG:CG	2.28	0.42
1:D:331:ASP:OD1	1:D:331:ASP:C	2.62	0.42
1:B:74:GLU:HG2	1:B:78:TYR:CZ	2.54	0.42
1:C:202:ASN:O	1:C:202:ASN:CG	2.62	0.42
1:D:187:ARG:HA	1:D:324:ASP:HB2	2.00	0.42
1:A:408:TYR:O	1:A:414:ASP:HB3	2.19	0.42
1:A:178:ILE:HD12	1:A:198:CYS:SG	2.59	0.42
1:A:237:LYS:NZ	2:A:701:PLP:O3	2.49	0.42
1:A:266:PHE:CD1	1:A:267:PRO:HA	2.54	0.42
1:A:162:ASP:OD2	1:A:165:SER:CB	2.67	0.42
1:B:236:HIS:O	1:B:237:LYS:HB2	2.18	0.42
1:D:122:MET:O	1:D:128:GLY:HA3	2.20	0.42
1:A:393:MET:HE2	1:A:393:MET:HB2	1.91	0.42
1:A:404:LEU:HB2	1:A:415:PHE:CE1	2.55	0.42
1:C:161:VAL:HG12	1:C:163:MET:HE2	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:183:THR:HG22	1:A:184:SER:HB3	2.02	0.41
1:A:314:ARG:HH22	1:A:388:GLU:CD	2.28	0.41
1:A:357:THR:HG23	1:A:361:ASP:HB2	2.01	0.41
1:B:241:GLY:HA2	1:B:284:GLN:HG2	2.02	0.41
1:D:214:SER:OG	1:D:239:LEU:HA	2.20	0.41
1:A:220:LEU:HD12	1:A:297:THR:HG22	2.02	0.41
1:D:107:GLN:HE22	1:D:270:GLN:HE21	1.67	0.41
1:D:357:THR:HG22	1:D:367:PRO:HB3	2.02	0.41
1:A:383:LYS:HE2	4:A:803:HOH:O	2.19	0.41
1:B:118:LYS:HE3	1:B:174:GLN:O	2.20	0.41
1:B:122:MET:O	1:B:128:GLY:HA3	2.19	0.41
1:B:356:ASN:ND2	4:B:823:HOH:O	2.53	0.41
1:C:56:GLU:HB3	1:C:266:PHE:CZ	2.56	0.41
1:D:195:ARG:HD2	1:D:227:TYR:O	2.20	0.41
1:A:86:ASN:HB3	1:C:406:GLN:OE1	2.20	0.41
1:A:163:MET:HA	1:A:163:MET:HE2	2.01	0.41
1:D:132:HIS:HE1	1:D:181:GLY:O	2.04	0.41
1:A:276:ASN:OD1	1:A:276:ASN:N	2.52	0.41
1:C:60:HIS:O	1:C:61:LYS:HG3	2.21	0.41
1:D:335:TYR:CD1	1:D:335:TYR:N	2.88	0.41
1:A:58:TYR:HB3	1:A:59:PRO:HD2	2.03	0.41
1:A:349:ILE:HG21	1:A:428:LEU:HB2	2.02	0.41
1:C:34:SER:HB2	1:C:237:LYS:HD3	2.03	0.41
1:C:183:THR:HG22	1:C:184:SER:HB3	2.02	0.41
1:D:120:MET:SD	1:D:169:LEU:HD23	2.61	0.41
1:D:258:ASP:OD1	1:D:258:ASP:O	2.37	0.41
1:B:405:GLN:HE21	1:B:410:LYS:HG3	1.84	0.41
1:C:59:PRO:O	1:C:60:HIS:HB2	2.21	0.41
1:C:266:PHE:CD1	1:C:267:PRO:HA	2.55	0.41
1:B:107:GLN:HE22	1:B:270:GLN:HE21	1.67	0.41
1:C:302:LEU:HA	1:C:302:LEU:HD23	1.78	0.41
1:D:161:VAL:HG12	1:D:163:MET:HE2	1.99	0.41
1:D:209:ILE:HG21	1:D:224:PRO:HG3	2.03	0.41
1:A:20:LYS:HD3	1:B:69:TYR:CE1	2.55	0.41
1:A:358:ILE:CD1	1:A:368:SER:HB2	2.42	0.41
1:B:181:GLY:C	1:B:182:TYR:CG	2.98	0.41
1:B:264:SER:O	1:B:268:SER:HB2	2.21	0.41
1:C:307:LEU:HD11	1:C:390:ILE:HG22	2.03	0.41
1:D:18:LEU:O	1:D:21:GLU:HB3	2.20	0.41
1:D:193:GLY:O	1:D:197:ILE:HD12	2.21	0.41
1:B:195:ARG:HD2	1:B:227:TYR:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:299:GLN:HB3	1:B:378:THR:HG23	2.03	0.41
1:B:352:ALA:O	1:B:353:LEU:HB3	2.20	0.41
1:A:137:GLU:HB3	1:A:138:LYS:H	1.48	0.40
1:B:237:LYS:HB3	1:B:238:ILE:H	1.67	0.40
1:C:241:GLY:CA	1:C:284:GLN:HG2	2.51	0.40
1:D:102:SER:N	3:D:501:PO4:O3	2.54	0.40
1:B:185:TYR:HA	1:B:186:PRO:HD3	1.83	0.40
1:C:176:LYS:O	1:C:176:LYS:HG3	2.21	0.40
1:D:98:PRO:HA	1:D:270:GLN:NE2	2.36	0.40
1:A:56:GLU:H	1:A:56:GLU:HG3	1.70	0.40
1:A:127:GLY:O	1:A:182:TYR:HB3	2.20	0.40
1:B:237:LYS:NZ	2:B:701:PLP:O3	2.54	0.40
1:D:264:SER:O	1:D:268:SER:HB2	2.21	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	446/480 (93%)	422 (95%)	23 (5%)	1 (0%)	43	73
1	B	443/480 (92%)	426 (96%)	17 (4%)	0	100	100
1	C	439/480 (92%)	415 (94%)	22 (5%)	2 (0%)	24	58
1	D	443/480 (92%)	428 (97%)	15 (3%)	0	100	100
All	All	1771/1920 (92%)	1691 (96%)	77 (4%)	3 (0%)	43	73

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	364	CYS
1	C	156	ASN

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Mol	Chain	Res	Type
1	A	374	THR

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	395/421 (94%)	367 (93%)	28 (7%)	13	41
1	B	392/421 (93%)	361 (92%)	31 (8%)	11	37
1	C	389/421 (92%)	363 (93%)	26 (7%)	15	44
1	D	392/421 (93%)	354 (90%)	38 (10%)	8	29
All	All	1568/1684 (93%)	1445 (92%)	123 (8%)	11	37

All (123) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	-5	HIS
1	A	3	ASN
1	A	27	GLU
1	A	53	LYS
1	A	54	TYR
1	A	88	SER
1	A	100	SER
1	A	136	ASP
1	A	137	GLU
1	A	151	LYS
1	A	212	ILE
1	A	286	LYS
1	A	290	THR
1	A	320	THR
1	A	337	ILE
1	A	340	SER
1	A	353	LEU
1	A	354	ASN
1	A	358	ILE
1	A	364	CYS

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Mol	Chain	Res	Type
1	A	365	VAL
1	A	371	ARG
1	A	372	ILE
1	A	388	GLU
1	A	400	LEU
1	A	413	VAL
1	A	426	ASP
1	A	438	ASN
1	B	36	ASN
1	B	53	LYS
1	B	55	SER
1	B	89	GLU
1	B	100	SER
1	B	120	MET
1	B	122	MET
1	B	131	THR
1	B	151	LYS
1	B	165	SER
1	B	172	SER
1	B	183	THR
1	B	189	ILE
1	B	234	THR
1	B	252	LYS
1	B	253	ARG
1	B	254	ASN
1	B	261	ILE
1	B	298	LYS
1	B	315	ASN
1	B	327	LEU
1	B	328	ILE
1	B	337	ILE
1	B	355	LYS
1	B	390	ILE
1	B	405	GLN
1	B	413	VAL
1	B	420	VAL
1	B	425	ILE
1	B	427	GLU
1	B	431	GLU
1	C	3	ASN
1	C	37	LEU
1	C	38	THR

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Mol	Chain	Res	Type
1	C	53	LYS
1	C	54	TYR
1	C	83	GLU
1	C	88	SER
1	C	100	SER
1	C	136	ASP
1	C	144	THR
1	C	151	LYS
1	C	212	ILE
1	C	240	ARG
1	C	286	LYS
1	C	290	THR
1	C	320	THR
1	C	337	ILE
1	C	340	SER
1	C	353	LEU
1	C	358	ILE
1	C	371	ARG
1	C	372	ILE
1	C	385	LYS
1	C	413	VAL
1	C	425	ILE
1	C	438	ASN
1	D	36	ASN
1	D	40	THR
1	D	53	LYS
1	D	55	SER
1	D	89	GLU
1	D	100	SER
1	D	120	MET
1	D	122	MET
1	D	131	THR
1	D	138	LYS
1	D	139	LYS
1	D	144	THR
1	D	151	LYS
1	D	165	SER
1	D	172	SER
1	D	174	GLN
1	D	183	THR
1	D	189	ILE
1	D	234	THR

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Mol	Chain	Res	Type
1	D	252	LYS
1	D	253	ARG
1	D	254	ASN
1	D	258	ASP
1	D	261	ILE
1	D	298	LYS
1	D	315	ASN
1	D	321	ASN
1	D	327	LEU
1	D	328	ILE
1	D	337	ILE
1	D	355	LYS
1	D	390	ILE
1	D	405	GLN
1	D	413	VAL
1	D	420	VAL
1	D	425	ILE
1	D	427	GLU
1	D	431	GLU

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	30	ASN
1	A	86	ASN
1	A	107	GLN
1	A	123	HIS
1	A	202	ASN
1	A	259	GLN
1	A	270	GLN
1	A	274	HIS
1	A	284	GLN
1	A	350	ASN
1	B	36	ASN
1	B	67	ASN
1	B	123	HIS
1	B	174	GLN
1	B	250	ASN
1	B	270	GLN
1	B	274	HIS
1	B	289	ASN
1	B	315	ASN

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Mol	Chain	Res	Type
1	B	356	ASN
1	B	405	GLN
1	C	107	GLN
1	C	202	ASN
1	C	270	GLN
1	C	275	ASN
1	C	284	GLN
1	D	36	ASN
1	D	67	ASN
1	D	123	HIS
1	D	132	HIS
1	D	174	GLN
1	D	250	ASN
1	D	270	GLN
1	D	274	HIS
1	D	289	ASN
1	D	405	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

4 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	PO4	D	501	-	4,4,4	0.67	0	6,6,6	1.09	0
2	PLP	C	701	1	15,15,16	0.78	0	21,22,23	1.54	5 (23%)
2	PLP	A	701	1	15,15,16	0.80	1 (6%)	21,22,23	1.91	8 (38%)
2	PLP	B	701	1	15,15,16	1.19	1 (6%)	21,22,23	1.83	7 (33%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	PLP	C	701	1	-	2/6/6/8	0/1/1/1
2	PLP	A	701	1	-	5/6/6/8	0/1/1/1
2	PLP	B	701	1	-	1/6/6/8	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	701	PLP	C3-C2	-3.93	1.36	1.41
2	A	701	PLP	C3-C2	-2.44	1.38	1.41

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	701	PLP	C4A-C4-C5	4.83	125.91	120.94
2	B	701	PLP	O4P-C5A-C5	4.14	117.12	109.36
2	A	701	PLP	C4A-C4-C3	-3.20	115.19	120.52
2	C	701	PLP	O2P-P-O4P	-3.03	98.76	106.67
2	B	701	PLP	C5A-C5-C4	2.95	128.46	122.64
2	B	701	PLP	C5A-C5-C6	-2.85	114.71	119.36
2	C	701	PLP	O3P-P-O2P	2.74	118.07	107.80
2	C	701	PLP	O4P-C5A-C5	2.71	114.43	109.36
2	B	701	PLP	O3P-P-O2P	2.57	117.45	107.80
2	A	701	PLP	O4P-C5A-C5	2.37	113.81	109.36
2	B	701	PLP	C4A-C4-C3	-2.33	116.63	120.52
2	A	701	PLP	O3-C3-C2	2.31	122.37	117.58
2	A	701	PLP	O2P-P-O4P	-2.29	100.71	106.67
2	A	701	PLP	O3P-P-O2P	2.20	116.05	107.80
2	C	701	PLP	C5-C6-N1	-2.19	120.27	123.83
2	B	701	PLP	C4A-C4-C5	2.17	123.18	120.94
2	B	701	PLP	C6-N1-C2	2.11	123.02	119.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	701	PLP	O4P-P-O1P	-2.09	100.78	106.44
2	C	701	PLP	C6-C5-C4	2.08	119.81	118.10
2	A	701	PLP	C5A-C5-C6	-2.06	116.00	119.36

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	701	PLP	C4-C5-C5A-O4P
2	A	701	PLP	C6-C5-C5A-O4P
2	A	701	PLP	C5A-O4P-P-O1P
2	A	701	PLP	C5A-O4P-P-O2P
2	A	701	PLP	C5A-O4P-P-O3P
2	C	701	PLP	C5A-O4P-P-O2P
2	C	701	PLP	C5A-O4P-P-O3P
2	B	701	PLP	C5A-O4P-P-O1P

There are no ring outliers.

4 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	501	PO4	2	0
2	C	701	PLP	3	0
2	A	701	PLP	1	0
2	B	701	PLP	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

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	448/480 (93%)	-0.50	0	100	100	24, 46, 88, 108	0
1	B	445/480 (92%)	-0.53	0	100	100	25, 42, 73, 112	0
1	C	441/480 (91%)	0.67	39 (8%)	15	10	31, 77, 119, 120	0
1	D	445/480 (92%)	1.75	152 (34%)	1	1	55, 116, 120, 120	0
All	All	1779/1920 (92%)	0.35	191 (10%)	11	7	24, 63, 120, 120	0

All (191) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	239	LEU	6.8
1	D	34	SER	5.2
1	D	29	ILE	5.2
1	D	428	LEU	5.2
1	D	33	ALA	5.1
1	D	238	ILE	4.9
1	D	31	LEU	4.8
1	D	213	SER	4.8
1	D	395	LEU	4.7
1	D	376	ALA	4.2
1	D	82	LEU	4.2
1	D	279	ALA	4.2
1	C	269	PHE	4.2
1	D	425	ILE	4.1
1	D	345	THR	4.1
1	C	276	ASN	4.1
1	D	285	LEU	4.0
1	D	375	PRO	4.0
1	D	132	HIS	4.0
1	C	73	ILE	4.0
1	D	325	ASN	4.0

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Mol	Chain	Res	Type	RSRZ
1	D	152	LEU	3.9
1	D	373	GLY	3.9
1	C	63	TYR	3.9
1	D	233	THR	3.8
1	D	206	PHE	3.8
1	D	191	TYR	3.8
1	D	326	HIS	3.8
1	D	342	LEU	3.8
1	D	165	SER	3.7
1	D	143	ILE	3.7
1	C	272	GLY	3.7
1	D	30	ASN	3.7
1	D	84	ALA	3.6
1	D	371	ARG	3.6
1	D	353	LEU	3.6
1	D	377	LEU	3.6
1	C	271	GLY	3.5
1	D	209	ILE	3.5
1	D	131	THR	3.5
1	D	346	CYS	3.5
1	D	241	GLY	3.5
1	C	47	GLY	3.4
1	D	306	ALA	3.4
1	D	348	ALA	3.4
1	D	329	VAL	3.4
1	D	235	THR	3.3
1	D	14	LEU	3.3
1	C	46	LEU	3.3
1	D	181	GLY	3.3
1	D	211	HIS	3.2
1	D	397	ALA	3.2
1	D	124	LEU	3.2
1	D	275	ASN	3.2
1	D	220	LEU	3.2
1	C	69	TYR	3.2
1	D	212	ILE	3.2
1	D	130	LEU	3.2
1	C	270	GLN	3.2
1	D	207	ALA	3.2
1	D	441	PHE	3.1
1	D	300	VAL	3.1
1	D	349	ILE	3.1

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Mol	Chain	Res	Type	RSRZ
1	D	372	ILE	3.1
1	D	169	LEU	3.0
1	D	217	ALA	3.0
1	D	96	VAL	3.0
1	D	398	ILE	3.0
1	D	417	LYS	3.0
1	D	76	LEU	3.0
1	D	210	SER	3.0
1	C	256	GLY	3.0
1	D	221	LEU	3.0
1	D	216	VAL	3.0
1	D	230	VAL	3.0
1	C	52	ASN	2.9
1	D	364	CYS	2.9
1	D	73	ILE	2.9
1	D	370	ILE	2.9
1	D	393	MET	2.9
1	C	275	ASN	2.9
1	D	161	VAL	2.9
1	D	282	ALA	2.9
1	D	261	ILE	2.9
1	D	328	ILE	2.9
1	C	70	VAL	2.8
1	D	17	LEU	2.8
1	D	153	TYR	2.8
1	C	419	LEU	2.8
1	C	81	ALA	2.8
1	C	65	GLY	2.8
1	D	197	ILE	2.7
1	D	416	LYS	2.7
1	D	424	LYS	2.7
1	D	335	TYR	2.7
1	D	374	THR	2.7
1	D	394	LEU	2.7
1	D	396	LYS	2.7
1	D	311	LEU	2.7
1	D	419	LEU	2.7
1	D	215	PHE	2.6
1	D	129	HIS	2.6
1	D	185	TYR	2.6
1	D	186	PRO	2.6
1	D	242	PRO	2.6

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Mol	Chain	Res	Type	RSRZ
1	D	36	ASN	2.6
1	D	46	LEU	2.6
1	D	337	ILE	2.6
1	D	154	LYS	2.6
1	D	412	LEU	2.6
1	C	6	PRO	2.5
1	D	64	TYR	2.5
1	D	399	LEU	2.5
1	D	163	MET	2.5
1	D	387	MET	2.5
1	D	183	THR	2.5
1	D	81	ALA	2.5
1	D	258	ASP	2.5
1	C	56	GLU	2.5
1	C	277	LYS	2.5
1	D	426	ASP	2.5
1	C	51	SER	2.5
1	D	423	PRO	2.5
1	D	179	ILE	2.5
1	D	267	PRO	2.4
1	D	413	VAL	2.4
1	D	234	THR	2.4
1	D	283	CYS	2.4
1	D	70	VAL	2.4
1	D	280	ALA	2.4
1	D	400	LEU	2.4
1	D	228	ALA	2.4
1	D	270	GLN	2.4
1	C	110	TYR	2.4
1	C	345	THR	2.4
1	D	257	ILE	2.4
1	D	330	VAL	2.4
1	D	134	PHE	2.4
1	D	301	LEU	2.4
1	D	302	LEU	2.4
1	C	7	LEU	2.3
1	C	99	LEU	2.3
1	D	112	LEU	2.3
1	C	114	GLY	2.3
1	D	45	CYS	2.3
1	C	268	SER	2.3
1	C	261	ILE	2.3

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Mol	Chain	Res	Type	RSRZ
1	C	42	VAL	2.3
1	D	104	ALA	2.3
1	C	97	GLN	2.3
1	D	214	SER	2.3
1	C	54	TYR	2.3
1	D	319	VAL	2.3
1	C	53	LYS	2.2
1	D	237	LYS	2.2
1	C	266	PHE	2.2
1	D	262	ASN	2.2
1	D	269	PHE	2.2
1	D	315	ASN	2.2
1	D	26	ILE	2.2
1	D	240	ARG	2.2
1	D	357	THR	2.2
1	D	415	PHE	2.2
1	D	219	ASN	2.2
1	C	96	VAL	2.2
1	D	101	GLY	2.2
1	D	389	PHE	2.2
1	D	32	ILE	2.1
1	C	274	HIS	2.1
1	D	231	VAL	2.1
1	D	338	THR	2.1
1	C	249	PHE	2.1
1	D	332	LEU	2.1
1	D	166	VAL	2.1
1	D	99	LEU	2.1
1	D	366	SER	2.1
1	D	382	CYS	2.1
1	C	94	VAL	2.1
1	D	433	VAL	2.1
1	D	194	PHE	2.1
1	D	442	ALA	2.1
1	D	281	VAL	2.1
1	D	208	ASP	2.1
1	D	391	ALA	2.1
1	C	273	PRO	2.0
1	D	37	LEU	2.0
1	D	225	PHE	2.0
1	D	278	ILE	2.0
1	C	108	ALA	2.0

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Mol	Chain	Res	Type	RSRZ
1	D	284	GLN	2.0
1	D	180	CYS	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	PO4	D	501	5/5	0.80	0.11	67,73,88,88	0
2	PLP	C	701	15/16	0.88	0.12	73,79,89,90	0
2	PLP	A	701	15/16	0.97	0.07	26,31,33,36	0
2	PLP	B	701	15/16	0.99	0.06	26,28,29,29	0

6.5 Other polymers [i](#)

There are no such residues in this entry.