



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 01:26 PM UTC

PDB ID : 2I22 / pdb_00002i22
Title : Crystal structure of Escherichia coli phosphoheptose isomerase in complex with reaction substrate sedoheptulose 7-phosphate
Authors : Blakely, K.; Zhang, K.; DeLeon, G.; Wright, G.; Junop, M.
Deposited on : 2006-08-15
Resolution : 2.80 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
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

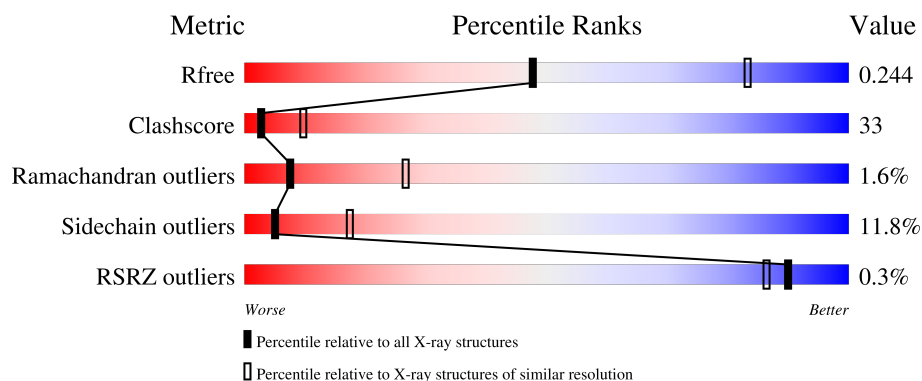
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	212	28% 37% 18% • 16%
1	B	212	29% 38% 16% 16%
1	C	212	32% 38% 12% • 16%
1	D	212	31% 35% 16% • 16%

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	I22	B	900	-	X	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5665 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 Phosphoheptose isomerase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	179	Total	C	N	O	S	0	0	0
			1365	853	245	260	7			
1	B	178	Total	C	N	O	S	0	0	0
			1357	849	244	257	7			
1	C	178	Total	C	N	O	S	0	0	0
			1357	849	244	257	7			
1	D	178	Total	C	N	O	S	0	0	0
			1357	849	244	257	7			

There are 80 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	expression tag	UNP P63224
A	-18	GLY	-	expression tag	UNP P63224
A	-17	SER	-	expression tag	UNP P63224
A	-16	SER	-	expression tag	UNP P63224
A	-15	HIS	-	expression tag	UNP P63224
A	-14	HIS	-	expression tag	UNP P63224
A	-13	HIS	-	expression tag	UNP P63224
A	-12	HIS	-	expression tag	UNP P63224
A	-11	HIS	-	expression tag	UNP P63224
A	-10	HIS	-	expression tag	UNP P63224
A	-9	SER	-	expression tag	UNP P63224
A	-8	SER	-	expression tag	UNP P63224
A	-7	GLY	-	expression tag	UNP P63224
A	-6	LEU	-	expression tag	UNP P63224
A	-5	VAL	-	expression tag	UNP P63224
A	-4	PRO	-	expression tag	UNP P63224
A	-3	ARG	-	expression tag	UNP P63224
A	-2	GLY	-	expression tag	UNP P63224
A	-1	SER	-	expression tag	UNP P63224
A	0	HIS	-	expression tag	UNP P63224
B	-19	MET	-	expression tag	UNP P63224

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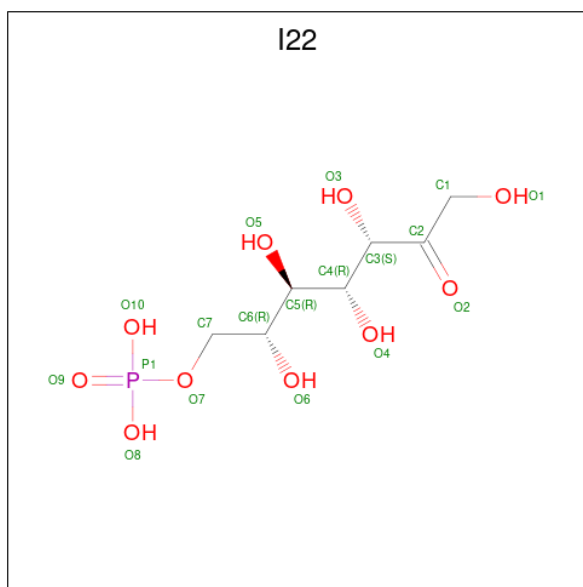
Chain	Residue	Modelled	Actual	Comment	Reference
B	-18	GLY	-	expression tag	UNP P63224
B	-17	SER	-	expression tag	UNP P63224
B	-16	SER	-	expression tag	UNP P63224
B	-15	HIS	-	expression tag	UNP P63224
B	-14	HIS	-	expression tag	UNP P63224
B	-13	HIS	-	expression tag	UNP P63224
B	-12	HIS	-	expression tag	UNP P63224
B	-11	HIS	-	expression tag	UNP P63224
B	-10	HIS	-	expression tag	UNP P63224
B	-9	SER	-	expression tag	UNP P63224
B	-8	SER	-	expression tag	UNP P63224
B	-7	GLY	-	expression tag	UNP P63224
B	-6	LEU	-	expression tag	UNP P63224
B	-5	VAL	-	expression tag	UNP P63224
B	-4	PRO	-	expression tag	UNP P63224
B	-3	ARG	-	expression tag	UNP P63224
B	-2	GLY	-	expression tag	UNP P63224
B	-1	SER	-	expression tag	UNP P63224
B	0	HIS	-	expression tag	UNP P63224
C	-19	MET	-	expression tag	UNP P63224
C	-18	GLY	-	expression tag	UNP P63224
C	-17	SER	-	expression tag	UNP P63224
C	-16	SER	-	expression tag	UNP P63224
C	-15	HIS	-	expression tag	UNP P63224
C	-14	HIS	-	expression tag	UNP P63224
C	-13	HIS	-	expression tag	UNP P63224
C	-12	HIS	-	expression tag	UNP P63224
C	-11	HIS	-	expression tag	UNP P63224
C	-10	HIS	-	expression tag	UNP P63224
C	-9	SER	-	expression tag	UNP P63224
C	-8	SER	-	expression tag	UNP P63224
C	-7	GLY	-	expression tag	UNP P63224
C	-6	LEU	-	expression tag	UNP P63224
C	-5	VAL	-	expression tag	UNP P63224
C	-4	PRO	-	expression tag	UNP P63224
C	-3	ARG	-	expression tag	UNP P63224
C	-2	GLY	-	expression tag	UNP P63224
C	-1	SER	-	expression tag	UNP P63224
C	0	HIS	-	expression tag	UNP P63224
D	-19	MET	-	expression tag	UNP P63224
D	-18	GLY	-	expression tag	UNP P63224
D	-17	SER	-	expression tag	UNP P63224

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-16	SER	-	expression tag	UNP P63224
D	-15	HIS	-	expression tag	UNP P63224
D	-14	HIS	-	expression tag	UNP P63224
D	-13	HIS	-	expression tag	UNP P63224
D	-12	HIS	-	expression tag	UNP P63224
D	-11	HIS	-	expression tag	UNP P63224
D	-10	HIS	-	expression tag	UNP P63224
D	-9	SER	-	expression tag	UNP P63224
D	-8	SER	-	expression tag	UNP P63224
D	-7	GLY	-	expression tag	UNP P63224
D	-6	LEU	-	expression tag	UNP P63224
D	-5	VAL	-	expression tag	UNP P63224
D	-4	PRO	-	expression tag	UNP P63224
D	-3	ARG	-	expression tag	UNP P63224
D	-2	GLY	-	expression tag	UNP P63224
D	-1	SER	-	expression tag	UNP P63224
D	0	HIS	-	expression tag	UNP P63224

- Molecule 2 is D-ALTRO-HEPT-2-ULOSE 7-PHOSPHATE (CCD ID: I22) (formula: $C_7H_{15}O_{10}P$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	B	1	Total	C	O	P	0	0
			18	7	10	1		

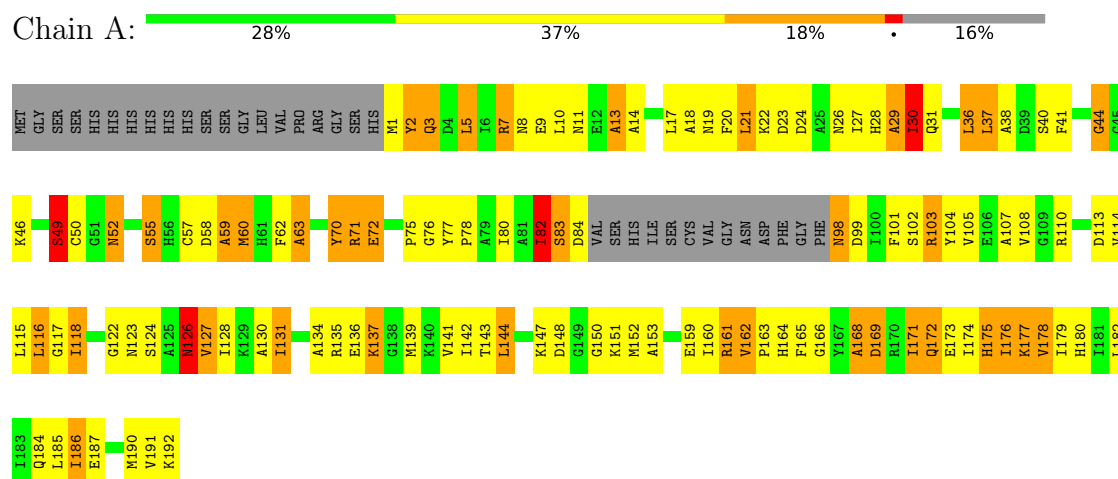
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	50	Total 50	O 50	0	0
3	B	52	Total 52	O 52	0	0
3	C	66	Total 66	O 66	0	0
3	D	43	Total 43	O 43	0	0

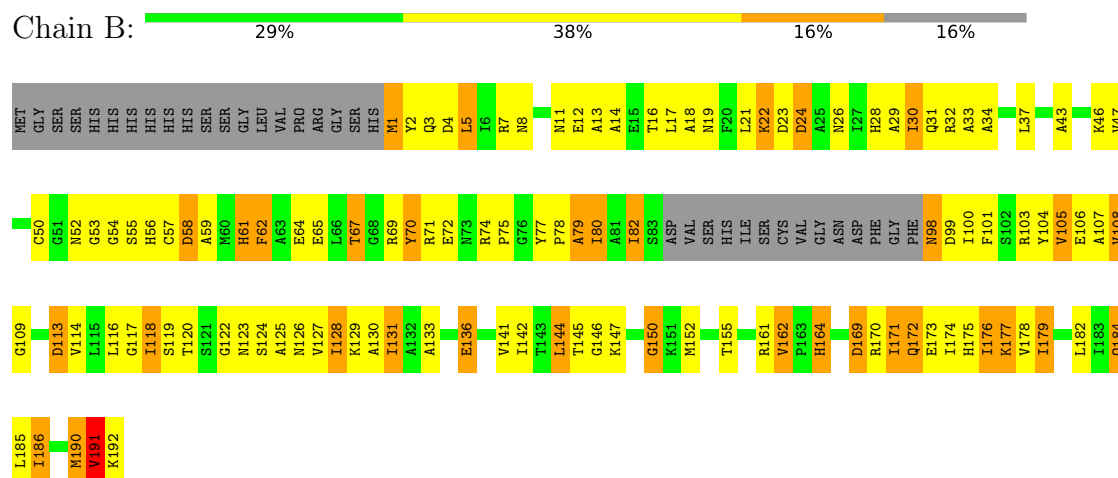
3 Residue-property plots

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

• Molecule 1: Phosphoheptose isomerase

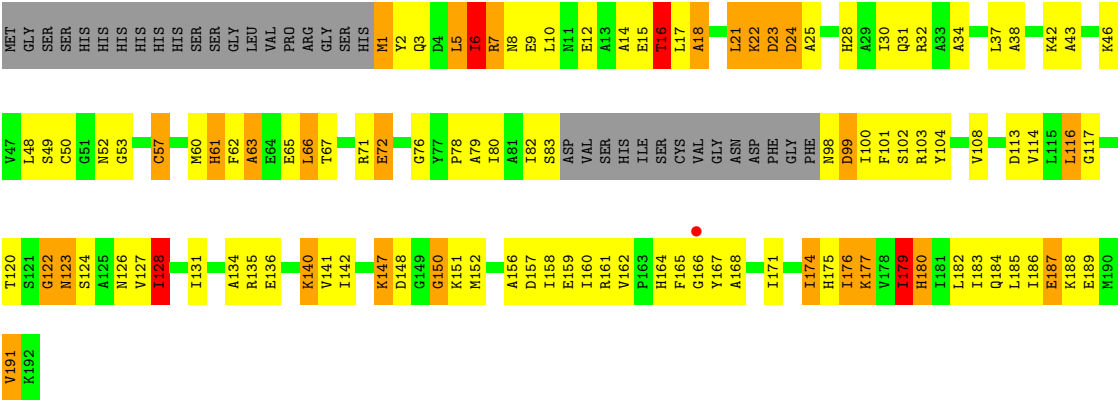


• Molecule 1: Phosphoheptose isomerase

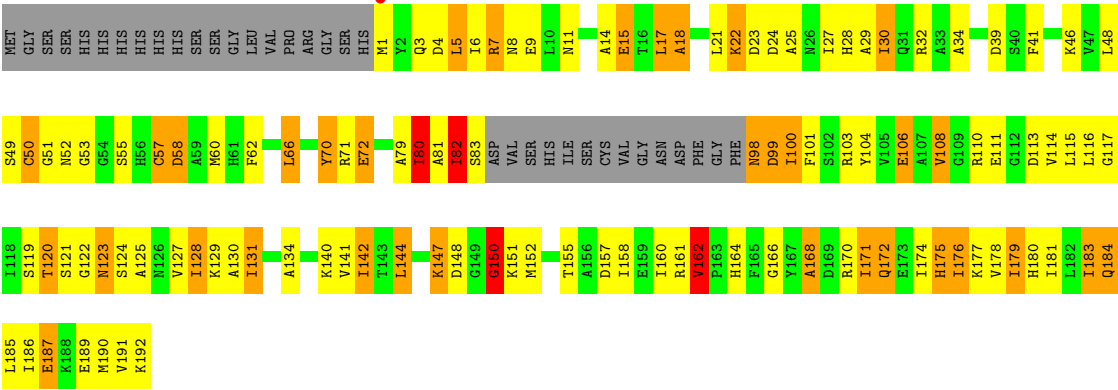
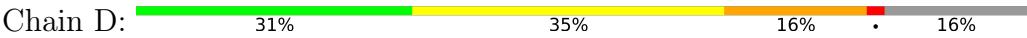


• Molecule 1: Phosphoheptose isomerase





● Molecule 1: Phosphoheptose isomerase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	73.05Å 76.54Å 78.32Å 90.00° 106.13° 90.00°	Depositor
Resolution (Å)	25.00 – 2.80 25.00 – 2.80	Depositor EDS
% Data completeness (in resolution range)	100.0 (25.00-2.80) 98.4 (25.00-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.71 (at 2.77Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.203 , 0.257 0.192 , 0.244	Depositor DCC
R_{free} test set	1499 reflections (7.30%)	wwPDB-VP
Wilson B-factor (Å ²)	54.0	Xtriage
Anisotropy	0.283	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 53.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5665	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.45% 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:
I22

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	2.18	42/1382 (3.0%)	1.80	16/1856 (0.9%)
1	B	2.27	55/1374 (4.0%)	1.78	20/1845 (1.1%)
1	C	2.27	50/1374 (3.6%)	1.84	28/1845 (1.5%)
1	D	2.19	41/1374 (3.0%)	1.83	20/1845 (1.1%)
All	All	2.23	188/5504 (3.4%)	1.81	84/7391 (1.1%)

All (188) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	63	ALA	CA-CB	-14.24	1.30	1.53
1	A	178	VAL	C-O	10.41	1.35	1.24
1	D	14	ALA	CA-CB	-10.24	1.37	1.53
1	C	49	SER	N-CA	-10.08	1.33	1.45
1	B	176	ILE	N-CA	-9.81	1.34	1.46
1	D	15	GLU	C-O	8.90	1.34	1.24
1	C	30	ILE	CA-CB	-8.70	1.44	1.54
1	A	117	GLY	N-CA	8.66	1.57	1.45
1	D	162	VAL	CA-C	-8.49	1.45	1.53
1	B	133	ALA	CA-CB	-8.43	1.40	1.53
1	A	176	ILE	N-CA	-8.42	1.36	1.46
1	A	82	ILE	CA-CB	-8.42	1.43	1.54
1	B	117	GLY	CA-C	-8.36	1.46	1.52
1	C	174	ILE	N-CA	-8.30	1.35	1.46
1	B	59	ALA	CA-CB	-8.27	1.39	1.53
1	A	191	VAL	CA-CB	8.20	1.65	1.54
1	C	127	VAL	CA-CB	-8.09	1.44	1.54
1	B	130	ALA	CA-CB	-8.01	1.41	1.53
1	B	30	ILE	CA-CB	-8.00	1.44	1.54
1	A	175	HIS	N-CA	7.97	1.56	1.46
1	C	117	GLY	CA-C	-7.84	1.46	1.52
1	B	74	ARG	CA-C	7.82	1.59	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	11	ASN	C-O	7.77	1.33	1.24
1	D	50	CYS	N-CA	7.62	1.55	1.46
1	C	136	GLU	C-O	-7.61	1.15	1.24
1	A	13	ALA	CA-CB	7.60	1.65	1.53
1	C	126	ASN	C-O	7.55	1.32	1.24
1	D	62	PHE	C-O	-7.53	1.15	1.24
1	A	134	ALA	CA-CB	-7.49	1.41	1.53
1	C	160	ILE	CA-CB	-7.33	1.45	1.53
1	B	105	VAL	CA-CB	-7.16	1.45	1.54
1	D	120	THR	CA-CB	-7.16	1.40	1.53
1	C	104	TYR	C-O	-7.14	1.15	1.24
1	D	101	PHE	C-O	-7.13	1.14	1.24
1	D	142	ILE	CG1-CD1	7.08	1.79	1.51
1	B	130	ALA	C-O	7.06	1.32	1.24
1	B	191	VAL	CA-CB	7.06	1.64	1.54
1	C	57	CYS	N-CA	-6.97	1.38	1.46
1	B	79	ALA	C-O	6.93	1.31	1.23
1	C	83	SER	C-O	6.91	1.37	1.23
1	D	147	LYS	N-CA	6.89	1.55	1.46
1	C	166	GLY	C-O	6.89	1.31	1.24
1	C	120	THR	CA-CB	-6.85	1.41	1.53
1	D	117	GLY	CA-C	-6.85	1.47	1.52
1	B	80	ILE	CA-CB	-6.82	1.45	1.54
1	C	101	PHE	C-O	-6.80	1.15	1.24
1	C	34	ALA	CA-CB	-6.79	1.42	1.53
1	D	125	ALA	CA-CB	-6.75	1.42	1.53
1	D	25	ALA	CA-C	6.74	1.61	1.52
1	A	171	ILE	CA-CB	-6.73	1.45	1.54
1	D	30	ILE	CA-CB	-6.73	1.47	1.54
1	A	75	PRO	N-CA	6.71	1.55	1.47
1	D	120	THR	N-CA	-6.69	1.37	1.46
1	C	134	ALA	CA-CB	-6.63	1.43	1.53
1	A	175	HIS	CA-C	6.61	1.62	1.52
1	B	144	LEU	N-CA	-6.59	1.38	1.46
1	D	18	ALA	C-O	6.59	1.32	1.24
1	A	59	ALA	CA-CB	-6.58	1.42	1.53
1	D	131	ILE	CA-C	-6.56	1.43	1.52
1	B	43	ALA	CA-CB	-6.54	1.43	1.53
1	B	75	PRO	CA-C	6.53	1.60	1.52
1	C	120	THR	N-CA	-6.51	1.37	1.46
1	D	9	GLU	C-O	6.50	1.31	1.24
1	D	100	ILE	C-O	-6.48	1.16	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	16	THR	CA-CB	-6.44	1.43	1.53
1	B	107	ALA	CA-CB	-6.40	1.43	1.53
1	B	190	MET	N-CA	-6.39	1.38	1.46
1	D	116	LEU	C-N	6.38	1.37	1.33
1	C	43	ALA	CA-CB	6.38	1.63	1.53
1	B	125	ALA	CA-CB	-6.35	1.43	1.53
1	B	128	ILE	N-CA	6.34	1.54	1.46
1	B	170	ARG	CA-C	-6.28	1.44	1.52
1	C	6	ILE	CG1-CD1	-6.28	1.27	1.51
1	A	174	ILE	CA-C	-6.26	1.44	1.52
1	B	178	VAL	CA-C	6.26	1.60	1.52
1	B	118	ILE	N-CA	-6.26	1.39	1.46
1	C	134	ALA	C-O	6.24	1.31	1.24
1	D	184	GLN	N-CA	6.20	1.53	1.46
1	C	14	ALA	CA-CB	-6.17	1.43	1.53
1	A	52	ASN	C-O	-6.17	1.15	1.23
1	C	140	LYS	CE-NZ	6.15	1.67	1.49
1	C	65	GLU	CA-CB	-6.14	1.43	1.53
1	A	127	VAL	CA-CB	-6.09	1.45	1.54
1	B	34	ALA	CA-CB	-6.05	1.43	1.53
1	A	168	ALA	CA-C	6.05	1.60	1.52
1	B	65	GLU	C-O	-6.04	1.17	1.24
1	B	64	GLU	N-CA	-6.01	1.39	1.46
1	B	178	VAL	C-O	6.00	1.31	1.24
1	B	142	ILE	C-O	5.99	1.30	1.24
1	C	48	LEU	C-O	-5.99	1.16	1.24
1	A	126	ASN	CA-C	-5.99	1.45	1.52
1	B	24	ASP	N-CA	5.98	1.53	1.46
1	A	128	ILE	CA-CB	5.98	1.62	1.54
1	A	80	ILE	C-O	-5.97	1.17	1.24
1	A	118	ILE	CA-CB	-5.96	1.47	1.54
1	D	125	ALA	C-O	5.95	1.30	1.24
1	D	121	SER	C-O	-5.94	1.16	1.24
1	C	160	ILE	C-O	-5.91	1.17	1.24
1	B	179	ILE	CA-CB	-5.91	1.47	1.54
1	A	60	MET	N-CA	5.90	1.53	1.46
1	A	14	ALA	N-CA	5.87	1.53	1.46
1	B	14	ALA	CA-CB	-5.86	1.44	1.53
1	C	114	VAL	N-CA	-5.84	1.39	1.46
1	D	180	HIS	N-CA	5.83	1.53	1.46
1	B	174	ILE	C-O	5.82	1.31	1.24
1	A	63	ALA	CA-CB	-5.80	1.44	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	141	VAL	CA-CB	-5.77	1.46	1.54
1	D	30	ILE	N-CA	-5.76	1.40	1.46
1	D	49	SER	N-CA	-5.76	1.38	1.45
1	B	18	ALA	CA-C	5.74	1.60	1.52
1	A	30	ILE	CA-CB	-5.73	1.46	1.54
1	A	14	ALA	CA-CB	-5.73	1.44	1.53
1	B	14	ALA	N-CA	5.73	1.53	1.46
1	D	130	ALA	N-CA	5.72	1.53	1.46
1	D	82	ILE	CA-CB	-5.71	1.48	1.54
1	D	57	CYS	N-CA	-5.70	1.39	1.46
1	B	185	LEU	C-O	5.69	1.30	1.24
1	C	176	ILE	N-CA	-5.69	1.40	1.46
1	B	70	TYR	N-CA	-5.64	1.39	1.46
1	B	113	ASP	C-O	-5.61	1.16	1.23
1	B	176	ILE	CA-CB	-5.60	1.47	1.54
1	B	171	ILE	CA-CB	-5.59	1.47	1.54
1	D	144	LEU	N-CA	-5.58	1.40	1.46
1	C	99	ASP	CA-C	5.56	1.60	1.52
1	D	186	ILE	CA-C	5.53	1.59	1.52
1	A	135	ARG	C-O	-5.52	1.17	1.24
1	C	180	HIS	N-CA	5.50	1.53	1.46
1	C	177	LYS	N-CA	5.49	1.53	1.46
1	C	128	ILE	CA-CB	-5.49	1.48	1.54
1	A	187	GLU	N-CA	-5.48	1.40	1.46
1	A	130	ALA	CA-CB	-5.46	1.44	1.53
1	C	43	ALA	N-CA	5.46	1.53	1.46
1	C	25	ALA	CA-C	5.46	1.59	1.52
1	B	116	LEU	C-N	5.45	1.37	1.33
1	C	6	ILE	C-O	5.44	1.30	1.24
1	C	165	PHE	CA-CB	-5.44	1.45	1.53
1	A	186	ILE	CA-C	5.44	1.59	1.52
1	A	134	ALA	CA-C	5.44	1.59	1.52
1	D	80	ILE	CA-CB	-5.42	1.46	1.54
1	A	49	SER	N-CA	-5.42	1.39	1.45
1	A	9	GLU	C-O	5.41	1.30	1.24
1	B	178	VAL	CA-CB	-5.40	1.47	1.54
1	C	174	ILE	C-O	5.39	1.30	1.23
1	D	119	SER	C-O	5.39	1.30	1.23
1	D	14	ALA	C-O	5.39	1.30	1.24
1	C	116	LEU	N-CA	-5.38	1.40	1.46
1	B	171	ILE	CA-C	5.37	1.59	1.52
1	A	176	ILE	CA-C	-5.37	1.46	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	80	ILE	CA-C	5.37	1.59	1.52
1	A	29	ALA	C-O	5.35	1.30	1.24
1	B	126	ASN	CA-C	-5.34	1.46	1.52
1	D	27	ILE	CA-CB	5.34	1.61	1.54
1	C	174	ILE	CA-CB	5.33	1.62	1.54
1	A	159	GLU	C-O	-5.31	1.17	1.23
1	B	82	ILE	C-O	-5.31	1.17	1.24
1	A	102	SER	CA-C	5.30	1.59	1.52
1	A	144	LEU	N-CA	-5.30	1.40	1.46
1	C	148	ASP	CA-C	-5.30	1.45	1.52
1	B	33	ALA	C-O	5.28	1.30	1.24
1	A	50	CYS	CA-C	-5.27	1.46	1.52
1	C	147	LYS	CE-NZ	5.26	1.65	1.49
1	D	58	ASP	CA-C	-5.26	1.45	1.52
1	C	165	PHE	CA-C	-5.26	1.46	1.52
1	B	13	ALA	C-O	5.26	1.30	1.24
1	C	18	ALA	C-O	5.25	1.30	1.24
1	B	164	HIS	C-O	5.24	1.30	1.23
1	B	179	ILE	N-CA	-5.24	1.40	1.46
1	C	116	LEU	C-N	5.24	1.36	1.33
1	C	23	ASP	C-O	-5.24	1.17	1.24
1	D	171	ILE	C-O	5.18	1.30	1.24
1	B	55	SER	CA-C	-5.16	1.45	1.52
1	B	61	HIS	N-CA	-5.16	1.40	1.46
1	B	114	VAL	C-O	-5.15	1.18	1.23
1	C	160	ILE	N-CA	-5.15	1.39	1.46
1	A	168	ALA	C-O	5.14	1.30	1.24
1	C	16	THR	N-CA	5.14	1.52	1.46
1	B	177	LYS	N-CA	5.13	1.52	1.46
1	A	18	ALA	N-CA	5.13	1.52	1.46
1	D	150	GLY	C-O	5.12	1.30	1.23
1	B	128	ILE	CA-CB	5.10	1.60	1.54
1	C	66	LEU	N-CA	-5.10	1.39	1.46
1	A	124	SER	N-CA	5.09	1.52	1.46
1	D	81	ALA	CA-CB	-5.09	1.46	1.53
1	B	136	GLU	CG-CD	5.07	1.64	1.52
1	D	62	PHE	CA-CB	-5.07	1.45	1.53
1	D	110	ARG	CA-C	-5.07	1.46	1.52
1	A	122	GLY	C-O	5.06	1.31	1.24
1	B	50	CYS	CA-C	-5.00	1.46	1.52

All (84) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	174	ILE	N-CA-C	-10.48	100.47	111.58
1	D	7	ARG	N-CA-C	-9.12	101.25	111.82
1	A	123	ASN	N-CA-C	8.92	125.25	113.30
1	D	123	ASN	N-CA-C	8.69	123.19	112.59
1	C	174	ILE	N-CA-C	-8.38	102.70	111.58
1	D	122	GLY	N-CA-C	-8.35	102.95	114.64
1	C	7	ARG	N-CA-C	-8.35	102.11	111.71
1	B	179	ILE	N-CA-C	-8.21	102.81	110.53
1	D	162	VAL	N-CA-C	-8.20	98.39	107.73
1	D	51	GLY	CA-C-O	-8.07	113.74	121.49
1	A	62	PHE	N-CA-C	-8.00	102.50	111.14
1	D	108	VAL	N-CA-C	7.98	121.35	113.53
1	C	67	THR	N-CA-C	-7.53	102.74	112.23
1	A	2	TYR	N-CA-C	-7.50	101.21	112.54
1	D	179	ILE	N-CA-C	-7.37	103.60	110.53
1	C	122	GLY	N-CA-C	-7.34	104.43	115.00
1	B	32	ARG	N-CA-C	-7.30	103.35	111.82
1	C	61	HIS	N-CA-C	7.20	120.04	111.33
1	C	61	HIS	N-CA-CB	7.17	120.81	110.06
1	B	186	ILE	N-CA-C	-7.13	103.35	110.62
1	C	17	LEU	N-CA-C	-7.07	102.77	111.40
1	D	48	LEU	CA-C-O	-7.05	112.70	120.38
1	A	122	GLY	N-CA-C	-7.04	105.60	115.32
1	B	190	MET	N-CA-C	-6.81	103.09	111.40
1	D	176	ILE	N-CA-C	6.69	117.45	110.62
1	A	37	LEU	N-CA-C	-6.64	103.73	110.97
1	B	164	HIS	N-CA-C	-6.62	98.97	109.23
1	C	60	MET	CG-SD-CE	6.57	115.35	100.90
1	C	160	ILE	N-CA-C	-6.53	97.58	106.85
1	A	161	ARG	NE-CZ-NH1	-6.44	115.06	121.50
1	D	66	LEU	N-CA-C	6.25	118.17	111.36
1	B	145	THR	CA-C-N	6.21	126.92	121.58
1	B	145	THR	C-N-CA	6.21	126.92	121.58
1	C	82	ILE	N-CA-C	6.11	116.31	108.12
1	A	177	LYS	CA-C-N	-6.10	112.87	120.56
1	A	177	LYS	C-N-CA	-6.10	112.87	120.56
1	C	78	PRO	N-CD-CG	-6.07	94.09	103.20
1	B	62	PHE	N-CA-C	-6.05	104.80	111.82
1	C	80	ILE	N-CA-C	6.03	116.56	108.11
1	A	161	ARG	NE-CZ-NH2	6.02	124.62	119.20
1	D	39	ASP	N-CA-C	-5.97	104.84	111.71
1	D	117	GLY	CA-C-O	-5.95	115.73	120.91
1	B	162	VAL	CB-CA-C	-5.89	104.09	111.04

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	58	ASP	N-CA-C	-5.87	104.96	111.71
1	C	16	THR	OG1-CB-CG2	-5.86	97.57	109.30
1	C	101	PHE	N-CA-C	-5.86	104.30	112.45
1	C	179	ILE	N-CA-C	-5.76	104.72	111.00
1	C	17	LEU	O-C-N	5.72	129.09	122.17
1	B	29	ALA	N-CA-C	-5.68	104.99	111.07
1	A	137	LYS	N-CA-C	-5.67	106.12	113.16
1	D	111	GLU	N-CA-C	-5.66	102.02	110.23
1	B	77	TYR	CA-C-N	-5.66	113.92	119.64
1	B	77	TYR	C-N-CA	-5.66	113.92	119.64
1	C	176	ILE	N-CA-C	5.61	115.81	110.42
1	B	136	GLU	N-CA-C	-5.58	105.34	111.82
1	C	24	ASP	N-CA-C	5.57	118.12	111.71
1	C	60	MET	CA-C-O	-5.57	114.59	120.55
1	B	119	SER	CB-CA-C	-5.53	102.59	110.34
1	A	102	SER	N-CA-C	5.47	117.05	111.14
1	B	61	HIS	N-CA-C	-5.45	105.34	111.28
1	A	70	TYR	N-CA-C	5.41	118.89	112.72
1	C	62	PHE	N-CA-C	-5.41	105.08	110.97
1	D	57	CYS	CA-CB-SG	-5.36	102.07	114.40
1	A	71	ARG	N-CA-C	-5.35	105.11	111.69
1	C	102	SER	N-CA-C	5.33	117.17	111.36
1	C	156	ALA	CA-C-O	-5.33	115.34	121.47
1	D	160	ILE	N-CA-C	-5.29	98.31	106.72
1	C	135	ARG	NE-CZ-NH1	-5.28	116.22	121.50
1	D	99	ASP	N-CA-C	-5.28	106.61	113.16
1	B	80	ILE	N-CA-C	5.28	115.72	108.12
1	C	82	ILE	CB-CA-C	-5.27	105.61	111.35
1	C	127	VAL	N-CA-C	5.26	115.99	110.62
1	D	141	VAL	CB-CA-C	-5.25	102.67	110.33
1	B	122	GLY	N-CA-C	-5.23	107.32	114.64
1	C	28	HIS	N-CA-C	-5.23	105.64	112.23
1	D	171	ILE	N-CA-C	-5.21	105.01	112.35
1	A	44	GLY	N-CA-C	5.20	123.12	113.76
1	A	126	ASN	N-CA-CB	5.17	117.51	110.01
1	D	106	GLU	CG-CD-OE2	-5.15	106.55	118.40
1	D	70	TYR	N-CA-C	5.15	119.25	112.92
1	C	123	ASN	N-CA-C	5.14	120.19	112.94
1	B	61	HIS	O-C-N	-5.10	116.71	122.12
1	C	72	GLU	N-CA-C	5.05	118.11	110.64
1	B	55	SER	N-CA-C	-5.03	107.15	113.28

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	1365	0	1375	122	0
1	B	1357	0	1371	91	0
1	C	1357	0	1371	76	0
1	D	1357	0	1371	99	0
2	B	18	0	13	14	0
3	A	50	0	0	12	0
3	B	52	0	0	10	0
3	C	66	0	0	4	0
3	D	43	0	0	1	0
All	All	5665	0	5501	362	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 33.

All (362) 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:142:ILE:CG1	1:D:142:ILE:CD1	1.79	1.52
1:C:140:LYS:CE	1:C:140:LYS:NZ	1.67	1.52
1:A:83:SER:HA	3:A:200:HOH:O	1.44	1.15
1:C:2:TYR:O	1:C:6:ILE:HD13	1.49	1.11
1:A:172:GLN:C	1:A:172:GLN:HE21	1.55	1.11
1:B:28:HIS:HB3	3:B:933:HOH:O	1.55	1.05
1:B:169:ASP:HB2	3:B:926:HOH:O	1.57	1.04
1:A:163:PRO:HD3	3:A:241:HOH:O	1.56	1.02
3:A:196:HOH:O	1:D:3:GLN:HG3	1.59	1.00
1:B:54:GLY:HA3	2:B:900:I22:O5	1.62	1.00
1:A:103:ARG:HG2	1:A:103:ARG:HH11	1.27	0.99
1:D:3:GLN:HG2	1:D:7:ARG:HH12	1.26	0.97
1:C:22:LYS:HG2	1:C:23:ASP:N	1.80	0.95
1:B:24:ASP:O	1:B:28:HIS:HD2	1.50	0.95
1:A:172:GLN:C	1:A:172:GLN:NE2	2.26	0.94

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:3:GLN:HG2	1:D:7:ARG:NH1	1.83	0.91
2:B:900:I22:O1	2:B:900:I22:H5	1.69	0.90
1:B:57:CYS:HG	1:C:57:CYS:HG	0.94	0.89
1:D:3:GLN:CG	1:D:7:ARG:HH12	1.86	0.88
1:A:31:GLN:HE22	1:D:1:MET:HA	1.38	0.87
2:B:900:I22:O1	2:B:900:I22:C5	2.22	0.87
1:A:127:VAL:O	1:A:131:ILE:CG1	2.25	0.85
1:B:70:TYR:CE2	1:B:190:MET:HG3	2.13	0.83
1:A:22:LYS:HG2	1:A:23:ASP:N	1.93	0.83
1:B:131:ILE:HG23	1:B:141:VAL:HG11	1.58	0.82
1:C:22:LYS:HG2	1:C:23:ASP:H	1.42	0.82
1:B:131:ILE:HD13	1:B:141:VAL:CG1	2.10	0.82
1:D:58:ASP:HB3	1:D:179:ILE:CD1	2.10	0.82
1:D:98:ASN:HD22	1:D:129:LYS:HZ3	1.25	0.82
1:C:2:TYR:O	1:C:6:ILE:CD1	2.28	0.82
1:A:28:HIS:HD2	3:A:196:HOH:O	1.62	0.81
1:A:172:GLN:O	1:A:175:HIS:HB2	1.81	0.81
1:D:98:ASN:HB3	1:D:129:LYS:HZ2	1.46	0.81
1:A:28:HIS:CD2	3:A:196:HOH:O	2.34	0.80
1:A:127:VAL:O	1:A:131:ILE:HG12	1.82	0.80
1:B:131:ILE:HD13	1:B:141:VAL:HG11	1.64	0.80
1:D:98:ASN:ND2	1:D:129:LYS:HZ3	1.80	0.80
1:D:172:GLN:HE21	1:D:172:GLN:C	1.90	0.79
1:B:56:HIS:CD2	3:B:938:HOH:O	2.33	0.79
1:C:6:ILE:N	1:C:6:ILE:HD12	1.96	0.79
1:B:162:VAL:HB	1:B:171:ILE:HG23	1.66	0.78
1:B:24:ASP:O	1:B:28:HIS:CD2	2.38	0.77
1:A:98:ASN:CB	3:A:220:HOH:O	2.33	0.76
1:A:177:LYS:HZ1	1:D:177:LYS:HZ1	1.33	0.76
1:C:21:LEU:HD12	1:C:21:LEU:O	1.85	0.76
2:B:900:I22:O5	2:B:900:I22:C1	2.34	0.76
1:C:6:ILE:HD12	1:C:6:ILE:H	1.49	0.75
1:D:147:LYS:O	1:D:161:ARG:HD3	1.85	0.75
1:B:46:LYS:O	1:B:113:ASP:HB3	1.86	0.75
1:B:144:LEU:HB3	1:B:175:HIS:CD2	2.21	0.75
1:B:177:LYS:NZ	1:C:177:LYS:HZ1	1.85	0.74
1:D:124:SER:O	1:D:128:ILE:HD12	1.86	0.74
1:A:98:ASN:HB3	3:A:220:HOH:O	1.86	0.74
1:B:177:LYS:HZ1	1:C:177:LYS:NZ	1.84	0.74
1:D:157:ASP:O	1:D:158:ILE:HG13	1.88	0.74
1:B:7:ARG:NH2	1:C:24:ASP:OD2	2.22	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:177:LYS:HZ1	1:C:177:LYS:HZ1	1.36	0.73
1:D:98:ASN:HB3	1:D:129:LYS:NZ	2.04	0.72
1:A:99:ASP:CB	3:A:204:HOH:O	2.37	0.72
1:A:127:VAL:O	1:A:131:ILE:HG13	1.87	0.72
1:C:22:LYS:CG	1:C:23:ASP:N	2.51	0.71
1:D:3:GLN:CD	1:D:7:ARG:HH12	1.98	0.71
1:D:127:VAL:O	1:D:131:ILE:HG13	1.91	0.70
1:D:22:LYS:HG2	1:D:23:ASP:N	2.06	0.70
2:B:900:I22:H3	1:C:180:HIS:NE2	2.06	0.70
1:D:98:ASN:CB	1:D:129:LYS:NZ	2.55	0.69
1:D:174:ILE:O	1:D:178:VAL:HG23	1.92	0.69
1:A:38:ALA:HA	1:A:190:MET:CE	2.23	0.69
1:B:2:TYR:HA	3:B:940:HOH:O	1.93	0.69
1:D:82:ILE:HG23	1:D:104:TYR:CD2	2.27	0.69
1:C:176:ILE:HA	1:C:179:ILE:HG13	1.74	0.69
1:B:103:ARG:HG2	1:B:103:ARG:HH11	1.56	0.69
1:C:6:ILE:CD1	1:C:6:ILE:H	2.06	0.69
1:A:71:ARG:HG3	1:A:72:GLU:N	2.07	0.68
1:A:177:LYS:NZ	1:D:177:LYS:NZ	2.42	0.68
1:A:31:GLN:NE2	1:D:1:MET:HA	2.08	0.68
1:B:182:LEU:O	1:B:186:ILE:HG13	1.93	0.67
1:B:26:ASN:O	1:B:30:ILE:HG13	1.95	0.67
1:A:57:CYS:O	1:A:58:ASP:C	2.34	0.67
1:A:22:LYS:HG2	1:A:23:ASP:H	1.59	0.67
1:D:98:ASN:C	1:D:129:LYS:HZ1	2.02	0.67
1:A:131:ILE:HG23	1:A:141:VAL:HG11	1.75	0.67
1:A:177:LYS:NZ	1:D:177:LYS:HZ1	1.92	0.67
1:A:103:ARG:HH11	1:A:103:ARG:CG	2.07	0.66
1:B:177:LYS:NZ	1:C:177:LYS:NZ	2.42	0.66
1:A:103:ARG:HG2	1:A:103:ARG:NH1	2.04	0.66
1:A:38:ALA:O	1:A:190:MET:HE1	1.96	0.66
1:B:8:ASN:ND2	3:B:913:HOH:O	2.30	0.65
1:C:175:HIS:O	1:C:179:ILE:HG12	1.97	0.64
1:C:15:GLU:O	1:C:18:ALA:HB3	1.97	0.64
1:C:122:GLY:O	1:C:152:MET:HG3	1.98	0.64
1:D:98:ASN:C	1:D:129:LYS:NZ	2.56	0.64
1:A:21:LEU:HD12	1:A:21:LEU:O	1.98	0.64
1:B:191:VAL:CG2	1:B:191:VAL:O	2.46	0.64
1:D:98:ASN:HD22	1:D:129:LYS:NZ	1.96	0.64
1:D:142:ILE:CD1	1:D:142:ILE:CB	2.72	0.64
1:D:28:HIS:O	1:D:32:ARG:HG3	1.97	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:191:VAL:O	1:B:191:VAL:HG23	1.98	0.63
1:B:54:GLY:HA3	2:B:900:I22:HO5	1.60	0.63
1:A:177:LYS:HZ1	1:D:177:LYS:NZ	1.97	0.63
1:A:175:HIS:O	1:A:179:ILE:HG13	1.99	0.62
1:D:98:ASN:ND2	1:D:129:LYS:NZ	2.46	0.62
1:A:37:LEU:O	1:A:40:SER:HB2	1.99	0.62
1:A:162:VAL:HG12	1:A:162:VAL:O	2.00	0.62
1:A:36:LEU:HD12	1:A:36:LEU:O	2.00	0.62
1:C:124:SER:O	1:C:128:ILE:HD12	1.99	0.62
1:D:191:VAL:HG23	1:D:191:VAL:O	1.97	0.61
1:C:191:VAL:HG23	1:C:191:VAL:O	1.98	0.61
1:A:17:LEU:HD12	1:A:17:LEU:O	2.01	0.61
1:C:21:LEU:HD12	1:C:21:LEU:C	2.25	0.61
1:B:184:GLN:NE2	3:B:909:HOH:O	2.34	0.61
1:A:5:LEU:O	1:A:8:ASN:HB3	2.01	0.60
1:D:147:LYS:HA	1:D:164:HIS:O	2.00	0.60
1:A:101:PHE:CD1	1:A:126:ASN:HB2	2.36	0.60
1:C:123:ASN:OD1	1:C:150:GLY:HA3	2.02	0.60
1:C:167:TYR:CD2	1:C:168:ALA:N	2.70	0.60
2:B:900:I22:O5	2:B:900:I22:H12	2.01	0.60
1:B:19:ASN:HA	1:B:22:LYS:HD2	1.84	0.60
1:D:144:LEU:O	1:D:175:HIS:CE1	2.55	0.60
1:D:98:ASN:N	3:D:211:HOH:O	2.34	0.59
1:A:98:ASN:HB2	3:A:220:HOH:O	1.98	0.59
1:A:22:LYS:NZ	3:A:212:HOH:O	2.34	0.59
1:B:37:LEU:HD13	1:B:62:PHE:HZ	1.65	0.59
1:A:172:GLN:NE2	1:A:173:GLU:N	2.51	0.59
1:A:22:LYS:CG	1:A:23:ASP:N	2.63	0.59
1:A:162:VAL:HB	1:A:171:ILE:HG23	1.85	0.58
1:B:104:TYR:O	1:B:105:VAL:C	2.45	0.58
1:A:23:ASP:C	1:A:23:ASP:OD1	2.47	0.57
1:A:164:HIS:CD2	1:A:166:GLY:H	2.21	0.57
1:A:147:LYS:HG2	1:A:165:PHE:HA	1.87	0.57
1:A:169:ASP:OD1	1:A:169:ASP:N	2.26	0.57
1:A:49:SER:OG	1:A:59:ALA:HB1	2.03	0.57
1:C:5:LEU:O	1:C:8:ASN:HB2	2.04	0.57
1:D:15:GLU:O	1:D:18:ALA:HB3	2.03	0.57
1:D:50:CYS:HB3	1:D:82:ILE:HG13	1.86	0.57
1:C:140:LYS:NZ	1:C:140:LYS:CD	2.62	0.57
1:C:159:GLU:OE1	1:C:161:ARG:NE	2.33	0.57
1:A:24:ASP:OD2	1:D:7:ARG:NH2	2.38	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:71:ARG:HG3	1:C:72:GLU:N	2.20	0.57
2:B:900:I22:C5	2:B:900:I22:C1	2.84	0.56
1:D:46:LYS:NZ	1:D:113:ASP:OD2	2.30	0.56
1:B:24:ASP:HB3	1:B:28:HIS:NE2	2.21	0.56
1:A:84:ASP:OD1	1:A:84:ASP:C	2.49	0.56
1:A:171:ILE:O	1:A:172:GLN:C	2.48	0.56
1:D:184:GLN:O	1:D:187:GLU:HB3	2.06	0.56
1:C:5:LEU:O	1:C:9:GLU:HG2	2.06	0.56
1:A:52:ASN:O	1:A:55:SER:HB2	2.06	0.55
1:B:162:VAL:CB	1:B:171:ILE:HG23	2.33	0.55
1:B:173:GLU:O	1:B:176:ILE:HG22	2.05	0.55
1:B:190:MET:C	1:B:192:LYS:H	2.13	0.55
1:B:7:ARG:O	1:B:11:ASN:HB2	2.06	0.55
1:C:16:THR:HG23	3:C:217:HOH:O	2.06	0.55
1:D:164:HIS:CE1	1:D:166:GLY:H	2.25	0.55
1:D:103:ARG:HA	1:D:106:GLU:OE1	2.07	0.55
1:A:103:ARG:CG	1:A:103:ARG:NH1	2.67	0.55
1:B:103:ARG:HG2	1:B:103:ARG:NH1	2.21	0.55
1:B:3:GLN:HB3	3:B:907:HOH:O	2.06	0.54
1:C:46:LYS:NZ	1:C:113:ASP:OD2	2.31	0.54
1:B:127:VAL:O	1:B:131:ILE:HG13	2.07	0.54
1:D:17:LEU:HD12	1:D:17:LEU:O	2.08	0.54
1:B:120:THR:O	1:B:146:GLY:HA3	2.08	0.53
1:B:54:GLY:HA3	2:B:900:I22:H12	1.89	0.53
1:D:150:GLY:C	1:D:152:MET:H	2.16	0.53
1:A:126:ASN:H	1:A:126:ASN:ND2	2.06	0.53
1:B:99:ASP:OD1	1:B:99:ASP:N	2.42	0.53
1:A:10:LEU:CD2	1:D:181:ILE:HD11	2.39	0.53
1:B:58:ASP:HB3	1:B:179:ILE:CD1	2.39	0.53
1:B:146:GLY:HA3	1:B:171:ILE:HD13	1.91	0.53
1:C:5:LEU:O	1:C:5:LEU:HD23	2.09	0.52
1:B:58:ASP:HB3	1:B:179:ILE:HD12	1.90	0.52
1:C:2:TYR:HA	3:C:202:HOH:O	2.09	0.52
1:A:10:LEU:O	1:A:13:ALA:HB3	2.09	0.52
1:A:57:CYS:O	1:A:60:MET:N	2.41	0.52
1:D:120:THR:HG21	1:D:168:ALA:HB1	1.91	0.52
1:A:10:LEU:HD21	1:D:181:ILE:CD1	2.38	0.52
1:D:21:LEU:O	1:D:21:LEU:HD12	2.09	0.52
1:D:148:ASP:N	1:D:148:ASP:OD1	2.39	0.52
1:B:4:ASP:O	1:B:8:ASN:HB2	2.10	0.52
1:C:3:GLN:HB3	3:C:210:HOH:O	2.08	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:10:LEU:HD23	1:D:181:ILE:HD11	1.91	0.52
1:A:52:ASN:ND2	1:A:126:ASN:OD1	2.34	0.52
1:B:147:LYS:O	1:B:161:ARG:HD3	2.10	0.52
1:C:5:LEU:C	1:C:5:LEU:CD2	2.83	0.51
1:C:103:ARG:HG2	1:C:103:ARG:HH11	1.76	0.51
1:C:76:GLY:HA2	1:D:103:ARG:O	2.11	0.51
1:C:162:VAL:HB	1:C:171:ILE:HG23	1.91	0.51
1:A:99:ASP:HB3	3:A:204:HOH:O	2.05	0.51
1:A:108:VAL:HG12	1:B:108:VAL:HB	1.92	0.51
1:A:162:VAL:CB	1:A:171:ILE:HG23	2.41	0.51
1:A:107:ALA:C	1:A:108:VAL:CG1	2.83	0.51
1:A:10:LEU:CD2	1:D:181:ILE:CD1	2.89	0.51
1:A:104:TYR:O	1:A:105:VAL:C	2.54	0.51
1:B:71:ARG:HG3	1:B:72:GLU:N	2.24	0.51
1:B:98:ASN:HD22	1:B:129:LYS:NZ	2.09	0.51
1:B:176:ILE:HG23	1:B:177:LYS:N	2.25	0.51
1:D:52:ASN:O	1:D:53:GLY:C	2.52	0.51
1:A:41:PHE:HD2	1:A:190:MET:HE3	1.75	0.51
1:B:69:ARG:HB2	1:B:70:TYR:CD1	2.45	0.51
1:D:17:LEU:HD22	1:D:177:LYS:HE2	1.93	0.51
1:D:123:ASN:OD1	1:D:151:LYS:HG3	2.11	0.51
1:A:172:GLN:O	1:A:175:HIS:N	2.44	0.50
1:A:1:MET:HA	1:A:3:GLN:HG2	1.94	0.50
1:A:168:ALA:O	1:A:171:ILE:N	2.44	0.50
1:D:58:ASP:HB3	1:D:179:ILE:HD13	1.92	0.50
1:C:32:ARG:HH11	1:C:32:ARG:HG3	1.77	0.50
1:D:98:ASN:O	1:D:129:LYS:NZ	2.45	0.50
1:D:152:MET:O	1:D:155:THR:OG1	2.27	0.49
1:A:177:LYS:HZ3	1:D:177:LYS:NZ	2.11	0.49
1:A:177:LYS:O	1:A:178:VAL:C	2.52	0.49
1:B:22:LYS:HG2	1:B:23:ASP:N	2.28	0.49
1:D:4:ASP:C	1:D:8:ASN:HD22	2.18	0.49
1:C:71:ARG:HG3	1:C:72:GLU:H	1.76	0.49
1:D:98:ASN:CB	1:D:129:LYS:HZ1	2.26	0.49
1:D:157:ASP:C	1:D:158:ILE:HG13	2.38	0.48
1:D:162:VAL:HB	1:D:171:ILE:HG23	1.95	0.48
1:A:41:PHE:O	1:A:44:GLY:N	2.44	0.48
1:A:182:LEU:O	1:A:186:ILE:HG13	2.14	0.48
1:A:107:ALA:C	1:A:108:VAL:HG13	2.37	0.48
1:B:5:LEU:HD13	1:C:185:LEU:CD2	2.44	0.48
1:A:110:ARG:HB2	1:A:113:ASP:OD2	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:168:ALA:O	1:A:169:ASP:C	2.56	0.48
1:B:144:LEU:HB3	1:B:175:HIS:NE2	2.29	0.48
1:C:5:LEU:HD23	1:C:5:LEU:C	2.39	0.48
1:D:98:ASN:CA	1:D:129:LYS:HZ1	2.25	0.48
1:B:61:HIS:O	1:B:62:PHE:C	2.49	0.48
1:D:5:LEU:O	1:D:8:ASN:HB2	2.14	0.48
1:B:54:GLY:N	1:C:61:HIS:CE1	2.82	0.48
1:B:67:THR:HA	1:B:78:PRO:HD2	1.96	0.48
1:D:99:ASP:OD2	1:D:103:ARG:NE	2.46	0.48
1:A:38:ALA:CA	1:A:190:MET:CE	2.92	0.47
1:A:184:GLN:O	1:A:185:LEU:C	2.54	0.47
1:D:71:ARG:HG3	1:D:72:GLU:N	2.29	0.47
1:A:41:PHE:CD2	1:A:190:MET:HE3	2.49	0.47
1:A:49:SER:O	1:A:82:ILE:HG13	2.14	0.47
1:B:24:ASP:OD2	1:C:7:ARG:NH2	2.48	0.47
2:B:900:I22:H3	1:C:180:HIS:CD2	2.49	0.46
1:C:12:GLU:O	1:C:16:THR:OG1	2.33	0.46
1:A:71:ARG:CG	1:A:72:GLU:N	2.76	0.46
1:D:172:GLN:O	1:D:175:HIS:HB2	2.15	0.46
1:A:28:HIS:O	1:A:31:GLN:N	2.48	0.46
1:A:46:LYS:N	1:A:113:ASP:OD1	2.36	0.46
1:D:98:ASN:O	1:D:129:LYS:CE	2.64	0.46
1:A:7:ARG:NH2	1:D:24:ASP:OD2	2.48	0.46
1:A:70:TYR:CE2	1:A:190:MET:HG3	2.50	0.46
1:B:147:LYS:HA	1:B:164:HIS:O	2.15	0.46
1:A:58:ASP:HB3	1:A:179:ILE:CD1	2.45	0.46
1:A:26:ASN:O	1:A:27:ILE:C	2.59	0.46
1:B:56:HIS:HD2	3:B:938:HOH:O	1.87	0.46
1:D:6:ILE:HG21	1:D:6:ILE:HD13	1.51	0.46
1:C:176:ILE:HG23	1:C:177:LYS:N	2.30	0.46
1:B:172:GLN:O	1:B:175:HIS:HB2	2.16	0.46
2:B:900:I22:C2	1:C:61:HIS:HD1	2.28	0.45
1:D:58:ASP:HB3	1:D:179:ILE:HD12	1.95	0.45
1:A:110:ARG:O	1:A:113:ASP:HB2	2.16	0.45
1:C:167:TYR:HD2	1:C:168:ALA:H	1.59	0.45
1:C:23:ASP:C	1:C:23:ASP:OD1	2.60	0.45
1:D:170:ARG:O	1:D:174:ILE:HG12	2.16	0.45
1:A:143:THR:O	1:A:144:LEU:HD23	2.17	0.45
1:A:114:VAL:HG12	1:A:115:LEU:N	2.31	0.45
1:B:21:LEU:O	1:B:21:LEU:HD12	2.16	0.45
1:B:103:ARG:HA	1:B:106:GLU:OE1	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:17:LEU:HD12	1:D:17:LEU:C	2.41	0.45
1:D:100:ILE:HG21	1:D:100:ILE:HD13	1.48	0.45
1:A:162:VAL:O	1:A:164:HIS:N	2.44	0.45
1:B:1:MET:N	3:B:906:HOH:O	2.50	0.45
1:B:2:TYR:OH	1:C:189:GLU:HG2	2.17	0.45
1:C:116:LEU:HA	1:C:142:ILE:O	2.17	0.45
1:C:32:ARG:HG3	1:C:32:ARG:NH1	2.31	0.45
1:A:29:ALA:HB1	1:A:160:ILE:HG12	1.99	0.44
1:C:10:LEU:HA	1:C:10:LEU:HD23	1.61	0.44
1:D:114:VAL:HG12	1:D:115:LEU:N	2.30	0.44
1:D:150:GLY:C	1:D:152:MET:N	2.75	0.44
1:C:99:ASP:N	1:C:99:ASP:OD1	2.48	0.44
1:A:152:MET:O	1:A:153:ALA:C	2.61	0.44
1:B:124:SER:O	1:B:128:ILE:HG13	2.17	0.44
1:B:47:VAL:HG12	1:B:78:PRO:O	2.17	0.44
1:A:21:LEU:O	1:A:27:ILE:HD11	2.17	0.44
1:B:23:ASP:OD1	1:B:23:ASP:C	2.60	0.44
1:B:100:ILE:HG23	1:B:101:PHE:CD2	2.53	0.44
1:C:131:ILE:HD11	1:C:152:MET:HE3	2.00	0.44
1:B:7:ARG:HG3	1:B:7:ARG:HH11	1.82	0.44
1:B:52:ASN:O	1:B:53:GLY:C	2.61	0.44
1:C:38:ALA:O	1:C:42:LYS:HG3	2.18	0.44
1:D:172:GLN:C	1:D:172:GLN:NE2	2.68	0.44
1:A:190:MET:C	1:A:192:LYS:H	2.25	0.43
1:D:131:ILE:O	1:D:134:ALA:HB3	2.17	0.43
1:A:57:CYS:SG	1:D:57:CYS:HA	2.58	0.43
1:A:60:MET:O	1:A:63:ALA:N	2.51	0.43
1:D:129:LYS:HB2	1:D:129:LYS:HE2	1.56	0.43
1:A:110:ARG:NH2	1:B:109:GLY:O	2.51	0.43
1:C:37:LEU:HD23	1:C:37:LEU:HA	1.76	0.43
1:C:50:CYS:HA	3:C:204:HOH:O	2.18	0.43
1:D:29:ALA:O	1:D:30:ILE:C	2.60	0.43
1:B:57:CYS:SG	1:C:57:CYS:HA	2.59	0.43
1:B:176:ILE:CG2	1:B:177:LYS:N	2.81	0.43
2:B:900:I22:O5	2:B:900:I22:O1	2.37	0.43
1:D:98:ASN:CG	1:D:129:LYS:NZ	2.77	0.43
1:A:127:VAL:O	1:A:127:VAL:HG12	2.18	0.43
1:A:20:PHE:HZ	1:A:30:ILE:HD11	1.83	0.43
1:A:177:LYS:NZ	1:D:177:LYS:HZ3	2.17	0.43
1:C:147:LYS:HA	1:C:164:HIS:O	2.18	0.43
1:A:58:ASP:HB3	1:A:179:ILE:HD11	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:66:LEU:HD21	1:D:183:ILE:CG1	2.48	0.42
1:D:70:TYR:CE2	1:D:190:MET:HG3	2.54	0.42
1:B:98:ASN:HD22	1:B:129:LYS:HZ2	1.66	0.42
2:B:900:I22:O5	2:B:900:I22:C2	2.67	0.42
1:C:157:ASP:C	1:C:158:ILE:HG13	2.44	0.42
1:A:77:TYR:HA	1:A:78:PRO:HD3	1.85	0.42
1:B:144:LEU:C	1:B:175:HIS:NE2	2.77	0.42
1:C:100:ILE:HD13	1:C:100:ILE:HG21	1.64	0.42
1:A:114:VAL:O	1:A:139:MET:CE	2.67	0.42
1:A:1:MET:HE2	1:A:1:MET:HB3	1.88	0.42
1:A:57:CYS:SG	1:D:60:MET:HE2	2.60	0.42
1:C:66:LEU:CD2	1:C:183:ILE:HG12	2.49	0.42
1:C:186:ILE:HD13	1:C:186:ILE:HG21	1.71	0.42
1:B:186:ILE:HD13	1:B:186:ILE:HG21	1.65	0.42
1:A:2:TYR:OH	1:D:189:GLU:HG2	2.19	0.42
1:A:164:HIS:HB3	1:A:171:ILE:HG12	2.01	0.42
1:D:41:PHE:N	1:D:41:PHE:CD1	2.86	0.42
1:D:82:ILE:HG23	1:D:104:TYR:CG	2.55	0.42
1:B:17:LEU:HD22	1:B:177:LYS:HE2	2.02	0.42
1:D:79:ALA:C	1:D:80:ILE:HG12	2.44	0.42
1:D:98:ASN:O	1:D:129:LYS:HE3	2.20	0.42
1:A:28:HIS:O	1:A:29:ALA:C	2.63	0.41
1:A:137:LYS:HA	1:A:137:LYS:HD3	1.85	0.41
1:B:28:HIS:CB	3:B:933:HOH:O	2.36	0.41
1:B:54:GLY:CA	2:B:900:I22:O5	2.50	0.41
1:A:179:ILE:O	1:A:180:HIS:C	2.61	0.41
1:C:52:ASN:O	1:C:53:GLY:C	2.63	0.41
1:C:184:GLN:O	1:C:187:GLU:HB3	2.20	0.41
1:D:34:ALA:CB	1:D:185:LEU:HB2	2.50	0.41
1:B:123:ASN:OD1	1:B:150:GLY:HA3	2.20	0.41
1:A:116:LEU:HA	1:A:142:ILE:O	2.20	0.41
1:A:118:ILE:HD13	1:A:118:ILE:HG21	1.75	0.41
1:B:79:ALA:C	1:B:80:ILE:HG13	2.44	0.41
1:A:165:PHE:N	3:A:199:HOH:O	2.54	0.41
1:A:29:ALA:CB	1:A:160:ILE:HG12	2.50	0.41
1:A:76:GLY:O	1:B:103:ARG:NH1	2.54	0.41
1:B:12:GLU:OE2	1:B:164:HIS:HE1	2.04	0.41
1:D:144:LEU:O	1:D:175:HIS:HE1	2.02	0.41
1:D:162:VAL:HG21	1:D:174:ILE:HB	2.02	0.41
1:A:19:ASN:HA	1:A:22:LYS:HD2	2.02	0.41
1:A:131:ILE:CG2	1:A:141:VAL:HG11	2.46	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:107:ALA:CB	1:A:108:VAL:HG13	2.51	0.41
1:B:69:ARG:HB2	1:B:70:TYR:CE1	2.56	0.41
1:B:118:ILE:HD13	1:B:118:ILE:HG21	1.89	0.41
1:B:152:MET:O	1:B:155:THR:OG1	2.35	0.41
1:A:99:ASP:OD1	1:A:99:ASP:N	2.41	0.40
1:D:66:LEU:HD21	1:D:183:ILE:HG13	2.02	0.40
1:D:147:LYS:CA	1:D:164:HIS:O	2.69	0.40
1:D:191:VAL:O	1:D:191:VAL:CG2	2.68	0.40
1:B:28:HIS:CD2	1:B:28:HIS:N	2.89	0.40
1:B:31:GLN:HE22	1:C:1:MET:HA	1.86	0.40
1:C:174:ILE:HD12	1:C:174:ILE:HG23	1.89	0.40
1:C:63:ALA:HA	1:C:79:ALA:HB1	2.03	0.40
1:C:103:ARG:HG2	1:C:103:ARG:NH1	2.36	0.40
1:C:182:LEU:O	1:C:186:ILE:HG13	2.21	0.40
1:A:148:ASP:O	1:A:161:ARG:NH1	2.55	0.40
1:B:24:ASP:C	1:B:28:HIS:CD2	2.99	0.40
1:C:66:LEU:HD21	1:C:183:ILE:HG12	2.04	0.40
1:C:99:ASP:OD2	1:C:103:ARG:NE	2.54	0.40
1:A:29:ALA:O	1:A:30:ILE:C	2.65	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	175/212 (82%)	146 (83%)	26 (15%)	3 (2%)	7	25
1	B	174/212 (82%)	154 (88%)	16 (9%)	4 (2%)	5	18
1	C	174/212 (82%)	161 (92%)	11 (6%)	2 (1%)	11	36
1	D	174/212 (82%)	157 (90%)	15 (9%)	2 (1%)	11	36
All	All	697/848 (82%)	618 (89%)	68 (10%)	11 (2%)	7	27

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	82	ILE
1	A	169	ASP
1	C	150	GLY
1	A	150	GLY
1	B	150	GLY
1	B	169	ASP
1	B	82	ILE
1	D	150	GLY
1	C	191	VAL
1	D	168	ALA
1	B	191	VAL

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	140/168 (83%)	118 (84%)	22 (16%)	2	9
1	B	139/168 (83%)	128 (92%)	11 (8%)	11	35
1	C	139/168 (83%)	125 (90%)	14 (10%)	7	23
1	D	139/168 (83%)	120 (86%)	19 (14%)	3	13
All	All	557/672 (83%)	491 (88%)	66 (12%)	5	17

All (66) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	GLN
1	A	5	LEU
1	A	7	ARG
1	A	11	ASN
1	A	21	LEU
1	A	30	ILE
1	A	36	LEU
1	A	49	SER
1	A	55	SER

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Mol	Chain	Res	Type
1	A	72	GLU
1	A	82	ILE
1	A	83	SER
1	A	98	ASN
1	A	103	ARG
1	A	116	LEU
1	A	126	ASN
1	A	131	ILE
1	A	136	GLU
1	A	151	LYS
1	A	162	VAL
1	A	172	GLN
1	A	176	ILE
1	B	1	MET
1	B	5	LEU
1	B	22	LYS
1	B	67	THR
1	B	98	ASN
1	B	108	VAL
1	B	131	ILE
1	B	136	GLU
1	B	172	GLN
1	B	184	GLN
1	B	191	VAL
1	C	1	MET
1	C	5	LEU
1	C	6	ILE
1	C	16	THR
1	C	21	LEU
1	C	22	LYS
1	C	31	GLN
1	C	98	ASN
1	C	108	VAL
1	C	128	ILE
1	C	151	LYS
1	C	179	ILE
1	C	187	GLU
1	C	188	LYS
1	D	5	LEU
1	D	17	LEU
1	D	22	LYS
1	D	55	SER

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Mol	Chain	Res	Type
1	D	72	GLU
1	D	80	ILE
1	D	82	ILE
1	D	83	SER
1	D	98	ASN
1	D	108	VAL
1	D	128	ILE
1	D	140	LYS
1	D	162	VAL
1	D	172	GLN
1	D	175	HIS
1	D	176	ILE
1	D	183	ILE
1	D	187	GLU
1	D	192	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	11	ASN
1	A	26	ASN
1	A	61	HIS
1	A	73	ASN
1	A	172	GLN
1	B	3	GLN
1	B	8	ASN
1	B	26	ASN
1	B	28	HIS
1	B	61	HIS
1	B	98	ASN
1	B	164	HIS
1	C	3	GLN
1	C	8	ASN
1	C	19	ASN
1	C	26	ASN
1	C	28	HIS
1	C	98	ASN
1	D	11	ASN
1	D	28	HIS
1	D	98	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

1 ligand is 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	I22	B	900	-	17,17,17	3.54	9 (52%)	19,24,24	3.29	10 (52%)

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	I22	B	900	-	-	10/24/24/24	-

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	900	I22	P1-O7	8.96	1.88	1.60
2	B	900	I22	C7-C6	5.13	1.58	1.51
2	B	900	I22	O1-C1	4.71	1.57	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	900	I22	O5-C5	4.61	1.54	1.43
2	B	900	I22	O3-C3	4.06	1.50	1.42
2	B	900	I22	O6-C6	3.40	1.50	1.43
2	B	900	I22	C4-C5	3.09	1.59	1.53
2	B	900	I22	O7-C7	2.79	1.55	1.44
2	B	900	I22	P1-O10	2.63	1.64	1.54

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	900	I22	C6-C5-C4	-6.18	102.99	112.48
2	B	900	I22	O4-C4-C5	-5.81	95.88	109.46
2	B	900	I22	O5-C5-C6	-5.78	95.80	108.93
2	B	900	I22	C5-C4-C3	-4.61	105.47	113.62
2	B	900	I22	O3-C3-C4	-4.19	101.28	110.38
2	B	900	I22	O6-C6-C7	-4.00	101.17	109.99
2	B	900	I22	O7-C7-C6	-3.99	98.70	109.36
2	B	900	I22	O6-C6-C5	3.81	118.16	109.25
2	B	900	I22	O5-C5-C4	-2.76	103.02	109.46
2	B	900	I22	O7-P1-O9	2.02	111.90	106.44

There are no chirality outliers.

All (10) torsion outliers are listed below:

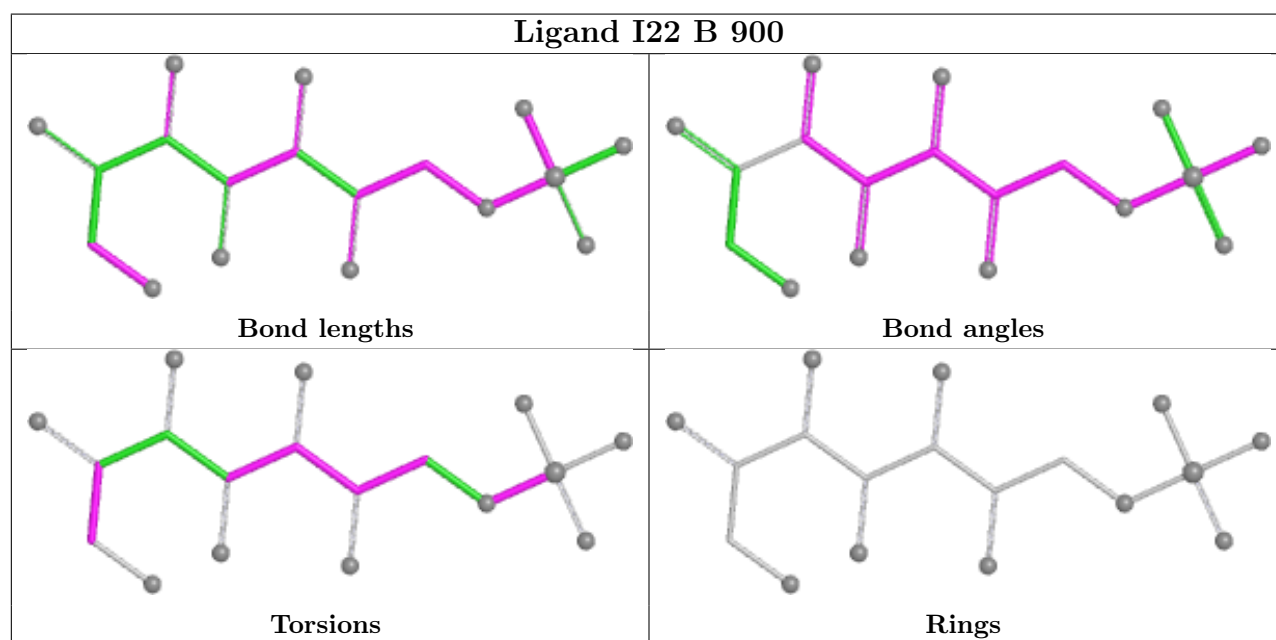
Mol	Chain	Res	Type	Atoms
2	B	900	I22	C7-O7-P1-O8
2	B	900	I22	C7-O7-P1-O10
2	B	900	I22	O1-C1-C2-C3
2	B	900	I22	O1-C1-C2-O2
2	B	900	I22	C4-C5-C6-C7
2	B	900	I22	C7-O7-P1-O9
2	B	900	I22	C4-C5-C6-O6
2	B	900	I22	O6-C6-C7-O7
2	B	900	I22	C3-C4-C5-O5
2	B	900	I22	O4-C4-C5-O5

There are no ring outliers.

1 monomer is involved in 14 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	900	I22	14	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	179/212 (84%)	-0.52	0 100 100	27, 43, 68, 84	0
1	B	178/212 (83%)	-0.57	0 100 100	22, 41, 68, 79	0
1	C	178/212 (83%)	-0.47	1 (0%) 85 80	22, 41, 67, 82	0
1	D	178/212 (83%)	-0.42	1 (0%) 85 80	22, 43, 69, 84	0
All	All	713/848 (84%)	-0.49	2 (0%) 90 86	22, 42, 68, 84	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	166	GLY	2.7
1	D	1	MET	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](#)

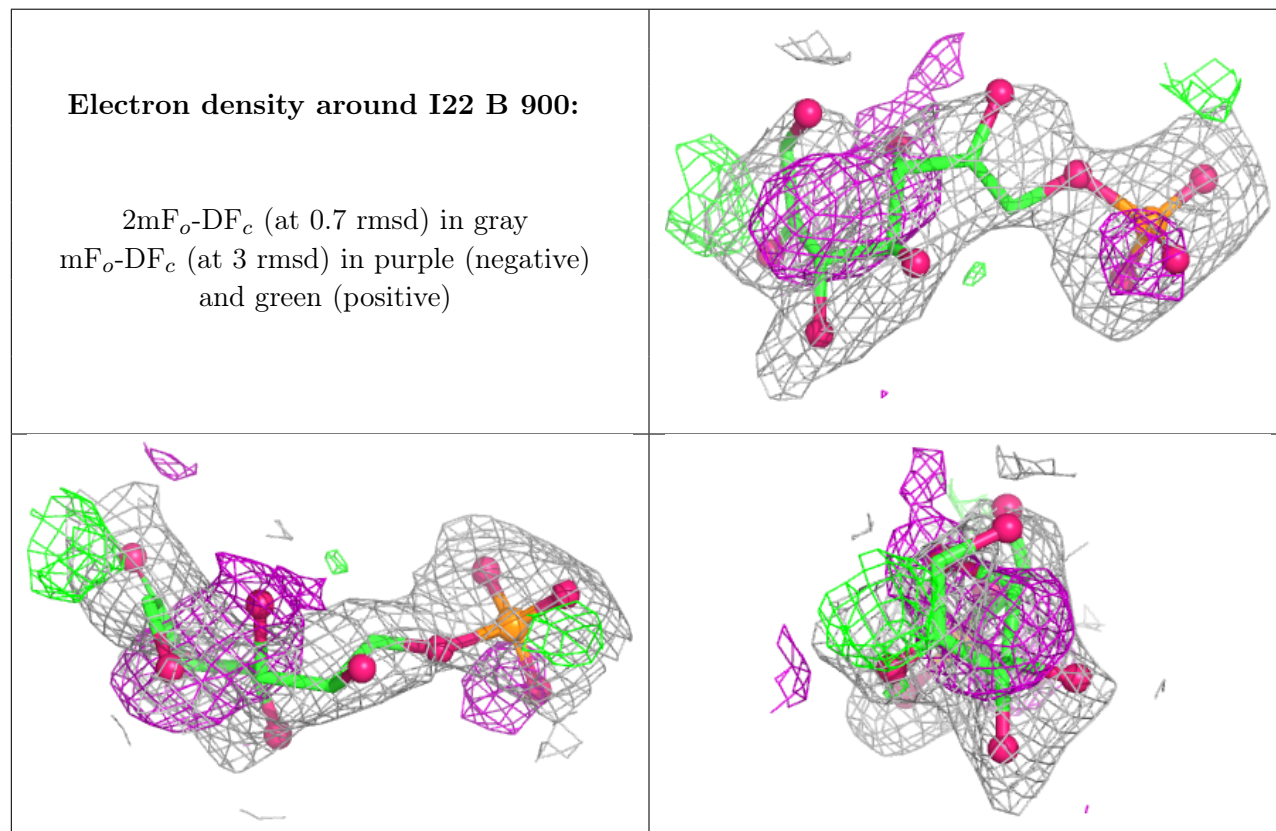
There are no oligosaccharides in this entry.

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($\text{\AA}^2$)	Q<0.9
2	I22	B	900	18/18	0.76	0.18	22,28,32,36	0

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



6.5 Other polymers ⓘ

There are no such residues in this entry.