



Full wwPDB X-ray Structure Validation Report ⓘ

May 26, 2020 – 04:55 pm BST

PDB ID : 2I22
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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

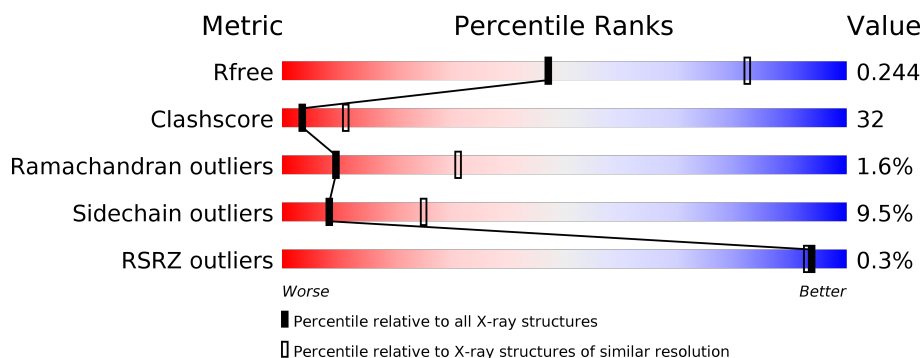
1 Overall quality at a glance

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}	130704	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (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 on 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	
1	B	212	
1	C	212	
1	D	212	

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 crite-

ria:

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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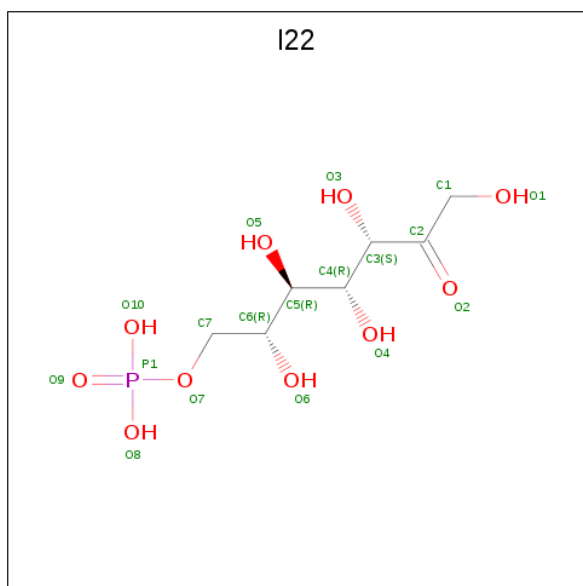
Chain	Residue	Modelled	Actual	Comment	Reference
B	-18	GLY	-	EXPRESSION TAG	UNP P63224
B	-17	SER	-	EXPRESSION TAG	UNP P63224
B	-16	SER	-	EXPRESSION TAG	UNP P63224
B	-15	HIS	-	EXPRESSION TAG	UNP P63224
B	-14	HIS	-	EXPRESSION TAG	UNP P63224
B	-13	HIS	-	EXPRESSION TAG	UNP P63224
B	-12	HIS	-	EXPRESSION TAG	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 (three-letter code: 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		

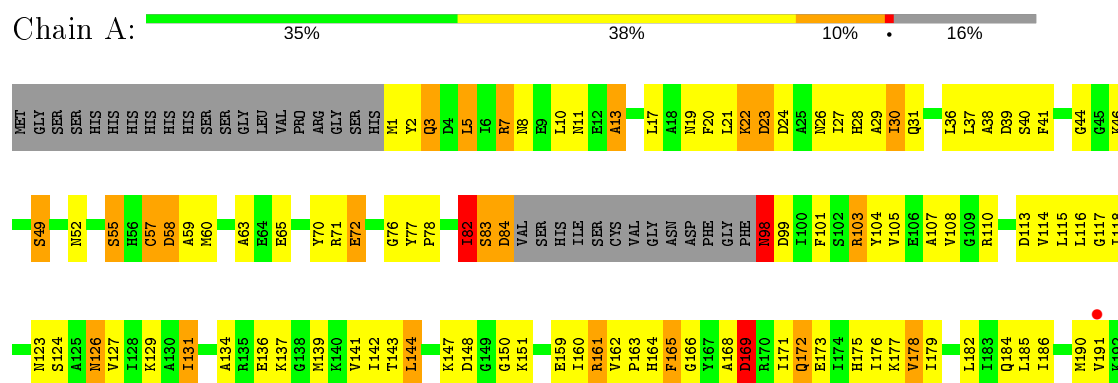
- Molecule 3 is water.

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

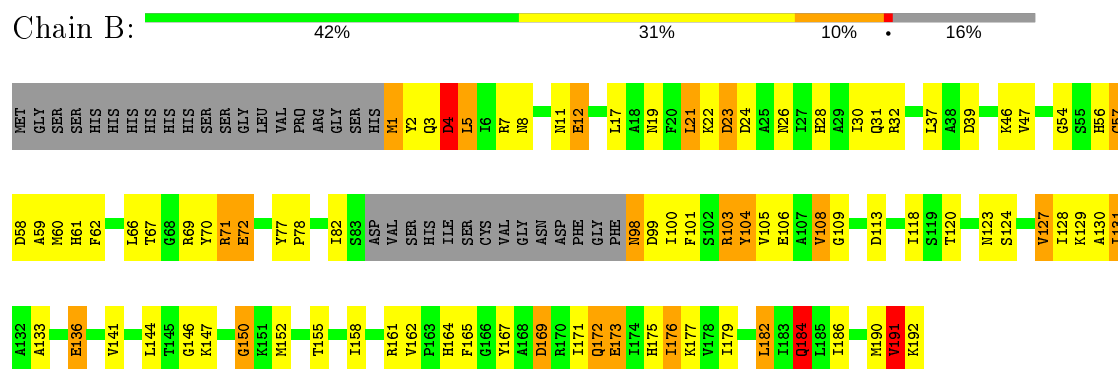
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($\text{RSRZ} > 2$). 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.

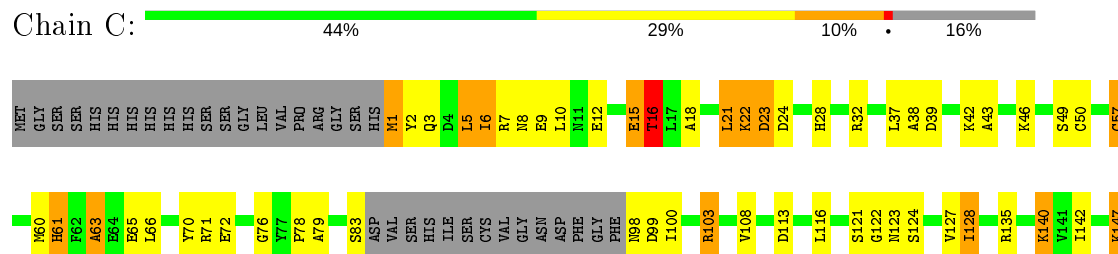
- Molecule 1: Phosphoheptose isomerase



- Molecule 1: Phosphoheptose isomerase

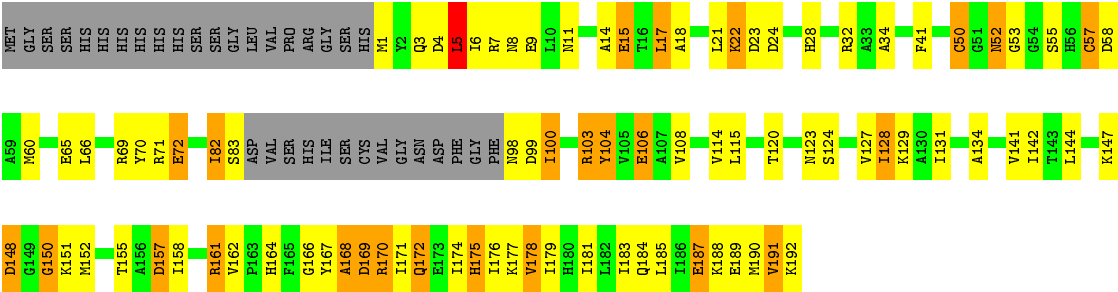


- Molecule 1: Phosphoheptose isomerase





● 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 47.39 – 2.79	Depositor EDS
% Data completeness (in resolution range)	100.0 (25.00-2.80) 98.5 (47.39-2.79)	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)	0.33 , 53.1	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 [i](#)

5.1 Standard geometry [i](#)

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	1.73	16/1382 (1.2%)	1.46	13/1856 (0.7%)
1	B	1.76	18/1374 (1.3%)	1.43	12/1845 (0.7%)
1	C	1.78	17/1374 (1.2%)	1.47	14/1845 (0.8%)
1	D	1.73	17/1374 (1.2%)	1.49	16/1845 (0.9%)
All	All	1.75	68/5504 (1.2%)	1.46	55/7391 (0.7%)

All (68) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	63	ALA	CA-CB	-10.38	1.30	1.52
1	B	136	GLU	CG-CD	8.52	1.64	1.51
1	A	117	GLY	N-CA	7.75	1.57	1.46
1	A	165	PHE	CE1-CZ	7.52	1.51	1.37
1	C	140	LYS	CE-NZ	7.50	1.67	1.49
1	A	65	GLU	CD-OE1	7.46	1.33	1.25
1	C	83	SER	C-O	7.38	1.37	1.23
1	D	14	ALA	CA-CB	-7.24	1.37	1.52
1	D	50	CYS	CB-SG	7.19	1.94	1.82
1	B	173	GLU	CD-OE1	7.12	1.33	1.25
1	C	159	GLU	CD-OE2	6.87	1.33	1.25
1	D	104	TYR	CD2-CE2	6.78	1.49	1.39
1	C	65	GLU	CD-OE1	6.78	1.33	1.25
1	A	178	VAL	C-O	6.60	1.35	1.23
1	B	165	PHE	CE1-CZ	6.51	1.49	1.37
1	C	49	SER	N-CA	-6.47	1.33	1.46
1	C	167	TYR	CG-CD2	6.46	1.47	1.39
1	C	147	LYS	CE-NZ	6.44	1.65	1.49
1	D	178	VAL	CB-CG2	6.33	1.66	1.52
1	D	141	VAL	CB-CG1	6.26	1.66	1.52
1	A	13	ALA	CA-CB	6.21	1.65	1.52
1	B	59	ALA	CA-CB	-6.14	1.39	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	136	GLU	CD-OE2	6.05	1.32	1.25
1	C	15	GLU	CB-CG	6.03	1.63	1.52
1	B	167	TYR	CD2-CE2	5.94	1.48	1.39
1	D	191	VAL	CB-CG1	5.92	1.65	1.52
1	B	133	ALA	CA-CB	-5.92	1.40	1.52
1	D	106	GLU	CD-OE2	-5.87	1.19	1.25
1	A	159	GLU	CD-OE2	5.85	1.32	1.25
1	B	77	TYR	CG-CD2	-5.79	1.31	1.39
1	D	15	GLU	C-O	5.78	1.34	1.23
1	D	167	TYR	CG-CD2	5.78	1.46	1.39
1	B	72	GLU	CG-CD	5.77	1.60	1.51
1	B	176	ILE	N-CA	-5.76	1.34	1.46
1	A	57	CYS	CB-SG	5.71	1.92	1.82
1	D	9	GLU	CB-CG	5.69	1.62	1.52
1	B	184	GLN	CG-CD	5.55	1.63	1.51
1	B	57	CYS	CB-SG	5.55	1.91	1.82
1	C	127	VAL	CB-CG1	5.53	1.64	1.52
1	B	130	ALA	CA-CB	-5.50	1.41	1.52
1	A	129	LYS	CE-NZ	5.46	1.62	1.49
1	B	127	VAL	CB-CG1	-5.45	1.41	1.52
1	C	43	ALA	CA-CB	5.44	1.63	1.52
1	C	174	ILE	N-CA	-5.43	1.35	1.46
1	D	106	GLU	CB-CG	-5.42	1.41	1.52
1	C	191	VAL	CB-CG1	5.42	1.64	1.52
1	C	159	GLU	CD-OE1	5.42	1.31	1.25
1	D	15	GLU	CG-CD	5.41	1.60	1.51
1	D	52	ASN	CB-CG	5.32	1.63	1.51
1	A	159	GLU	CD-OE1	5.24	1.31	1.25
1	D	187	GLU	CG-CD	5.21	1.59	1.51
1	D	167	TYR	CG-CD1	5.20	1.46	1.39
1	B	104	TYR	CD1-CE1	-5.20	1.31	1.39
1	B	136	GLU	CB-CG	5.19	1.62	1.52
1	B	167	TYR	CZ-OH	-5.18	1.29	1.37
1	A	191	VAL	CA-CB	5.17	1.65	1.54
1	A	82	ILE	CA-CB	-5.15	1.43	1.54
1	A	22	LYS	CB-CG	5.14	1.66	1.52
1	D	11	ASN	C-O	5.14	1.33	1.23
1	C	173	GLU	CD-OE2	5.13	1.31	1.25
1	A	134	ALA	CA-CB	-5.10	1.41	1.52
1	C	70	TYR	CD1-CE1	-5.07	1.31	1.39
1	B	12	GLU	CG-CD	5.05	1.59	1.51
1	A	124	SER	CB-OG	5.04	1.48	1.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	98	ASN	CB-CG	5.03	1.62	1.51
1	C	176	ILE	CB-CG2	-5.03	1.37	1.52
1	D	65	GLU	CD-OE2	5.02	1.31	1.25
1	A	165	PHE	CD2-CE2	5.01	1.49	1.39

All (55) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	103	ARG	NE-CZ-NH1	11.08	125.84	120.30
1	A	161	ARG	NE-CZ-NH1	-10.47	115.06	120.30
1	C	103	ARG	NE-CZ-NH1	9.83	125.22	120.30
1	C	60	MET	CG-SD-CE	9.47	115.35	100.20
1	C	39	ASP	CB-CG-OD1	9.26	126.64	118.30
1	A	58	ASP	CB-CG-OD1	9.26	126.64	118.30
1	A	169	ASP	CB-CG-OD1	-9.25	109.98	118.30
1	C	5	LEU	CB-CG-CD1	-8.86	95.94	111.00
1	D	169	ASP	CB-CG-OD1	-8.65	110.52	118.30
1	A	161	ARG	NE-CZ-NH2	8.64	124.62	120.30
1	D	69	ARG	NE-CZ-NH1	8.53	124.56	120.30
1	C	135	ARG	NE-CZ-NH1	-8.16	116.22	120.30
1	B	32	ARG	NE-CZ-NH1	8.15	124.38	120.30
1	D	60	MET	CG-SD-CE	6.92	111.27	100.20
1	B	182	LEU	CB-CG-CD2	-6.88	99.30	111.00
1	A	103	ARG	NE-CZ-NH2	-6.79	116.90	120.30
1	C	151	LYS	CD-CE-NZ	-6.70	96.29	111.70
1	D	57	CYS	CA-CB-SG	-6.63	102.07	114.00
1	D	188	LYS	CD-CE-NZ	-6.50	96.74	111.70
1	D	5	LEU	CB-CG-CD1	-6.46	100.03	111.00
1	C	169	ASP	CB-CG-OD1	-6.42	112.52	118.30
1	D	169	ASP	CB-CG-OD2	6.40	124.06	118.30
1	B	103	ARG	NE-CZ-NH1	6.40	123.50	120.30
1	B	21	LEU	CB-CG-CD2	-6.37	100.17	111.00
1	A	82	ILE	CG1-CB-CG2	-6.30	97.53	111.40
1	C	169	ASP	CB-CG-OD2	6.21	123.89	118.30
1	C	78	PRO	N-CD-CG	-6.07	94.09	103.20
1	A	23	ASP	CB-CG-OD2	-6.04	112.86	118.30
1	C	57	CYS	CA-CB-SG	-6.00	103.20	114.00
1	A	103	ARG	NE-CZ-NH1	5.95	123.28	120.30
1	B	158	ILE	CG1-CB-CG2	-5.93	98.35	111.40
1	B	4	ASP	CB-CG-OD2	-5.92	112.97	118.30
1	D	106	GLU	CG-CD-OE2	-5.88	106.55	118.30
1	B	39	ASP	CB-CG-OD2	-5.76	113.11	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	100	ILE	CG1-CB-CG2	-5.70	98.86	111.40
1	D	161	ARG	NE-CZ-NH1	5.69	123.14	120.30
1	C	61	HIS	N-CA-CB	5.67	120.81	110.60
1	A	39	ASP	CB-CG-OD2	-5.63	113.23	118.30
1	A	144	LEU	CB-CG-CD1	-5.63	101.43	111.00
1	B	103	ARG	NE-CZ-NH2	-5.60	117.50	120.30
1	B	66	LEU	CB-CG-CD1	-5.58	101.52	111.00
1	D	170	ARG	NE-CZ-NH1	-5.56	117.52	120.30
1	D	106	GLU	CG-CD-OE1	5.41	129.11	118.30
1	C	16	THR	OG1-CB-CG2	-5.40	97.57	110.00
1	D	50	CYS	CA-CB-SG	-5.39	104.30	114.00
1	B	191	VAL	CG1-CB-CG2	-5.36	102.33	110.90
1	A	123	ASN	N-CA-C	5.28	125.25	111.00
1	D	157	ASP	CB-CG-OD2	-5.24	113.58	118.30
1	B	23	ASP	CB-CG-OD1	5.20	122.98	118.30
1	D	148	ASP	CB-CG-OD1	-5.18	113.64	118.30
1	B	71	ARG	NE-CZ-NH1	5.17	122.89	120.30
1	A	169	ASP	CB-CG-OD2	5.16	122.95	118.30
1	C	159	GLU	OE1-CD-OE2	5.12	129.45	123.30
1	A	84	ASP	CB-CG-OD2	5.09	122.88	118.30
1	C	23	ASP	CB-CG-OD2	-5.08	113.73	118.30

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	116	0
1	B	1357	0	1371	88	0
1	C	1357	0	1371	75	0
1	D	1357	0	1371	93	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

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	D	43	0	0	1	0
All	All	5665	0	5501	346	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 32.

All (346) 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:CD1	1:D:142:ILE:CG1	1.79	1.52
1:C:140:LYS:NZ	1:C:140:LYS:CE	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	1.06
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: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:A:172:GLN:C	1:A:172:GLN:NE2	2.26	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:98:ASN:HB3	1:D:129:LYS:HZ2	1.46	0.81
1:A:127:VAL:O	1:A:131:ILE:HG12	1.82	0.80
1:A:28:HIS:CD2	3:A:196:HOH:O	2.34	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:B:56:HIS:CD2	3:B:938:HOH:O	2.33	0.79
1:C:6:ILE:HD12	1:C:6:ILE:N	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:O	1:C:21:LEU:HD12	1.85	0.76
2:B:900:I22:C1	2:B:900:I22:O5	2.34	0.76
1:C:6:ILE:H	1:C:6:ILE:HD12	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:D:172:GLN:C	1:D:172:GLN:HE21	1.90	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: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
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: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:A:71:ARG:HG3	1:A:72:GLU:N	2.07	0.69
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:HH11	1:B:103:ARG:HG2	1.56	0.69
1:C:6:ILE:H	1:C:6:ILE:CD1	2.06	0.69
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:177:LYS:NZ	1:D:177:LYS:HZ1	1.91	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:22:LYS:HG2	1:A:23:ASP:H	1.59	0.67
1:A:131:ILE:HG23	1:A:141:VAL:HG11	1.75	0.67
1:D:3:GLN:CD	1:D:7:ARG:HH12	1.98	0.67
1:B:177:LYS:NZ	1:C:177:LYS:NZ	2.42	0.67
1:A:103:ARG:CG	1:A:103:ARG:HH11	2.07	0.66
1:A:103:ARG:HG2	1:A:103:ARG:NH1	2.04	0.66
1:B:8:ASN:ND2	3:B:913:HOH:O	2.30	0.65
1:D:191:VAL:O	1:D:191:VAL:HG23	1.97	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:A:21:LEU:O	1:A:21:LEU:HD12	1.98	0.64
1:B:191:VAL:CG2	1:B:191:VAL:O	2.46	0.64
1:D:142:ILE:CD1	1:D:142:ILE:CB	2.72	0.64
1:D:98:ASN:HD22	1:D:129:LYS:NZ	1.96	0.64
1:D:28:HIS:O	1:D:32:ARG:HG3	1.97	0.64
1:B:191:VAL:HG23	1:B:191:VAL:O	1.98	0.63
1:B:54:GLY:HA3	2:B:900:I22:HO5	1.60	0.63
1:A:175:HIS:O	1:A:179:ILE:HG13	1.99	0.62
1:A:57:CYS:O	1:A:58:ASP:C	2.34	0.62
1:D:98:ASN:C	1:D:129:LYS:HZ1	2.02	0.62
1:D:98:ASN:ND2	1:D:129:LYS:NZ	2.46	0.62
1:A:162:VAL:HG12	1:A:162:VAL:O	2.00	0.62
1:A:37:LEU:O	1:A:40:SER:HB2	1.99	0.62
1:A:36:LEU:O	1:A:36:LEU:HD12	2.00	0.62
1:A:169:ASP:OD1	1:A:169:ASP:N	2.26	0.62
1:C:124:SER:O	1:C:128:ILE:HD12	1.99	0.62
1:A:177:LYS:HZ1	1:D:177:LYS:NZ	1.98	0.62
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: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:H12	2:B:900:I22:O5	2.01	0.60
1:B:19:ASN:HA	1:B:22:LYS:HD2	1.84	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:B:177:LYS:HZ1	1:C:177:LYS:HZ1	1.49	0.59
1:D:98:ASN:C	1:D:129:LYS:NZ	2.56	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: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: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:CD	1:C:140:LYS:NZ	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
1:C:71:ARG:HG3	1:C:72:GLU:N	2.20	0.57
1:C:21:LEU:C	1:C:21:LEU:HD12	2.25	0.56
2:B:900:I22:C1	2:B:900:I22:C5	2.84	0.56
1:B:24:ASP:HB3	1:B:28:HIS:NE2	2.21	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:177:LYS:HZ3	1:C:177:LYS:NZ	2.04	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:NH1	1:A:103:ARG:CG	2.67	0.55
1:B:103:ARG:NH1	1:B:103:ARG:HG2	2.21	0.55
1:B:177:LYS:HZ1	1:C:177:LYS:NZ	2.04	0.54
1:D:148:ASP:N	1:D:148:ASP:OD1	2.39	0.54
1:B:3:GLN:HB3	3:B:907:HOH:O	2.06	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:104:TYR:O	1:B:105:VAL:C	2.45	0.53
1:B:54:GLY:HA3	2:B:900:I22:H12	1.89	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:23:ASP:C	1:A:23:ASP:OD1	2.47	0.53
1:B:58:ASP:HB3	1:B:179:ILE:CD1	2.39	0.53
1:A:10:LEU:CD2	1:D:181:ILE:HD11	2.39	0.53
1:B:146:GLY:HA3	1:B:171:ILE:HD13	1.91	0.53
1:C:5:LEU:HD23	1:C:5:LEU:O	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: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
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:B:176:ILE:HG23	1:B:177:LYS:N	2.25	0.52
1:C:103:ARG:HG2	1:C:103:ARG:HH11	1.76	0.51
1:C:162:VAL:HB	1:C:171:ILE:HG23	1.91	0.51
1:C:76:GLY:HA2	1:D:103:ARG:O	2.11	0.51
1:A:162:VAL:CB	1:A:171:ILE:HG23	2.41	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:171:ILE:O	1:A:172:GLN:C	2.48	0.51
1:A:84:ASP:C	1:A:84:ASP:OD1	2.49	0.51
1:A:10:LEU:CD2	1:D:181:ILE:CD1	2.89	0.51
1:B:190:MET:C	1:B:192:LYS:H	2.13	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:A:177:LYS:HZ3	1:D:177:LYS:NZ	2.09	0.51
1:B:69:ARG:HB2	1:B:70:TYR:CD1	2.45	0.51
1:D:123:ASN:OD1	1:D:151:LYS:HG3	2.11	0.51
1:D:17:LEU:HD22	1:D:177:LYS:HE2	1.93	0.51
1:A:38:ALA:CA	1:A:190:MET:HE1	2.41	0.50
1:A:172:GLN:O	1:A:175:HIS:N	2.44	0.50
1:A:168:ALA:O	1:A:171:ILE:N	2.44	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1:MET:HA	1:A:3:GLN:HG2	1.94	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:A:38:ALA:HB1	1:A:190:MET:HE1	1.93	0.50
1:D:152:MET:O	1:D:155:THR:OG1	2.27	0.49
1:B:22:LYS:HG2	1:B:23:ASP:N	2.28	0.49
1:C:46:LYS:NZ	1:C:113:ASP:OD2	2.31	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:150:GLY:C	1:D:152:MET:H	2.16	0.48
1:D:162:VAL:HB	1:D:171:ILE:HG23	1.95	0.48
1:A:182:LEU:O	1:A:186:ILE:HG13	2.14	0.48
1:A:41:PHE:O	1:A:44:GLY:N	2.44	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
1:B:144:LEU:HB3	1:B:175:HIS:NE2	2.29	0.48
1:C:152:MET:O	1:C:155:THR:OG1	2.28	0.48
1:D:98:ASN:CA	1:D:129:LYS:HZ1	2.25	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:61:HIS:O	1:B:62:PHE:C	2.49	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:D:71:ARG:HG3	1:D:72:GLU:N	2.29	0.47
1:A:107:ALA:C	1:A:108:VAL:CG1	2.83	0.47
1:A:177:LYS:O	1:A:178:VAL:C	2.52	0.47
1:A:41:PHE:HB2	1:A:190:MET:HE3	1.97	0.47
1:A:49:SER:O	1:A:82:ILE:HG13	2.14	0.47
1:D:52:ASN:O	1:D:53:GLY:C	2.52	0.47
1:B:24:ASP:OD2	1:C:7:ARG:NH2	2.48	0.47
1:C:12:GLU:O	1:C:16:THR:OG1	2.33	0.46
2:B:900:I22:H3	1:C:180:HIS:CD2	2.49	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:46:LYS:N	1:A:113:ASP:OD1	2.36	0.46
1:A:28:HIS:O	1:A:31:GLN:N	2.48	0.46
1:D:98:ASN:O	1:D:129:LYS:CE	2.64	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:5:LEU:CD2	1:C:5:LEU:C	2.83	0.46
1:A:7:ARG:NH2	1:D:24:ASP:OD2	2.48	0.46
1:A:104:TYR:O	1:A:105:VAL:C	2.54	0.46
1:A:58:ASP:HB3	1:A:179:ILE:CD1	2.45	0.46
1:B:56:HIS:HD2	3:B:938:HOH:O	1.87	0.46
1:D:6:ILE:HD13	1:D:6:ILE:HG21	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: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:103:ARG:HA	1:B:106:GLU:OE1	2.17	0.45
1:B:21:LEU:HD12	1:B:21:LEU:O	2.16	0.45
1:D:100:ILE:HD13	1:D:100:ILE:HG21	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:NH1	1:C:32:ARG:HG3	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:C:99:ASP:N	1:C:99:ASP:OD1	2.48	0.44
1:B:124:SER:O	1:B:128:ILE:HG13	2.17	0.44
1:D:4:ASP:C	1:D:8:ASN:HD22	2.18	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:100:ILE:HG23	1:B:101:PHE:CD2	2.53	0.44
1:A:184:GLN:O	1:A:185:LEU:C	2.54	0.44
1:A:38:ALA:CB	1:A:190:MET:HE1	2.48	0.44
1:B:7:ARG:HG3	1:B:7:ARG:HH11	1.82	0.44
1:C:38:ALA:O	1:C:42:LYS:HG3	2.18	0.44
1:D:157:ASP:C	1:D:158:ILE:HG13	2.38	0.44
1:A:107:ALA:C	1:A:108:VAL:HG13	2.37	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:129:LYS:HB2	1:D:129:LYS:HE2	1.56	0.43
1:A:168:ALA:O	1:A:169:ASP:C	2.56	0.43
1:A:110:ARG:NH2	1:B:109:GLY:O	2.51	0.43
1:C:37:LEU:HA	1:C:37:LEU:HD23	1.76	0.43
1:C:50:CYS:HA	3:C:204:HOH:O	2.18	0.43
1:B:176:ILE:CG2	1:B:177:LYS:N	2.81	0.43
1:B:57:CYS:SG	1:C:57:CYS:HA	2.59	0.43
2:B:900:I22:O1	2:B:900:I22:O5	2.37	0.43
1:A:127:VAL:HG12	1:A:127:VAL:O	2.18	0.43
1:C:5:LEU:HD23	1:C:5:LEU:C	2.39	0.43
1:A:20:PHE:HZ	1:A:30:ILE:HD11	1.83	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
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
2:B:900:I22:O5	2:B:900:I22:C2	2.67	0.42
1:B:98:ASN:HD22	1:B:129:LYS:HZ2	1.66	0.42
1:A:77:TYR:HA	1:A:78:PRO:HD3	1.85	0.42
1:C:100:ILE:HG21	1:C:100:ILE:HD13	1.64	0.42
1:A:114:VAL:O	1:A:139:MET:CE	2.67	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:164:HIS:HB3	1:A:171:ILE:HG12	2.01	0.42
1:D:82:ILE:HG23	1:D:104:TYR:CG	2.55	0.42
1:A:2:TYR:OH	1:D:189:GLU:HG2	2.19	0.42
1:D:41:PHE:N	1:D:41:PHE:CD1	2.86	0.42
1:B:17:LEU:HD22	1:B:177:LYS:HE2	2.02	0.42
1:D:98:ASN:O	1:D:129:LYS:HE3	2.20	0.42
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:C:184:GLN:O	1:C:187:GLU:HB3	2.20	0.41
1:A:177:LYS:NZ	1:D:177:LYS:HZ3	2.18	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:118:ILE:HD13	1:A:118:ILE:HG21	1.75	0.41
1:A:116:LEU:HA	1:A:142:ILE:O	2.20	0.41
1:A:26:ASN:O	1:A:27:ILE:C	2.59	0.41
1:D:169:ASP:N	1:D:169:ASP:OD1	2.54	0.41
1:A:165:PHE:N	3:A:199:HOH:O	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:131:ILE:CG2	1:A:141:VAL:HG11	2.46	0.41
1:A:19:ASN:HA	1:A:22:LYS:HD2	2.02	0.41
1:A:107:ALA:CB	1:A:108:VAL:HG13	2.51	0.41
1:B:118:ILE:HG21	1:B:118:ILE:HD13	1.89	0.41
1:B:152:MET:O	1:B:155:THR:OG1	2.35	0.41
1:B:69:ARG:HB2	1:B:70:TYR:CE1	2.56	0.41
1:A:99:ASP:N	1:A:99:ASP:OD1	2.41	0.40
1:B:60:MET:HB3	1:B:60:MET:HE3	1.92	0.40
1:C:23:ASP:OD1	1:C:23:ASP:C	2.60	0.40
1:D:147:LYS:CA	1:D:164:HIS:O	2.69	0.40
1:D:66:LEU:HD21	1:D:183:ILE:HG13	2.02	0.40
1:D:191:VAL:CG2	1:D:191:VAL:O	2.68	0.40
1:B:28:HIS:CD2	1:B:28:HIS:N	2.89	0.40
1:C:174:ILE:HD12	1:C:174:ILE:HG23	1.89	0.40
1:B:31:GLN:HE22	1:C:1:MET:HA	1.86	0.40
1:C:103:ARG:NH1	1:C:103:ARG:HG2	2.36	0.40
1:C:182:LEU:O	1:C:186:ILE:HG13	2.21	0.40
1:C:63:ALA:HA	1:C:79:ALA:HB1	2.03	0.40
1:A:148:ASP:O	1:A:161:ARG:NH1	2.55	0.40
1:C:99:ASP:OD2	1:C:103:ARG:NE	2.54	0.40
1:C:66:LEU:HD21	1:C:183:ILE:HG12	2.04	0.40
1:D:17:LEU:C	1:D:17:LEU:HD12	2.41	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%)	9	29
1	B	174/212 (82%)	154 (88%)	16 (9%)	4 (2%)	6	21
1	C	174/212 (82%)	161 (92%)	11 (6%)	2 (1%)	14	41
1	D	174/212 (82%)	157 (90%)	15 (9%)	2 (1%)	14	41
All	All	697/848 (82%)	618 (89%)	68 (10%)	11 (2%)	9	31

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%)	123 (88%)	17 (12%)	5	15
1	B	139/168 (83%)	129 (93%)	10 (7%)	14	38
1	C	139/168 (83%)	126 (91%)	13 (9%)	8	26
1	D	139/168 (83%)	126 (91%)	13 (9%)	8	26
All	All	557/672 (83%)	504 (90%)	53 (10%)	8	25

All (53) 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	30	ILE
1	A	49	SER
1	A	55	SER
1	A	72	GLU
1	A	82	ILE
1	A	83	SER
1	A	98	ASN
1	A	126	ASN
1	A	131	ILE
1	A	136	GLU
1	A	151	LYS
1	A	172	GLN
1	A	176	ILE
1	B	1	MET
1	B	4	ASP
1	B	5	LEU
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	6	ILE
1	C	16	THR
1	C	21	LEU
1	C	22	LYS
1	C	28	HIS
1	C	98	ASN
1	C	108	VAL
1	C	121	SER
1	C	128	ILE
1	C	179	ILE
1	C	187	GLU
1	C	188	LYS
1	D	5	LEU
1	D	17	LEU
1	D	22	LYS

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Mol	Chain	Res	Type
1	D	55	SER
1	D	72	GLU
1	D	82	ILE
1	D	83	SER
1	D	108	VAL
1	D	128	ILE
1	D	172	GLN
1	D	175	HIS
1	D	176	ILE
1	D	192	LYS

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

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

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

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	-	16,17,17	3.57	9 (56%)	20,24,24	3.16	9 (45%)

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	8.82	1.88	1.60
2	B	900	I22	C7-C6	4.94	1.58	1.51
2	B	900	I22	O5-C5	4.84	1.54	1.43
2	B	900	I22	O1-C1	4.45	1.57	1.41
2	B	900	I22	O3-C3	3.93	1.50	1.42
2	B	900	I22	O6-C6	3.36	1.50	1.43
2	B	900	I22	C4-C5	2.96	1.59	1.53
2	B	900	I22	O7-C7	2.70	1.55	1.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	900	I22	P1-O10	2.52	1.64	1.54

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	900	I22	C6-C5-C4	-6.06	102.99	112.47
2	B	900	I22	O4-C4-C5	-5.84	95.88	109.47
2	B	900	I22	O5-C5-C6	-5.39	95.80	108.81
2	B	900	I22	C5-C4-C3	-4.59	105.47	113.60
2	B	900	I22	O3-C3-C4	-4.33	101.28	110.45
2	B	900	I22	O7-C7-C6	-3.99	98.70	109.36
2	B	900	I22	O6-C6-C7	-3.89	101.17	109.92
2	B	900	I22	O6-C6-C5	3.73	118.16	109.10
2	B	900	I22	O5-C5-C4	-2.77	103.02	109.47

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	C4-C5-C6-O6
2	B	900	I22	C7-O7-P1-O9
2	B	900	I22	O6-C6-C7-O7
2	B	900	I22	C3-C4-C5-O5
2	B	900	I22	O4-C4-C5-O5

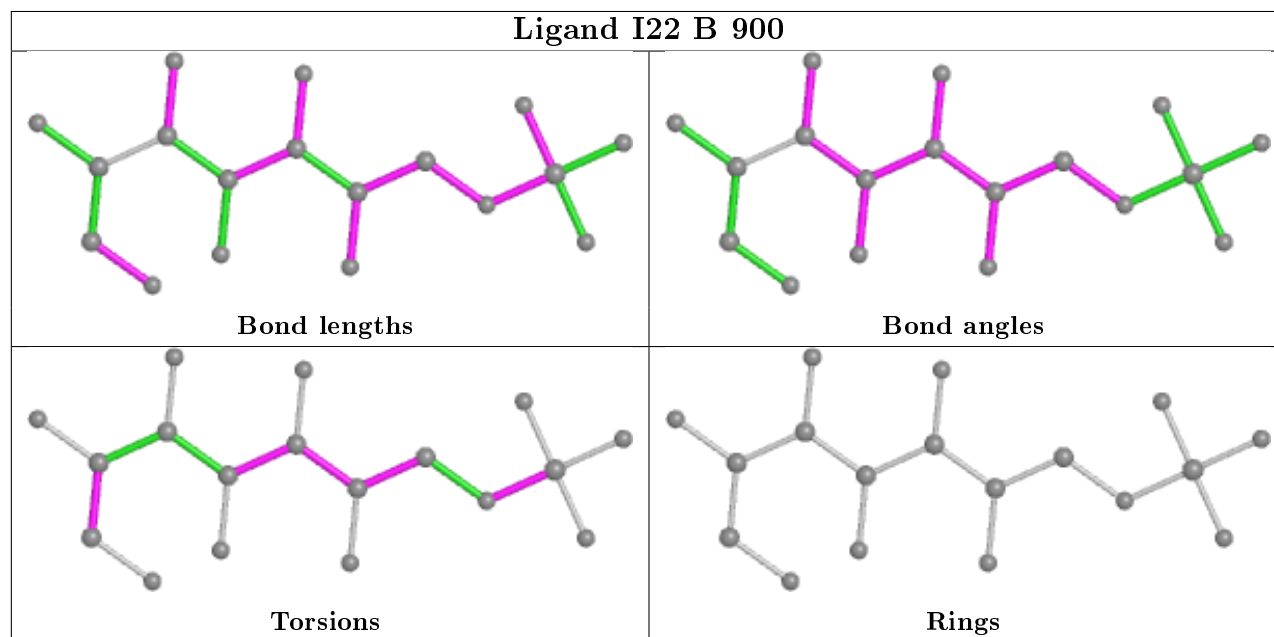
There are no ring outliers.

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%)	-0.51	1 (0%) 89 86	27, 43, 68, 84	0
1	B	178/212 (83%)	-0.55	0 100 100	22, 41, 68, 79	0
1	C	178/212 (83%)	-0.52	1 (0%) 89 86	22, 41, 67, 82	0
1	D	178/212 (83%)	-0.53	0 100 100	22, 43, 69, 84	0
All	All	713/848 (84%)	-0.53	2 (0%) 94 93	22, 42, 68, 84	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	166	GLY	2.7
1	A	191	VAL	2.3

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

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

6.3 Carbohydrates [i](#)

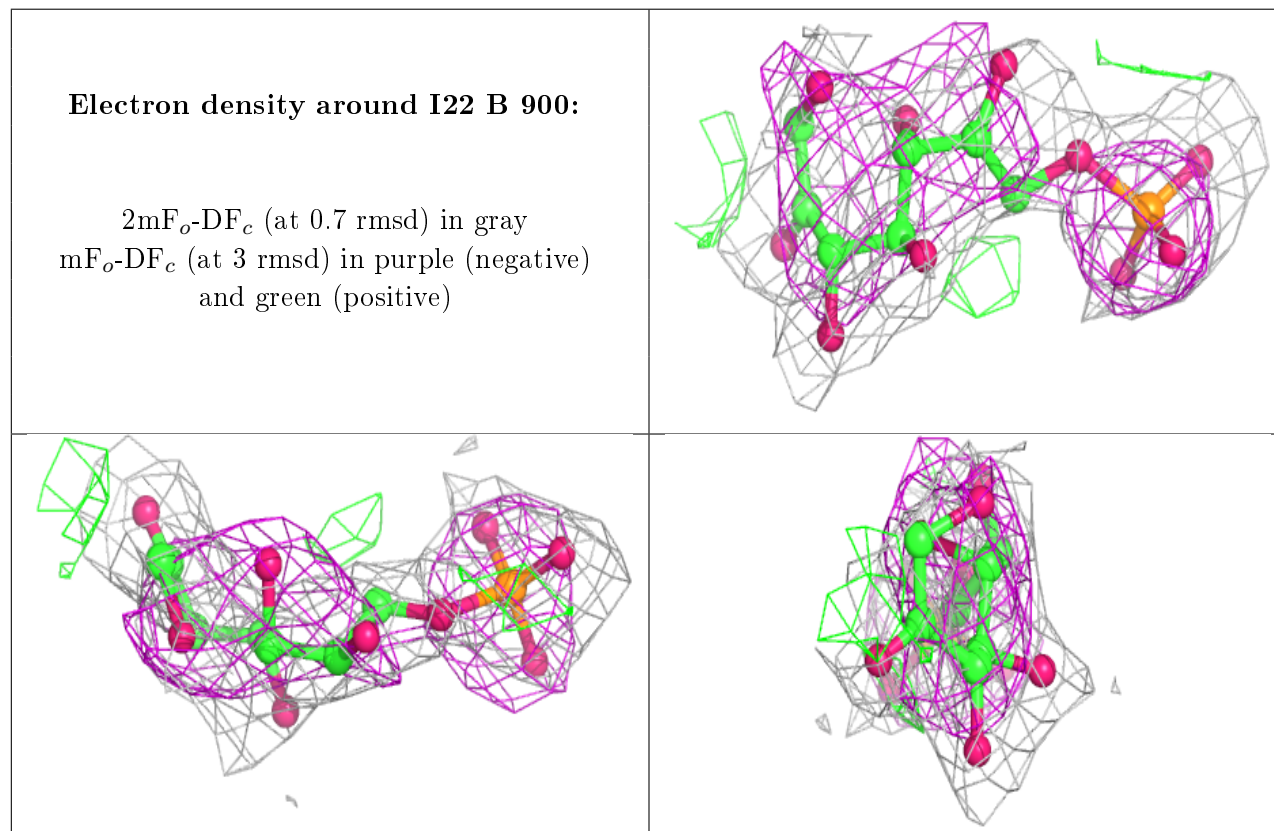
There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. 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(\AA^2)	Q<0.9
2	I22	B	900	18/18	0.88	0.32	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.