



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 02:32 PM UTC

PDB ID : 1HG3 / pdb_00001hg3
Title : Crystal structure of tetrameric TIM from Pyrococcus woesei.
Authors : Walden, H.; Bell, G.S.; Russell, R.J.M.; Siebers, B.; Hensel, R.; Taylor, G.L.
Deposited on : 2000-12-12
Resolution : 2.70 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)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

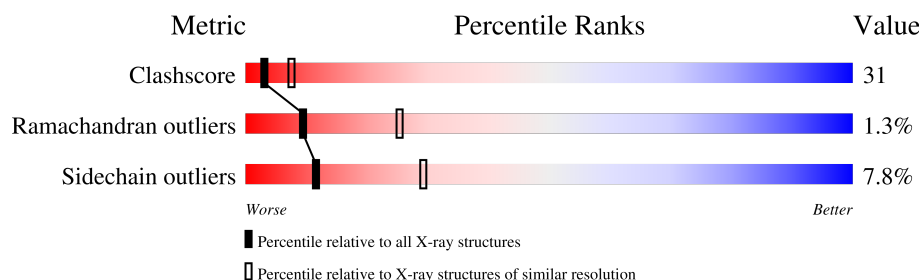
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)

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

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	225	58% 36% 5% •
1	B	225	55% 38% 7%
1	C	225	56% 37% 6% •
1	D	225	59% 34% 6% •
1	E	225	56% 36% 6% •
1	F	225	57% 36% 6% •
1	G	225	56% 37% 5% •
1	H	225	59% 34% 6% •

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	3PP	A	1226	-	-	X	-
2	3PP	B	1226	-	-	X	-
2	3PP	C	1226	-	-	X	-
2	3PP	D	1226	-	X	X	-
2	3PP	E	1226	-	-	X	-
2	3PP	F	1226	-	-	X	-
2	3PP	G	1226	-	-	X	-
2	3PP	H	1226	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 13650 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 TRIOSEPHOSPHATE ISOMERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	224	Total	C	N	O	S	0	0	0
			1654	1055	277	316	6			
1	B	224	Total	C	N	O	S	0	0	0
			1654	1055	277	316	6			
1	C	224	Total	C	N	O	S	0	0	0
			1654	1055	277	316	6			
1	D	224	Total	C	N	O	S	0	0	0
			1654	1055	277	316	6			
1	E	224	Total	C	N	O	S	0	0	0
			1654	1055	277	316	6			
1	F	224	Total	C	N	O	S	0	0	0
			1654	1055	277	316	6			
1	G	224	Total	C	N	O	S	0	0	0
			1654	1055	277	316	6			
1	H	224	Total	C	N	O	S	0	0	0
			1654	1055	277	316	6			

There are 112 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	211	ALA	GLN	SEE REMARK 999	UNP P95583
A	212	LYS	ARG	SEE REMARK 999	UNP P95583
A	213	ASP	SER	SEE REMARK 999	UNP P95583
A	214	PRO	ARG	SEE REMARK 999	UNP P95583
A	216	LYS	GLY	SEE REMARK 999	UNP P95583
A	217	ALA	ASP	SEE REMARK 999	UNP P95583
A	218	ILE	MET	SEE REMARK 999	UNP P95583
A	219	TRP	GLY	SEE REMARK 999	UNP P95583
A	220	ASP	SER	SEE REMARK 999	UNP P95583
A	221	LEU	CYS	SEE REMARK 999	UNP P95583
A	222	VAL	PHE	SEE REMARK 999	UNP P95583
A	223	SER	GLY	SEE REMARK 999	UNP P95583
A	224	GLY	ASN	SEE REMARK 999	UNP P95583

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Chain	Residue	Modelled	Actual	Comment	Reference
A	225	ILE	ASN	SEE REMARK 999	UNP P95583
B	211	ALA	GLN	SEE REMARK 999	UNP P95583
B	212	LYS	ARG	SEE REMARK 999	UNP P95583
B	213	ASP	SER	SEE REMARK 999	UNP P95583
B	214	PRO	ARG	SEE REMARK 999	UNP P95583
B	216	LYS	GLY	SEE REMARK 999	UNP P95583
B	217	ALA	ASP	SEE REMARK 999	UNP P95583
B	218	ILE	MET	SEE REMARK 999	UNP P95583
B	219	TRP	GLY	SEE REMARK 999	UNP P95583
B	220	ASP	SER	SEE REMARK 999	UNP P95583
B	221	LEU	CYS	SEE REMARK 999	UNP P95583
B	222	VAL	PHE	SEE REMARK 999	UNP P95583
B	223	SER	GLY	SEE REMARK 999	UNP P95583
B	224	GLY	ASN	SEE REMARK 999	UNP P95583
B	225	ILE	ASN	SEE REMARK 999	UNP P95583
C	211	ALA	GLN	SEE REMARK 999	UNP P95583
C	212	LYS	ARG	SEE REMARK 999	UNP P95583
C	213	ASP	SER	SEE REMARK 999	UNP P95583
C	214	PRO	ARG	SEE REMARK 999	UNP P95583
C	216	LYS	GLY	SEE REMARK 999	UNP P95583
C	217	ALA	ASP	SEE REMARK 999	UNP P95583
C	218	ILE	MET	SEE REMARK 999	UNP P95583
C	219	TRP	GLY	SEE REMARK 999	UNP P95583
C	220	ASP	SER	SEE REMARK 999	UNP P95583
C	221	LEU	CYS	SEE REMARK 999	UNP P95583
C	222	VAL	PHE	SEE REMARK 999	UNP P95583
C	223	SER	GLY	SEE REMARK 999	UNP P95583
C	224	GLY	ASN	SEE REMARK 999	UNP P95583
C	225	ILE	ASN	SEE REMARK 999	UNP P95583
D	211	ALA	GLN	SEE REMARK 999	UNP P95583
D	212	LYS	ARG	SEE REMARK 999	UNP P95583
D	213	ASP	SER	SEE REMARK 999	UNP P95583
D	214	PRO	ARG	SEE REMARK 999	UNP P95583
D	216	LYS	GLY	SEE REMARK 999	UNP P95583
D	217	ALA	ASP	SEE REMARK 999	UNP P95583
D	218	ILE	MET	SEE REMARK 999	UNP P95583
D	219	TRP	GLY	SEE REMARK 999	UNP P95583
D	220	ASP	SER	SEE REMARK 999	UNP P95583
D	221	LEU	CYS	SEE REMARK 999	UNP P95583
D	222	VAL	PHE	SEE REMARK 999	UNP P95583
D	223	SER	GLY	SEE REMARK 999	UNP P95583
D	224	GLY	ASN	SEE REMARK 999	UNP P95583

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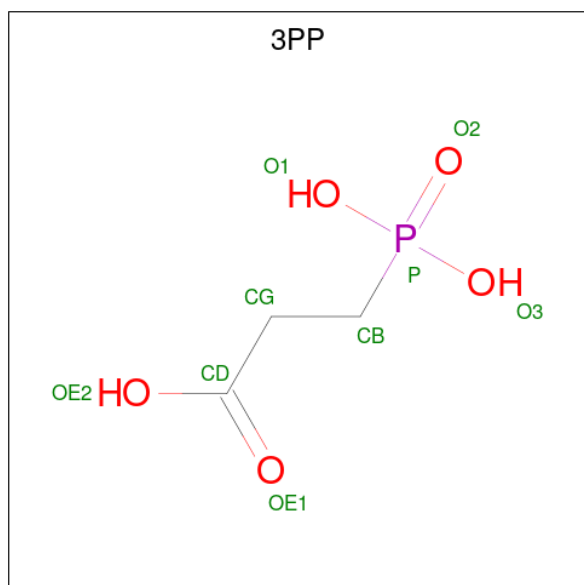
Chain	Residue	Modelled	Actual	Comment	Reference
D	225	ILE	ASN	SEE REMARK 999	UNP P95583
E	211	ALA	GLN	SEE REMARK 999	UNP P95583
E	212	LYS	ARG	SEE REMARK 999	UNP P95583
E	213	ASP	SER	SEE REMARK 999	UNP P95583
E	214	PRO	ARG	SEE REMARK 999	UNP P95583
E	216	LYS	GLY	SEE REMARK 999	UNP P95583
E	217	ALA	ASP	SEE REMARK 999	UNP P95583
E	218	ILE	MET	SEE REMARK 999	UNP P95583
E	219	TRP	GLY	SEE REMARK 999	UNP P95583
E	220	ASP	SER	SEE REMARK 999	UNP P95583
E	221	LEU	CYS	SEE REMARK 999	UNP P95583
E	222	VAL	PHE	SEE REMARK 999	UNP P95583
E	223	SER	GLY	SEE REMARK 999	UNP P95583
E	224	GLY	ASN	SEE REMARK 999	UNP P95583
E	225	ILE	ASN	SEE REMARK 999	UNP P95583
F	211	ALA	GLN	SEE REMARK 999	UNP P95583
F	212	LYS	ARG	SEE REMARK 999	UNP P95583
F	213	ASP	SER	SEE REMARK 999	UNP P95583
F	214	PRO	ARG	SEE REMARK 999	UNP P95583
F	216	LYS	GLY	SEE REMARK 999	UNP P95583
F	217	ALA	ASP	SEE REMARK 999	UNP P95583
F	218	ILE	MET	SEE REMARK 999	UNP P95583
F	219	TRP	GLY	SEE REMARK 999	UNP P95583
F	220	ASP	SER	SEE REMARK 999	UNP P95583
F	221	LEU	CYS	SEE REMARK 999	UNP P95583
F	222	VAL	PHE	SEE REMARK 999	UNP P95583
F	223	SER	GLY	SEE REMARK 999	UNP P95583
F	224	GLY	ASN	SEE REMARK 999	UNP P95583
F	225	ILE	ASN	SEE REMARK 999	UNP P95583
G	211	ALA	GLN	SEE REMARK 999	UNP P95583
G	212	LYS	ARG	SEE REMARK 999	UNP P95583
G	213	ASP	SER	SEE REMARK 999	UNP P95583
G	214	PRO	ARG	SEE REMARK 999	UNP P95583
G	216	LYS	GLY	SEE REMARK 999	UNP P95583
G	217	ALA	ASP	SEE REMARK 999	UNP P95583
G	218	ILE	MET	SEE REMARK 999	UNP P95583
G	219	TRP	GLY	SEE REMARK 999	UNP P95583
G	220	ASP	SER	SEE REMARK 999	UNP P95583
G	221	LEU	CYS	SEE REMARK 999	UNP P95583
G	222	VAL	PHE	SEE REMARK 999	UNP P95583
G	223	SER	GLY	SEE REMARK 999	UNP P95583
G	224	GLY	ASN	SEE REMARK 999	UNP P95583

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Chain	Residue	Modelled	Actual	Comment	Reference
G	225	ILE	ASN	SEE REMARK 999	UNP P95583
H	211	ALA	GLN	SEE REMARK 999	UNP P95583
H	212	LYS	ARG	SEE REMARK 999	UNP P95583
H	213	ASP	SER	SEE REMARK 999	UNP P95583
H	214	PRO	ARG	SEE REMARK 999	UNP P95583
H	216	LYS	GLY	SEE REMARK 999	UNP P95583
H	217	ALA	ASP	SEE REMARK 999	UNP P95583
H	218	ILE	MET	SEE REMARK 999	UNP P95583
H	219	TRP	GLY	SEE REMARK 999	UNP P95583
H	220	ASP	SER	SEE REMARK 999	UNP P95583
H	221	LEU	CYS	SEE REMARK 999	UNP P95583
H	222	VAL	PHE	SEE REMARK 999	UNP P95583
H	223	SER	GLY	SEE REMARK 999	UNP P95583
H	224	GLY	ASN	SEE REMARK 999	UNP P95583
H	225	ILE	ASN	SEE REMARK 999	UNP P95583

- Molecule 2 is 3-PHOSPHONOPROPANOIC ACID (CCD ID: 3PP) (formula: $C_3H_7O_5P$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	O	P	0	0
			9	3	5	1		
2	B	1	Total	C	O	P	0	0
			9	3	5	1		
2	C	1	Total	C	O	P	0	0
			9	3	5	1		
2	D	1	Total	C	O	P	0	0
			9	3	5	1		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	E	1	Total	C	O	P	0	0
			9	3	5	1		
2	F	1	Total	C	O	P	0	0
			9	3	5	1		
2	G	1	Total	C	O	P	0	0
			9	3	5	1		
2	H	1	Total	C	O	P	0	0
			9	3	5	1		

- Molecule 3 is water.

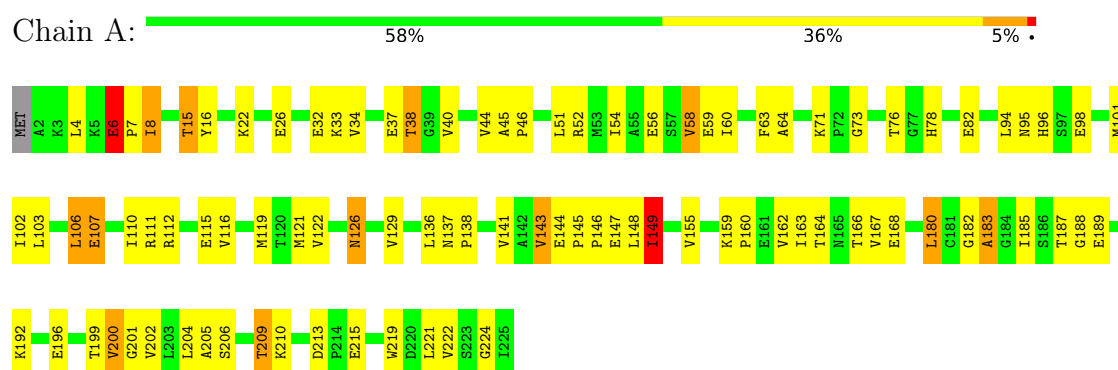
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	38	Total	O	0	0
			38	38		
3	B	36	Total	O	0	0
			36	36		
3	C	52	Total	O	0	0
			52	52		
3	D	46	Total	O	0	0
			46	46		
3	E	43	Total	O	0	0
			43	43		
3	F	41	Total	O	0	0
			41	41		
3	G	46	Total	O	0	0
			46	46		
3	H	44	Total	O	0	0
			44	44		

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. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

Note EDS was not executed.

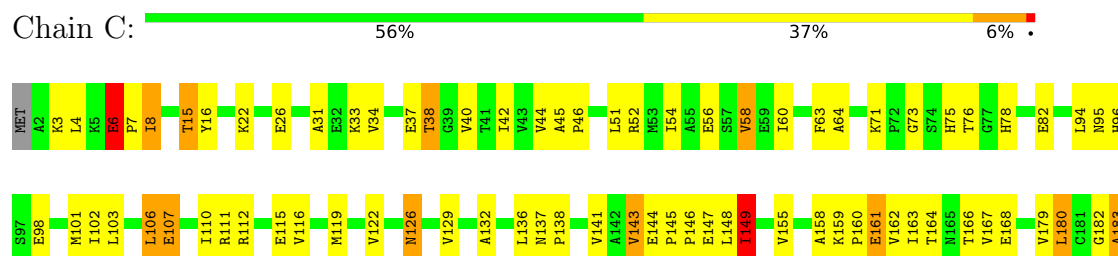
• Molecule 1: TRIOSEPHOSPHATE ISOMERASE



• Molecule 1: TRIOSEPHOSPHATE ISOMERASE



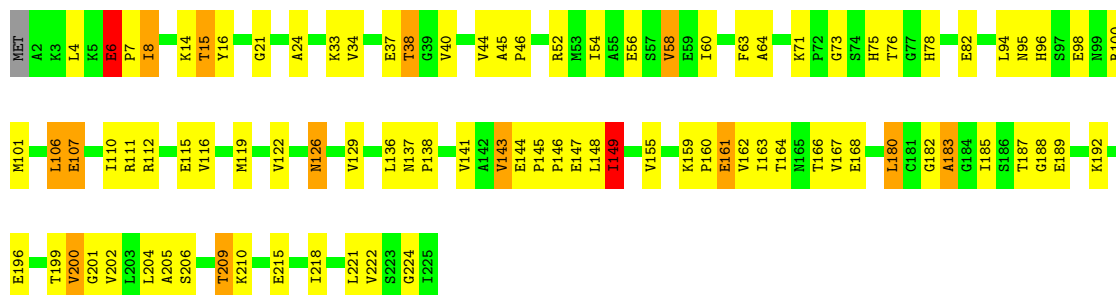
• Molecule 1: TRIOSEPHOSPHATE ISOMERASE





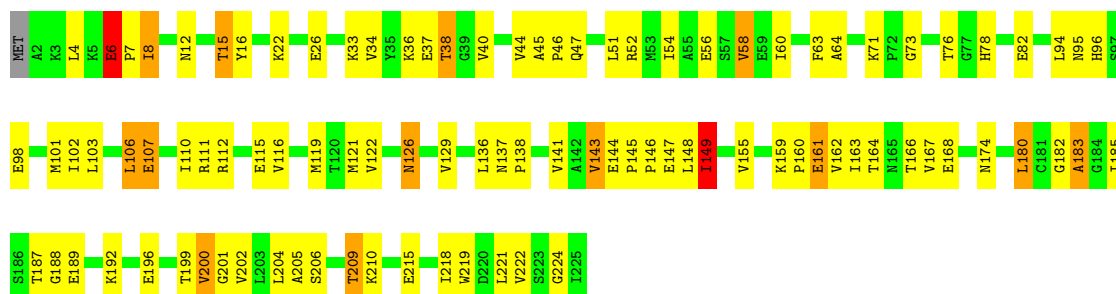
• Molecule 1: TRIOSEPHOSPHATE ISOMERASE

Chain D: 59% 34% 6% .



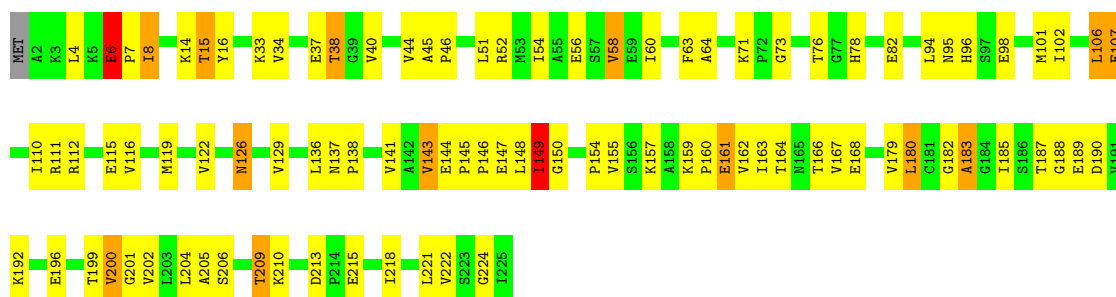
• Molecule 1: TRIOSEPHOSPHATE ISOMERASE

Chain E: 56% 36% 6% .



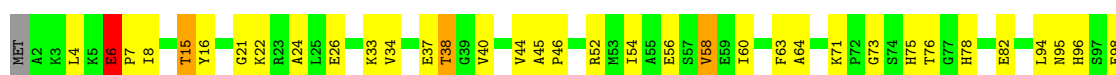
• Molecule 1: TRIOSEPHOSPHATE ISOMERASE

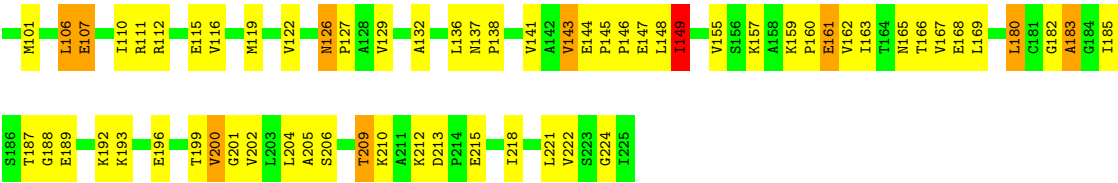
Chain F: 57% 36% 6% .



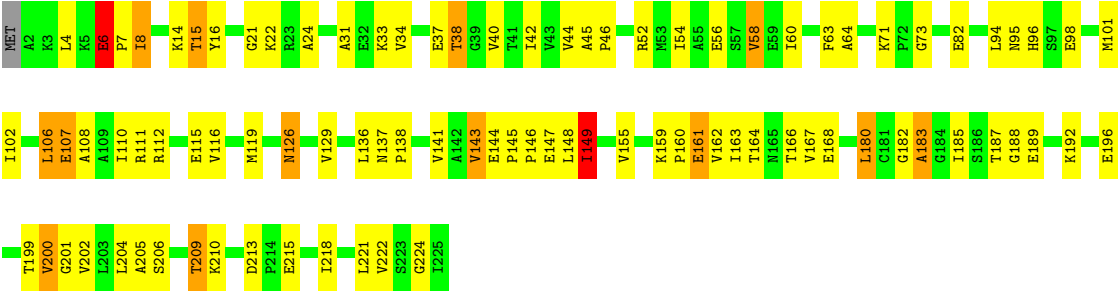
• Molecule 1: TRIOSEPHOSPHATE ISOMERASE

Chain G: 56% 37% 5% .





• Molecule 1: TRIOSEPHOSPHATE ISOMERASE



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	79.00Å 89.10Å 145.10Å 90.00° 92.80° 90.00°	Depositor
Resolution (Å)	6.00 – 2.70	Depositor
% Data completeness (in resolution range)	98.1 (6.00-2.70)	Depositor
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS 0.9	Depositor
R, R_{free}	0.254 , 0.286	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	13650	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: 3PP

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.46	0/1678	0.98	8/2282 (0.4%)
1	B	0.52	1/1678 (0.1%)	1.08	16/2282 (0.7%)
1	C	0.49	0/1678	0.98	7/2282 (0.3%)
1	D	0.46	0/1678	0.99	5/2282 (0.2%)
1	E	0.46	0/1678	0.98	8/2282 (0.4%)
1	F	0.46	0/1678	0.99	9/2282 (0.4%)
1	G	0.49	0/1678	0.99	7/2282 (0.3%)
1	H	0.46	0/1678	0.98	7/2282 (0.3%)
All	All	0.48	1/13424 (0.0%)	1.00	67/18256 (0.4%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	7	PRO	N-CD	5.44	1.55	1.47

All (67) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	8	ILE	N-CA-C	13.91	127.52	109.30
1	B	6	GLU	CA-C-N	7.46	129.16	119.84
1	B	6	GLU	C-N-CA	7.46	129.16	119.84
1	B	7	PRO	CA-C-N	6.94	130.82	120.69
1	B	7	PRO	C-N-CA	6.94	130.82	120.69
1	A	6	GLU	N-CA-C	6.83	124.91	109.81
1	C	6	GLU	N-CA-C	6.80	124.83	109.81
1	E	6	GLU	N-CA-C	6.79	124.82	109.81
1	F	8	ILE	N-CA-C	6.71	118.26	108.53
1	F	6	GLU	N-CA-C	6.65	124.51	109.81
1	A	8	ILE	N-CA-C	6.63	118.14	108.53
1	E	8	ILE	N-CA-C	6.63	118.14	108.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	6	GLU	N-CA-C	6.62	124.45	109.81
1	D	6	GLU	N-CA-C	6.58	124.36	109.81
1	H	8	ILE	N-CA-C	6.56	118.04	108.53
1	D	8	ILE	N-CA-C	6.50	117.95	108.53
1	C	8	ILE	N-CA-C	6.37	117.76	108.53
1	B	6	GLU	O-C-N	-6.36	114.00	121.32
1	C	64	ALA	N-CA-C	-6.34	102.36	110.53
1	F	64	ALA	N-CA-C	-6.13	102.62	110.53
1	G	180	LEU	N-CA-C	-6.02	100.19	109.52
1	E	64	ALA	N-CA-C	-5.94	102.53	110.55
1	B	4	LEU	CA-C-N	-5.93	112.70	121.24
1	B	4	LEU	C-N-CA	-5.93	112.70	121.24
1	G	64	ALA	N-CA-C	-5.92	102.56	110.55
1	B	64	ALA	N-CA-C	-5.83	102.68	110.55
1	G	8	ILE	N-CA-C	5.74	116.74	108.48
1	F	51	LEU	N-CA-C	5.70	117.95	111.11
1	A	64	ALA	N-CA-C	-5.65	102.92	110.55
1	B	180	LEU	N-CA-C	-5.65	100.77	109.52
1	H	180	LEU	N-CA-C	-5.58	100.87	109.52
1	D	64	ALA	N-CA-C	-5.57	103.02	110.55
1	C	180	LEU	N-CA-C	-5.53	100.94	109.52
1	B	51	LEU	N-CA-C	5.46	117.67	111.11
1	H	64	ALA	N-CA-C	-5.43	103.21	110.55
1	F	213	ASP	CA-C-N	5.43	124.89	119.24
1	F	213	ASP	C-N-CA	5.43	124.89	119.24
1	A	180	LEU	N-CA-C	-5.42	101.12	109.52
1	E	174	ASN	CA-C-N	5.40	125.07	119.56
1	E	174	ASN	C-N-CA	5.40	125.07	119.56
1	E	180	LEU	N-CA-C	-5.31	101.30	109.52
1	D	180	LEU	N-CA-C	-5.27	101.35	109.52
1	B	174	ASN	CA-C-N	5.27	124.93	119.56
1	B	174	ASN	C-N-CA	5.27	124.93	119.56
1	B	213	ASP	CA-C-N	5.24	124.69	119.24
1	B	213	ASP	C-N-CA	5.24	124.69	119.24
1	G	213	ASP	CA-C-N	5.23	124.67	119.24
1	G	213	ASP	C-N-CA	5.23	124.67	119.24
1	A	51	LEU	N-CA-C	5.22	117.38	111.11
1	F	180	LEU	N-CA-C	-5.17	101.50	109.52
1	F	149	ILE	N-CA-C	5.17	120.09	109.34
1	C	149	ILE	N-CA-C	5.16	120.07	109.34
1	H	213	ASP	CA-C-N	5.16	124.61	119.24
1	H	213	ASP	C-N-CA	5.16	124.61	119.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	149	ILE	N-CA-C	5.14	120.03	109.34
1	A	213	ASP	CA-C-N	5.14	124.58	119.24
1	A	213	ASP	C-N-CA	5.14	124.58	119.24
1	E	149	ILE	N-CA-C	5.13	120.01	109.34
1	H	149	ILE	N-CA-C	5.12	119.99	109.34
1	B	149	ILE	N-CA-C	5.12	119.99	109.34
1	D	149	ILE	N-CA-C	5.12	119.98	109.34
1	G	149	ILE	N-CA-C	5.10	119.94	109.34
1	C	179	VAL	N-CA-C	5.08	115.89	108.53
1	G	6	GLU	N-CA-C	5.06	121.00	109.81
1	F	179	VAL	N-CA-C	5.06	115.87	108.53
1	C	51	LEU	N-CA-C	5.04	117.16	111.11
1	E	51	LEU	N-CA-C	5.02	117.14	111.11

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	1654	0	1744	118	1
1	B	1654	0	1744	134	0
1	C	1654	0	1744	108	1
1	D	1654	0	1744	101	0
1	E	1654	0	1744	128	1
1	F	1654	0	1744	94	2
1	G	1654	0	1744	106	3
1	H	1654	0	1744	96	0
2	A	9	0	4	8	0
2	B	9	0	4	6	0
2	C	9	0	4	6	0
2	D	9	0	4	4	0
2	E	9	0	4	7	0
2	F	9	0	4	5	0
2	G	9	0	4	6	0
2	H	9	0	4	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	38	0	0	0	0
3	B	36	0	0	0	0
3	C	52	0	0	4	0
3	D	46	0	0	5	0
3	E	43	0	0	2	0
3	F	41	0	0	3	0
3	G	46	0	0	5	0
3	H	44	0	0	3	0
All	All	13650	0	13984	859	4

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 31.

All (859) 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:E:34:VAL:HG22	1:E:219:TRP:CE2	1.53	1.43
1:A:34:VAL:HG22	1:A:219:TRP:CE2	1.54	1.42
1:B:34:VAL:HG22	1:B:219:TRP:CE2	1.54	1.40
1:A:215:GLU:HG2	1:A:219:TRP:NE1	1.44	1.32
1:B:215:GLU:HG2	1:B:219:TRP:NE1	1.44	1.31
1:E:215:GLU:HG2	1:E:219:TRP:NE1	1.45	1.31
1:E:34:VAL:HA	1:E:219:TRP:CH2	1.75	1.21
1:B:34:VAL:HA	1:B:219:TRP:CH2	1.74	1.20
1:B:215:GLU:CG	1:B:219:TRP:HE1	1.55	1.19
1:A:34:VAL:HA	1:A:219:TRP:CH2	1.77	1.19
1:A:215:GLU:CG	1:A:219:TRP:HE1	1.55	1.18
1:E:215:GLU:CG	1:E:219:TRP:HE1	1.56	1.17
1:F:187:THR:HG22	1:F:189:GLU:H	1.19	1.06
1:E:37:GLU:CG	1:E:219:TRP:HH2	1.68	1.05
1:B:37:GLU:CG	1:B:219:TRP:HH2	1.69	1.03
1:E:187:THR:HG22	1:E:189:GLU:H	1.23	1.03
1:D:187:THR:HG22	1:D:189:GLU:H	1.19	1.03
1:A:187:THR:HG22	1:A:189:GLU:H	1.23	1.03
1:B:187:THR:HG22	1:B:189:GLU:H	1.23	1.02
1:C:187:THR:HG22	1:C:189:GLU:H	1.22	1.01
1:H:187:THR:HG22	1:H:189:GLU:H	1.22	1.01
1:B:162:VAL:O	1:B:166:THR:HG23	1.60	1.01
1:G:187:THR:HG22	1:G:189:GLU:H	1.23	1.00
1:E:34:VAL:CG2	1:E:219:TRP:CE2	2.47	0.98
1:A:34:VAL:CG2	1:A:219:TRP:CE2	2.48	0.97

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:215:GLU:O	1:E:219:TRP:CD1	2.21	0.94
1:A:34:VAL:HG22	1:A:219:TRP:CD2	2.01	0.94
1:E:34:VAL:HG22	1:E:219:TRP:CD2	2.02	0.94
1:E:37:GLU:HG3	1:E:219:TRP:CH2	2.03	0.94
1:E:34:VAL:HG22	1:E:219:TRP:CZ2	2.03	0.93
1:B:34:VAL:CG2	1:B:219:TRP:CE2	2.49	0.93
1:B:34:VAL:HG22	1:B:219:TRP:CD2	2.03	0.92
1:F:199:THR:HG22	1:F:201:GLY:H	1.33	0.92
1:A:199:THR:HG22	1:A:201:GLY:H	1.34	0.92
1:B:199:THR:HG22	1:B:201:GLY:H	1.35	0.92
1:E:199:THR:HG22	1:E:201:GLY:H	1.34	0.92
1:B:37:GLU:HG3	1:B:219:TRP:CH2	2.05	0.92
1:A:34:VAL:HG22	1:A:219:TRP:CZ2	2.04	0.92
1:B:34:VAL:HG22	1:B:219:TRP:CZ2	2.05	0.91
1:E:37:GLU:CG	1:E:219:TRP:CH2	2.53	0.91
1:C:199:THR:HG22	1:C:201:GLY:H	1.34	0.91
1:B:37:GLU:CG	1:B:219:TRP:CH2	2.54	0.91
1:H:199:THR:HG22	1:H:201:GLY:H	1.33	0.90
1:D:199:THR:HG22	1:D:201:GLY:H	1.34	0.90
1:G:199:THR:HG22	1:G:201:GLY:H	1.35	0.90
1:B:215:GLU:HG2	1:B:219:TRP:HE1	0.75	0.89
1:A:215:GLU:HG2	1:A:219:TRP:HE1	0.75	0.89
1:A:215:GLU:O	1:A:219:TRP:CD1	2.27	0.88
1:E:215:GLU:HG2	1:E:219:TRP:HE1	0.75	0.88
1:B:37:GLU:HG3	1:B:219:TRP:HH2	1.38	0.87
1:A:38:THR:HG22	1:A:40:VAL:H	1.40	0.87
1:E:38:THR:HG22	1:E:40:VAL:H	1.39	0.87
1:C:96:HIS:HE1	2:C:1226:3PP:OE1	1.56	0.87
1:B:215:GLU:O	1:B:219:TRP:CD1	2.27	0.86
1:A:34:VAL:HG13	1:A:219:TRP:CE3	2.11	0.86
1:B:37:GLU:CD	1:B:219:TRP:CH2	2.53	0.86
1:E:37:GLU:CD	1:E:219:TRP:CH2	2.54	0.86
1:B:38:THR:HG22	1:B:40:VAL:H	1.40	0.85
1:F:38:THR:HG22	1:F:40:VAL:H	1.40	0.85
1:C:6:GLU:O	1:C:8:ILE:HG13	1.77	0.85
1:G:38:THR:HG22	1:G:40:VAL:H	1.40	0.85
1:E:34:VAL:HG13	1:E:219:TRP:CE3	2.12	0.85
1:H:38:THR:HG22	1:H:40:VAL:H	1.40	0.85
1:B:5:LYS:HB2	1:B:41:THR:OG1	1.76	0.84
1:E:6:GLU:O	1:E:8:ILE:HG13	1.78	0.84
1:C:38:THR:HG22	1:C:40:VAL:H	1.42	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:37:GLU:HG3	1:E:219:TRP:HH2	1.36	0.84
1:H:6:GLU:O	1:H:8:ILE:HG13	1.79	0.83
1:B:34:VAL:HG13	1:B:219:TRP:CE3	2.13	0.83
1:D:38:THR:HG22	1:D:40:VAL:H	1.43	0.83
1:B:96:HIS:HD2	1:B:98:GLU:H	1.26	0.82
1:A:6:GLU:O	1:A:8:ILE:HG13	1.79	0.82
1:H:205:ALA:N	2:H:1226:3PP:HCB1	1.95	0.82
1:G:96:HIS:HD2	1:G:98:GLU:H	1.26	0.82
1:H:96:HIS:HD2	1:H:98:GLU:H	1.29	0.81
1:C:96:HIS:HD2	1:C:98:GLU:H	1.27	0.81
1:F:6:GLU:O	1:F:8:ILE:HG13	1.81	0.81
1:D:6:GLU:O	1:D:8:ILE:HG13	1.81	0.81
1:G:126:ASN:ND2	1:G:129:VAL:H	1.79	0.81
1:F:143:VAL:HG21	1:F:166:THR:HG21	1.64	0.80
1:H:58:VAL:CG1	1:H:60:ILE:HG12	2.12	0.80
1:B:126:ASN:ND2	1:B:129:VAL:H	1.79	0.79
1:E:96:HIS:HD2	1:E:98:GLU:H	1.28	0.79
1:F:96:HIS:HD2	1:F:98:GLU:H	1.29	0.79
1:H:126:ASN:ND2	1:H:129:VAL:H	1.81	0.79
1:D:96:HIS:HD2	1:D:98:GLU:H	1.30	0.79
1:F:126:ASN:ND2	1:F:129:VAL:H	1.80	0.79
1:A:126:ASN:ND2	1:A:129:VAL:H	1.80	0.79
1:D:58:VAL:CG1	1:D:60:ILE:HG12	2.13	0.79
1:D:126:ASN:ND2	1:D:129:VAL:H	1.81	0.79
1:C:126:ASN:ND2	1:C:129:VAL:H	1.81	0.79
1:F:187:THR:HG22	1:F:189:GLU:N	1.98	0.79
1:A:96:HIS:HD2	1:A:98:GLU:H	1.30	0.79
1:D:126:ASN:C	1:D:126:ASN:HD22	1.91	0.79
1:E:58:VAL:CG1	1:E:60:ILE:HG12	2.13	0.79
1:B:126:ASN:C	1:B:126:ASN:HD22	1.91	0.78
1:D:143:VAL:HG21	1:D:166:THR:HG21	1.65	0.78
1:H:126:ASN:C	1:H:126:ASN:HD22	1.92	0.78
1:E:126:ASN:ND2	1:E:129:VAL:H	1.81	0.78
1:E:126:ASN:C	1:E:126:ASN:HD22	1.92	0.77
1:G:58:VAL:CG1	1:G:60:ILE:HG12	2.14	0.77
1:A:58:VAL:CG1	1:A:60:ILE:HG12	2.15	0.77
1:C:58:VAL:CG1	1:C:60:ILE:HG12	2.15	0.77
1:C:187:THR:HG22	1:C:189:GLU:N	1.99	0.77
1:A:126:ASN:HD22	1:A:126:ASN:C	1.93	0.76
1:B:187:THR:HG22	1:B:189:GLU:N	2.01	0.76
1:A:187:THR:HG22	1:A:189:GLU:N	1.99	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:192:LYS:O	1:B:196:GLU:HG3	1.86	0.76
1:H:146:PRO:HA	1:H:149:ILE:HG12	1.67	0.76
1:F:126:ASN:HD22	1:F:126:ASN:C	1.92	0.76
1:H:143:VAL:HG21	1:H:166:THR:HG21	1.67	0.76
1:B:146:PRO:HA	1:B:149:ILE:HG12	1.66	0.76
1:E:187:THR:HG22	1:E:189:GLU:N	2.00	0.76
1:D:146:PRO:HA	1:D:149:ILE:HG12	1.67	0.76
1:G:187:THR:HG22	1:G:189:GLU:N	2.00	0.76
1:C:146:PRO:HA	1:C:149:ILE:HG12	1.66	0.75
1:E:143:VAL:HG21	1:E:166:THR:HG21	1.67	0.75
1:E:215:GLU:O	1:E:219:TRP:HD1	1.67	0.75
1:F:58:VAL:CG1	1:F:60:ILE:HG12	2.16	0.75
1:A:146:PRO:HA	1:A:149:ILE:HG12	1.67	0.75
1:F:38:THR:HG21	1:F:222:VAL:HG11	1.68	0.75
1:F:143:VAL:CG2	1:F:166:THR:HG21	2.16	0.75
1:E:143:VAL:CG2	1:E:166:THR:HG21	2.17	0.75
1:A:143:VAL:CG2	1:A:166:THR:HG21	2.17	0.75
1:B:205:ALA:N	2:B:1226:3PP:HCB1	2.02	0.75
1:D:143:VAL:CG2	1:D:166:THR:HG21	2.16	0.75
1:D:187:THR:HG22	1:D:189:GLU:N	1.97	0.75
1:A:38:THR:HG21	1:A:222:VAL:HG11	1.69	0.74
1:G:38:THR:HG21	1:G:222:VAL:HG11	1.68	0.74
1:B:162:VAL:O	1:B:166:THR:CG2	2.35	0.74
1:G:126:ASN:C	1:G:126:ASN:HD22	1.94	0.74
1:C:126:ASN:C	1:C:126:ASN:HD22	1.94	0.74
1:C:192:LYS:O	1:C:196:GLU:HG3	1.86	0.74
1:D:38:THR:HG21	1:D:222:VAL:HG11	1.69	0.74
1:A:143:VAL:HG21	1:A:166:THR:HG21	1.68	0.74
1:F:146:PRO:HA	1:F:149:ILE:HG12	1.68	0.74
1:B:38:THR:HG21	1:B:222:VAL:HG11	1.69	0.74
1:B:209:THR:HG22	1:B:210:LYS:HG2	1.69	0.74
1:C:143:VAL:HG21	1:C:166:THR:HG21	1.68	0.74
1:E:38:THR:HG21	1:E:222:VAL:HG11	1.70	0.74
1:E:146:PRO:HA	1:E:149:ILE:HG12	1.68	0.74
1:G:146:PRO:HA	1:G:149:ILE:HG12	1.68	0.74
1:D:209:THR:HG22	1:D:210:LYS:HG2	1.69	0.74
1:F:209:THR:HG22	1:F:210:LYS:HG2	1.70	0.74
1:H:38:THR:HG21	1:H:222:VAL:HG11	1.70	0.74
1:H:192:LYS:O	1:H:196:GLU:HG3	1.87	0.74
1:H:209:THR:HG22	1:H:210:LYS:HG2	1.69	0.74
1:G:192:LYS:O	1:G:196:GLU:HG3	1.86	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:143:VAL:HG11	1:B:166:THR:HG21	1.68	0.73
1:C:143:VAL:CG2	1:C:166:THR:HG21	2.17	0.73
1:C:38:THR:HG21	1:C:222:VAL:HG11	1.69	0.73
1:H:143:VAL:CG2	1:H:166:THR:HG21	2.18	0.73
1:E:205:ALA:N	2:E:1226:3PP:HCB1	2.03	0.73
1:F:205:ALA:N	2:F:1226:3PP:HCB1	2.03	0.73
1:B:215:GLU:HG2	1:B:219:TRP:CD1	2.24	0.73
1:D:192:LYS:O	1:D:196:GLU:HG3	1.89	0.73
1:F:192:LYS:O	1:F:196:GLU:HG3	1.89	0.73
1:H:187:THR:HG22	1:H:189:GLU:N	2.00	0.73
1:A:209:THR:HG22	1:A:210:LYS:HG2	1.70	0.72
1:A:96:HIS:HE1	2:A:1226:3PP:OE1	1.71	0.72
1:G:165:ASN:O	1:G:168:GLU:HG2	1.89	0.72
1:E:192:LYS:O	1:E:196:GLU:HG3	1.89	0.72
1:E:209:THR:HG22	1:E:210:LYS:HG2	1.71	0.72
1:G:96:HIS:HE1	2:G:1226:3PP:OE1	1.73	0.72
1:B:215:GLU:O	1:B:219:TRP:HD1	1.73	0.71
1:G:143:VAL:CG2	1:G:166:THR:HG21	2.19	0.71
1:B:96:HIS:CD2	1:B:98:GLU:H	2.08	0.71
1:E:34:VAL:HA	1:E:219:TRP:CZ2	2.24	0.71
1:G:209:THR:HG22	1:G:210:LYS:HG2	1.71	0.71
1:A:215:GLU:HG2	1:A:219:TRP:CD1	2.24	0.71
1:G:143:VAL:HG21	1:G:166:THR:HG21	1.70	0.71
1:B:34:VAL:HA	1:B:219:TRP:CZ2	2.25	0.71
1:D:46:PRO:HB3	1:D:54:ILE:HD12	1.73	0.71
1:B:6:GLU:O	1:B:200:VAL:O	2.08	0.71
1:G:46:PRO:HB3	1:G:54:ILE:HD12	1.73	0.71
1:A:192:LYS:O	1:A:196:GLU:HG3	1.90	0.71
1:F:159:LYS:N	1:F:160:PRO:HD3	2.06	0.71
1:H:46:PRO:HB3	1:H:54:ILE:HD12	1.73	0.71
1:H:96:HIS:CD2	1:H:98:GLU:H	2.09	0.71
1:E:215:GLU:HG2	1:E:219:TRP:CD1	2.25	0.71
1:G:96:HIS:CD2	1:G:98:GLU:H	2.08	0.71
1:B:159:LYS:N	1:B:160:PRO:HD3	2.06	0.70
1:C:96:HIS:CD2	1:C:98:GLU:H	2.08	0.70
1:A:159:LYS:N	1:A:160:PRO:HD3	2.06	0.70
1:C:193:LYS:HD2	3:C:2049:HOH:O	1.90	0.70
1:E:46:PRO:HB3	1:E:54:ILE:HD12	1.72	0.70
1:E:96:HIS:CD2	1:E:98:GLU:H	2.08	0.70
1:E:159:LYS:N	1:E:160:PRO:HD3	2.06	0.70
1:D:96:HIS:CD2	1:D:98:GLU:H	2.10	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:46:PRO:HB3	1:A:54:ILE:HD12	1.73	0.70
1:G:126:ASN:HD21	1:G:129:VAL:H	1.40	0.70
1:D:159:LYS:N	1:D:160:PRO:HD3	2.06	0.70
1:B:58:VAL:CG1	1:B:60:ILE:HG12	2.22	0.69
1:B:46:PRO:HB3	1:B:54:ILE:HD12	1.73	0.69
1:C:46:PRO:HB3	1:C:54:ILE:HD12	1.74	0.69
1:A:34:VAL:HA	1:A:219:TRP:CZ2	2.26	0.69
1:H:159:LYS:N	1:H:160:PRO:HD3	2.07	0.69
1:F:46:PRO:HB3	1:F:54:ILE:HD12	1.73	0.69
1:F:96:HIS:CD2	1:F:98:GLU:H	2.10	0.69
1:C:209:THR:HG22	1:C:210:LYS:HG2	1.74	0.69
1:A:96:HIS:CD2	1:A:98:GLU:H	2.09	0.69
1:A:126:ASN:HD21	1:A:129:VAL:H	1.40	0.69
1:B:126:ASN:HD21	1:B:129:VAL:H	1.38	0.69
1:C:73:GLY:O	1:D:15:THR:HG22	1.93	0.68
1:E:204:LEU:HA	2:E:1226:3PP:HCB2	1.75	0.68
1:E:107:GLU:HG2	1:G:129:VAL:HG23	1.75	0.68
1:E:126:ASN:HD21	1:E:129:VAL:H	1.40	0.68
1:A:215:GLU:O	1:A:219:TRP:HD1	1.72	0.68
1:D:205:ALA:N	2:D:1226:3PP:HCB1	2.08	0.68
1:A:107:GLU:HG2	1:C:129:VAL:HG23	1.76	0.68
1:C:126:ASN:HD21	1:C:129:VAL:H	1.43	0.67
1:A:205:ALA:CB	2:A:1226:3PP:HCB1	2.24	0.67
1:B:37:GLU:CD	1:B:219:TRP:HH2	1.97	0.67
1:E:15:THR:HG22	1:F:73:GLY:O	1.95	0.67
1:G:159:LYS:N	1:G:160:PRO:HD3	2.10	0.67
1:A:15:THR:HG22	1:B:73:GLY:O	1.96	0.66
1:C:159:LYS:N	1:C:160:PRO:HD3	2.10	0.66
1:A:34:VAL:HG13	1:A:219:TRP:CZ3	2.29	0.66
1:B:6:GLU:O	1:B:200:VAL:HG13	1.95	0.66
1:F:126:ASN:HD21	1:F:129:VAL:H	1.41	0.66
1:H:126:ASN:HD21	1:H:129:VAL:H	1.41	0.66
1:E:73:GLY:O	1:F:15:THR:HG22	1.95	0.65
1:E:34:VAL:HG13	1:E:219:TRP:CZ3	2.31	0.65
1:E:94:LEU:HD22	1:E:101:MET:HE1	1.79	0.65
1:H:162:VAL:HG13	3:H:2038:HOH:O	1.96	0.65
1:H:94:LEU:HD22	1:H:101:MET:HE1	1.79	0.65
1:D:126:ASN:HD21	1:D:129:VAL:H	1.43	0.64
1:B:94:LEU:HD22	1:B:101:MET:HE1	1.78	0.64
1:D:94:LEU:HD22	1:D:101:MET:HE1	1.80	0.64
1:D:58:VAL:HG13	1:D:60:ILE:HG12	1.79	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:73:GLY:O	1:H:15:THR:HG22	1.98	0.63
1:G:94:LEU:HB3	1:G:101:MET:CE	2.28	0.63
1:B:180:LEU:HA	1:B:199:THR:HG23	1.80	0.63
1:C:15:THR:HG22	1:D:73:GLY:O	1.98	0.63
1:E:180:LEU:HA	1:E:199:THR:HG23	1.80	0.63
1:G:7:PRO:HB2	3:G:2002:HOH:O	1.98	0.63
1:H:58:VAL:HG13	1:H:60:ILE:HG12	1.80	0.63
1:G:180:LEU:HA	1:G:199:THR:HG23	1.80	0.63
1:A:205:ALA:N	2:A:1226:3PP:HCB1	2.13	0.63
1:C:180:LEU:HA	1:C:199:THR:HG23	1.80	0.63
1:E:95:ASN:H	1:E:101:MET:HE2	1.63	0.63
1:F:58:VAL:HG13	1:F:60:ILE:HG12	1.79	0.63
1:A:94:LEU:HD22	1:A:101:MET:HE1	1.80	0.63
1:B:58:VAL:HG12	1:B:60:ILE:HG12	1.80	0.63
1:C:95:ASN:H	1:C:101:MET:HE2	1.64	0.62
1:A:15:THR:HG23	1:A:15:THR:O	1.97	0.62
1:A:180:LEU:HA	1:A:199:THR:HG23	1.81	0.62
1:C:205:ALA:N	2:C:1226:3PP:HCB1	2.15	0.62
1:E:159:LYS:HE2	1:E:162:VAL:HG21	1.81	0.62
1:E:215:GLU:C	1:E:219:TRP:HD1	2.07	0.62
1:F:95:ASN:H	1:F:101:MET:HE2	1.65	0.62
1:F:180:LEU:HA	1:F:199:THR:HG23	1.81	0.62
1:E:144:GLU:OE2	1:E:182:GLY:HA3	1.99	0.62
1:E:163:ILE:O	1:E:167:VAL:HG23	2.00	0.62
1:E:185:ILE:HG21	1:E:202:VAL:HG11	1.82	0.62
1:H:95:ASN:H	1:H:101:MET:HE2	1.63	0.62
1:H:159:LYS:HE2	1:H:162:VAL:HG21	1.81	0.62
1:H:180:LEU:HA	1:H:199:THR:HG23	1.80	0.62
1:E:15:THR:HG23	1:E:15:THR:O	1.98	0.62
1:D:15:THR:O	1:D:15:THR:HG23	1.99	0.62
1:H:111:ARG:NH2	1:H:136:LEU:O	2.27	0.62
1:A:159:LYS:HE2	1:A:162:VAL:HG21	1.82	0.61
1:B:34:VAL:HG13	1:B:219:TRP:CZ3	2.34	0.61
1:C:94:LEU:HB3	1:C:101:MET:CE	2.30	0.61
1:D:95:ASN:H	1:D:101:MET:HE2	1.64	0.61
1:F:15:THR:HG23	1:F:15:THR:O	1.98	0.61
1:A:163:ILE:O	1:A:167:VAL:HG23	2.00	0.61
1:G:193:LYS:HD2	3:G:2043:HOH:O	1.98	0.61
1:C:159:LYS:HE2	1:C:162:VAL:HG21	1.81	0.61
1:C:15:THR:HG23	1:C:15:THR:O	1.99	0.61
1:F:94:LEU:HD22	1:F:101:MET:HE1	1.82	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:94:LEU:HD22	1:C:101:MET:HE1	1.83	0.61
1:D:180:LEU:HA	1:D:199:THR:HG23	1.81	0.61
1:G:159:LYS:HE2	1:G:162:VAL:HG21	1.82	0.61
1:B:95:ASN:H	1:B:101:MET:HE2	1.65	0.60
1:A:58:VAL:HG13	1:A:60:ILE:HG12	1.83	0.60
1:B:159:LYS:HE2	1:B:162:VAL:HG21	1.82	0.60
1:E:58:VAL:HG13	1:E:60:ILE:HG12	1.83	0.60
1:B:163:ILE:O	1:B:167:VAL:HG23	2.00	0.60
1:D:111:ARG:NH2	1:D:136:LEU:O	2.29	0.60
1:G:15:THR:HG22	1:H:73:GLY:O	2.01	0.60
1:B:94:LEU:HB3	1:B:101:MET:CE	2.31	0.60
1:A:215:GLU:C	1:A:219:TRP:HD1	2.09	0.60
1:B:111:ARG:NH2	1:B:136:LEU:O	2.29	0.60
1:C:58:VAL:HG13	1:C:60:ILE:HG12	1.83	0.60
1:D:164:THR:HG22	3:D:2041:HOH:O	2.02	0.60
1:F:94:LEU:HB3	1:F:101:MET:CE	2.32	0.60
1:B:215:GLU:C	1:B:219:TRP:HD1	2.09	0.60
1:C:96:HIS:CE1	2:C:1226:3PP:OE1	2.48	0.60
1:A:144:GLU:OE2	1:A:182:GLY:HA3	2.01	0.59
1:F:159:LYS:HE2	1:F:162:VAL:HG21	1.83	0.59
1:G:182:GLY:O	1:G:183:ALA:HB3	2.03	0.59
1:H:163:ILE:O	1:H:167:VAL:HG23	2.03	0.59
1:A:111:ARG:NH2	1:A:136:LEU:O	2.26	0.59
1:C:185:ILE:HG21	1:C:202:VAL:HG11	1.84	0.59
1:D:94:LEU:HB3	1:D:101:MET:CE	2.33	0.59
1:G:15:THR:O	1:G:15:THR:HG23	2.02	0.59
1:E:94:LEU:HB3	1:E:101:MET:CE	2.33	0.59
1:D:159:LYS:HE2	1:D:162:VAL:HG21	1.83	0.59
1:G:58:VAL:HG13	1:G:60:ILE:HG12	1.84	0.59
1:A:94:LEU:HB3	1:A:101:MET:CE	2.33	0.59
1:B:185:ILE:HG21	1:B:202:VAL:HG11	1.84	0.59
1:D:144:GLU:OE2	1:D:182:GLY:HA3	2.02	0.59
1:H:94:LEU:HB3	1:H:101:MET:CE	2.33	0.59
1:A:185:ILE:HG21	1:A:202:VAL:HG11	1.85	0.58
1:F:163:ILE:O	1:F:167:VAL:HG23	2.02	0.58
1:G:94:LEU:HD22	1:G:101:MET:HE1	1.85	0.58
1:A:204:LEU:HA	2:A:1226:3PP:HCB2	1.84	0.58
1:H:15:THR:HG23	1:H:15:THR:O	2.02	0.58
1:B:34:VAL:HG22	1:B:219:TRP:NE1	2.15	0.58
1:E:34:VAL:CG2	1:E:219:TRP:CZ2	2.80	0.58
1:F:111:ARG:NH2	1:F:136:LEU:O	2.30	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:34:VAL:O	1:G:38:THR:HB	2.03	0.58
1:A:73:GLY:O	1:B:15:THR:HG22	2.03	0.58
1:F:185:ILE:HG21	1:F:202:VAL:HG11	1.85	0.58
1:F:199:THR:CG2	1:F:201:GLY:H	2.13	0.58
1:B:34:VAL:CG2	1:B:219:TRP:CZ2	2.82	0.58
1:B:146:PRO:HA	1:B:149:ILE:CG1	2.33	0.58
1:C:146:PRO:HA	1:C:149:ILE:CG1	2.34	0.58
1:G:163:ILE:O	1:G:167:VAL:HG23	2.04	0.57
1:G:95:ASN:H	1:G:101:MET:HE2	1.69	0.57
1:B:5:LYS:C	1:B:7:PRO:CD	2.77	0.57
1:B:37:GLU:HG3	1:B:219:TRP:CZ3	2.38	0.57
1:G:146:PRO:HA	1:G:149:ILE:CG1	2.34	0.57
1:A:95:ASN:H	1:A:101:MET:HE2	1.69	0.57
1:B:34:VAL:O	1:B:38:THR:HB	2.04	0.57
1:D:163:ILE:O	1:D:167:VAL:HG23	2.05	0.57
1:B:199:THR:CG2	1:B:201:GLY:H	2.14	0.57
1:F:144:GLU:OE2	1:F:182:GLY:HA3	2.04	0.57
1:E:37:GLU:HG3	1:E:219:TRP:CZ3	2.38	0.57
1:C:52:ARG:O	1:C:56:GLU:HG3	2.05	0.57
1:F:146:PRO:HA	1:F:149:ILE:CG1	2.34	0.57
1:H:146:PRO:HA	1:H:149:ILE:CG1	2.33	0.57
1:A:182:GLY:O	1:A:183:ALA:HB3	2.04	0.57
1:C:34:VAL:O	1:C:38:THR:HB	2.04	0.57
1:G:205:ALA:N	2:G:1226:3PP:HCB1	2.20	0.57
1:C:182:GLY:O	1:C:183:ALA:HB3	2.05	0.57
1:D:146:PRO:HA	1:D:149:ILE:CG1	2.34	0.57
1:H:182:GLY:O	1:H:183:ALA:HB3	2.04	0.57
1:H:185:ILE:HG21	1:H:202:VAL:HG11	1.87	0.57
1:E:146:PRO:HA	1:E:149:ILE:CG1	2.34	0.57
1:G:204:LEU:HD21	1:G:221:LEU:HD13	1.87	0.57
1:A:34:VAL:CG2	1:A:219:TRP:CZ2	2.82	0.56
1:B:15:THR:O	1:B:15:THR:HG23	2.04	0.56
1:B:148:LEU:HD12	1:B:155:VAL:HG12	1.87	0.56
1:B:215:GLU:C	1:B:219:TRP:CD1	2.83	0.56
1:D:155:VAL:HG13	1:D:183:ALA:O	2.05	0.56
1:A:146:PRO:HA	1:A:149:ILE:CG1	2.34	0.56
1:C:144:GLU:OE2	1:C:182:GLY:HA3	2.05	0.56
2:C:1226:3PP:HCB2	3:C:2048:HOH:O	2.04	0.56
1:D:185:ILE:HG21	1:D:202:VAL:HG11	1.87	0.56
1:E:34:VAL:O	1:E:38:THR:HB	2.05	0.56
1:E:205:ALA:CB	2:E:1226:3PP:HCB1	2.34	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:154:PRO:HB2	1:F:157:LYS:HG3	1.85	0.56
1:G:148:LEU:HD12	1:G:155:VAL:HG12	1.88	0.56
1:G:199:THR:CG2	1:G:201:GLY:H	2.15	0.56
1:B:5:LYS:HB2	1:B:41:THR:HG1	1.70	0.56
1:F:34:VAL:O	1:F:38:THR:HB	2.06	0.56
1:A:34:VAL:O	1:A:38:THR:HB	2.05	0.56
1:H:58:VAL:HG11	1:H:60:ILE:HG12	1.88	0.56
1:B:182:GLY:O	1:B:183:ALA:HB3	2.06	0.56
1:D:182:GLY:O	1:D:183:ALA:HB3	2.06	0.56
1:F:199:THR:CG2	1:F:200:VAL:N	2.69	0.56
1:A:155:VAL:HG13	1:A:183:ALA:O	2.06	0.56
1:D:164:THR:HG23	3:D:2040:HOH:O	2.06	0.56
1:F:14:LYS:HE3	2:F:1226:3PP:HCG1	1.86	0.56
1:E:155:VAL:HG13	1:E:183:ALA:O	2.06	0.56
1:G:6:GLU:OE1	1:G:6:GLU:HA	2.04	0.56
1:B:144:GLU:OE2	1:B:182:GLY:HA3	2.05	0.56
1:E:215:GLU:C	1:E:219:TRP:CD1	2.81	0.56
1:G:111:ARG:NH2	1:G:136:LEU:O	2.28	0.56
1:G:185:ILE:HG21	1:G:202:VAL:HG11	1.88	0.55
1:A:204:LEU:HD21	1:A:221:LEU:HD13	1.89	0.55
1:H:155:VAL:HG13	1:H:183:ALA:O	2.06	0.55
1:F:52:ARG:O	1:F:56:GLU:HG3	2.06	0.55
1:H:148:LEU:HD12	1:H:155:VAL:HG12	1.88	0.55
1:A:205:ALA:HB3	2:A:1226:3PP:HCB1	1.88	0.55
1:C:148:LEU:HD12	1:C:155:VAL:HG12	1.89	0.55
1:D:34:VAL:O	1:D:38:THR:HB	2.06	0.55
1:D:188:GLY:O	1:D:224:GLY:HA3	2.07	0.55
1:G:204:LEU:HA	2:G:1226:3PP:HCB2	1.88	0.55
1:A:148:LEU:HD12	1:A:155:VAL:HG12	1.88	0.55
1:F:204:LEU:HD21	1:F:221:LEU:HD13	1.89	0.55
1:B:6:GLU:N	1:B:7:PRO:CD	2.70	0.55
1:E:111:ARG:NH2	1:E:136:LEU:O	2.29	0.55
1:G:144:GLU:OE2	1:G:182:GLY:HA3	2.07	0.55
1:C:111:ARG:NH2	1:C:136:LEU:O	2.28	0.54
1:E:126:ASN:ND2	1:E:126:ASN:C	2.65	0.54
1:H:204:LEU:HD21	1:H:221:LEU:HD13	1.89	0.54
1:B:155:VAL:HG13	1:B:183:ALA:O	2.07	0.54
1:G:52:ARG:O	1:G:56:GLU:HG3	2.07	0.54
1:A:215:GLU:C	1:A:219:TRP:CD1	2.83	0.54
1:B:5:LYS:C	1:B:7:PRO:HD2	2.32	0.54
1:E:204:LEU:HD21	1:E:221:LEU:HD13	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:32:GLU:HB2	1:B:58:VAL:HG13	1.90	0.54
1:C:145:PRO:HD2	1:C:155:VAL:HG11	1.88	0.54
1:D:204:LEU:HD21	1:D:221:LEU:HD13	1.90	0.54
1:G:145:PRO:HD2	1:G:155:VAL:HG11	1.88	0.54
1:H:34:VAL:O	1:H:38:THR:HB	2.07	0.54
1:B:204:LEU:HD21	1:B:221:LEU:HD13	1.90	0.54
1:D:46:PRO:CB	1:D:54:ILE:HD12	2.37	0.54
1:F:155:VAL:HG13	1:F:183:ALA:O	2.07	0.54
1:G:200:VAL:CG1	1:G:200:VAL:O	2.56	0.54
1:H:144:GLU:OE2	1:H:182:GLY:HA3	2.07	0.54
1:H:46:PRO:CB	1:H:54:ILE:HD12	2.37	0.54
1:C:204:LEU:HD21	1:C:221:LEU:HD13	1.90	0.54
1:G:111:ARG:O	1:G:115:GLU:HG3	2.08	0.54
1:H:38:THR:CG2	1:H:40:VAL:HG23	2.38	0.54
1:B:215:GLU:CG	1:B:219:TRP:NE1	2.36	0.54
1:C:163:ILE:O	1:C:167:VAL:HG23	2.07	0.54
1:B:111:ARG:O	1:B:115:GLU:HG3	2.08	0.54
1:D:200:VAL:O	1:D:200:VAL:CG1	2.56	0.54
1:E:37:GLU:CD	1:E:219:TRP:HH2	1.99	0.54
1:E:148:LEU:HD12	1:E:155:VAL:HG12	1.90	0.54
1:A:52:ARG:O	1:A:56:GLU:HG3	2.08	0.53
1:C:155:VAL:HG13	1:C:183:ALA:O	2.07	0.53
1:F:188:GLY:O	1:F:224:GLY:HA3	2.08	0.53
1:D:148:LEU:HD12	1:D:155:VAL:HG12	1.90	0.53
1:D:95:ASN:N	1:D:101:MET:HE2	2.22	0.53
1:H:200:VAL:O	1:H:200:VAL:CG1	2.57	0.53
1:B:46:PRO:CB	1:B:54:ILE:HD12	2.37	0.53
1:C:204:LEU:HA	2:C:1226:3PP:HCB2	1.89	0.53
1:F:182:GLY:O	1:F:183:ALA:HB3	2.08	0.53
1:B:129:VAL:HG23	1:D:107:GLU:HG2	1.91	0.53
1:D:107:GLU:HG3	3:D:2029:HOH:O	2.08	0.53
1:G:58:VAL:HG11	1:G:60:ILE:HG12	1.87	0.53
1:A:38:THR:CG2	1:A:40:VAL:HG23	2.38	0.53
1:D:199:THR:CG2	1:D:200:VAL:N	2.71	0.53
1:H:188:GLY:O	1:H:224:GLY:HA3	2.09	0.53
1:D:111:ARG:O	1:D:115:GLU:HG3	2.09	0.53
1:E:58:VAL:HG11	1:E:60:ILE:HG12	1.87	0.53
1:F:111:ARG:O	1:F:115:GLU:HG3	2.08	0.53
1:F:199:THR:HG22	1:F:200:VAL:N	2.22	0.53
1:G:127:PRO:HD2	3:G:2038:HOH:O	2.07	0.53
1:E:46:PRO:CB	1:E:54:ILE:HD12	2.39	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:182:GLY:O	1:E:183:ALA:HB3	2.09	0.53
1:F:46:PRO:CB	1:F:54:ILE:HD12	2.38	0.53
1:C:46:PRO:CB	1:C:54:ILE:HD12	2.39	0.52
1:D:38:THR:CG2	1:D:40:VAL:HG23	2.39	0.52
1:D:199:THR:CG2	1:D:201:GLY:H	2.14	0.52
1:E:199:THR:CG2	1:E:200:VAL:N	2.72	0.52
1:B:34:VAL:HA	1:B:219:TRP:CZ3	2.37	0.52
1:E:95:ASN:N	1:E:101:MET:HE2	2.25	0.52
1:F:38:THR:CG2	1:F:40:VAL:HG23	2.39	0.52
1:F:95:ASN:N	1:F:101:MET:HE2	2.23	0.52
1:F:107:GLU:HG3	3:F:2028:HOH:O	2.09	0.52
1:G:126:ASN:ND2	1:G:126:ASN:C	2.67	0.52
1:B:188:GLY:O	1:B:224:GLY:HA3	2.09	0.52
1:E:52:ARG:O	1:E:56:GLU:HG3	2.10	0.52
1:F:148:LEU:HD12	1:F:155:VAL:HG12	1.92	0.52
1:G:107:GLU:O	1:G:111:ARG:HG2	2.09	0.52
1:G:155:VAL:HG13	1:G:183:ALA:O	2.09	0.52
1:A:46:PRO:CB	1:A:54:ILE:HD12	2.39	0.52
1:B:38:THR:CG2	1:B:40:VAL:HG23	2.40	0.52
1:B:199:THR:CG2	1:B:200:VAL:N	2.73	0.52
1:C:188:GLY:O	1:C:224:GLY:HA3	2.10	0.52
1:D:52:ARG:O	1:D:56:GLU:HG3	2.09	0.52
1:D:126:ASN:ND2	1:D:126:ASN:C	2.64	0.52
1:C:199:THR:CG2	1:C:200:VAL:N	2.72	0.52
1:E:34:VAL:HG22	1:E:219:TRP:NE1	2.15	0.52
1:B:107:GLU:O	1:B:111:ARG:HG2	2.10	0.52
1:C:111:ARG:O	1:C:115:GLU:HG3	2.10	0.52
1:A:215:GLU:CG	1:A:219:TRP:NE1	2.36	0.51
1:G:46:PRO:CB	1:G:54:ILE:HD12	2.38	0.51
1:H:199:THR:CG2	1:H:200:VAL:N	2.73	0.51
1:C:107:GLU:O	1:C:111:ARG:HG2	2.10	0.51
1:C:161:GLU:CD	1:C:161:GLU:H	2.17	0.51
1:F:149:ILE:O	2:F:1226:3PP:O2	2.28	0.51
1:H:107:GLU:O	1:H:111:ARG:HG2	2.10	0.51
1:B:58:VAL:HG11	1:B:60:ILE:HG12	1.93	0.51
1:B:200:VAL:O	1:B:200:VAL:CG1	2.58	0.51
1:F:161:GLU:H	1:F:161:GLU:CD	2.18	0.51
1:F:200:VAL:O	1:F:200:VAL:CG1	2.58	0.51
1:G:161:GLU:CD	1:G:161:GLU:H	2.17	0.51
1:H:95:ASN:N	1:H:101:MET:HE2	2.25	0.51
1:C:58:VAL:HG11	1:C:60:ILE:HG12	1.89	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:188:GLY:O	1:G:224:GLY:HA3	2.11	0.51
1:H:111:ARG:O	1:H:115:GLU:HG3	2.11	0.51
1:H:126:ASN:ND2	1:H:126:ASN:C	2.65	0.51
1:H:199:THR:HG22	1:H:200:VAL:N	2.25	0.51
1:B:52:ARG:O	1:B:56:GLU:HG3	2.10	0.51
1:D:199:THR:HG22	1:D:200:VAL:N	2.24	0.51
1:E:199:THR:CG2	1:E:201:GLY:H	2.15	0.51
1:A:107:GLU:O	1:A:111:ARG:HG2	2.11	0.51
1:A:199:THR:CG2	1:A:201:GLY:H	2.14	0.51
1:B:161:GLU:CD	1:B:161:GLU:H	2.18	0.51
1:C:38:THR:CG2	1:C:40:VAL:HG23	2.41	0.51
1:F:107:GLU:O	1:F:111:ARG:HG2	2.11	0.51
1:A:34:VAL:HG22	1:A:219:TRP:NE1	2.16	0.51
1:A:58:VAL:HG11	1:A:60:ILE:HG12	1.90	0.51
1:A:188:GLY:O	1:A:224:GLY:HA3	2.11	0.51
1:D:107:GLU:O	1:D:111:ARG:HG2	2.10	0.51
1:E:111:ARG:O	1:E:115:GLU:HG3	2.10	0.51
1:A:199:THR:CG2	1:A:200:VAL:N	2.74	0.51
1:A:200:VAL:O	1:A:200:VAL:CG1	2.58	0.51
1:B:206:SER:HA	1:B:209:THR:HG22	1.93	0.51
1:G:16:TYR:CD2	1:G:209:THR:HG21	2.46	0.51
1:G:199:THR:CG2	1:G:200:VAL:N	2.74	0.51
1:A:206:SER:HA	1:A:209:THR:HG22	1.92	0.51
1:C:206:SER:HA	1:C:209:THR:HG22	1.93	0.51
1:E:34:VAL:HA	1:E:219:TRP:CZ3	2.40	0.51
1:F:4:LEU:HD21	1:F:119:MET:HE1	1.93	0.51
1:F:145:PRO:HD2	1:F:155:VAL:HG11	1.91	0.51
1:H:6:GLU:HA	1:H:6:GLU:OE1	2.11	0.51
1:H:199:THR:CG2	1:H:201:GLY:H	2.16	0.51
1:B:95:ASN:N	1:B:101:MET:HE2	2.25	0.50
1:E:188:GLY:O	1:E:224:GLY:HA3	2.11	0.50
1:G:38:THR:CG2	1:G:40:VAL:HG23	2.41	0.50
1:A:111:ARG:O	1:A:115:GLU:HG3	2.11	0.50
1:F:16:TYR:CD2	1:F:209:THR:HG21	2.46	0.50
1:B:16:TYR:CD2	1:B:209:THR:HG21	2.46	0.50
1:C:200:VAL:O	1:C:200:VAL:CG1	2.59	0.50
1:E:200:VAL:CG1	1:E:200:VAL:O	2.58	0.50
1:G:206:SER:HA	1:G:209:THR:HG22	1.93	0.50
1:D:145:PRO:HD2	1:D:155:VAL:HG11	1.93	0.50
1:H:161:GLU:CD	1:H:161:GLU:H	2.19	0.50
1:A:145:PRO:HD2	1:A:155:VAL:HG11	1.92	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:6:GLU:N	1:B:7:PRO:HD3	2.27	0.50
1:B:145:PRO:HD2	1:B:155:VAL:HG11	1.92	0.50
1:D:161:GLU:H	1:D:161:GLU:CD	2.18	0.50
1:D:206:SER:HA	1:D:209:THR:HG22	1.94	0.50
1:E:145:PRO:HD2	1:E:155:VAL:HG11	1.92	0.50
1:G:106:LEU:O	1:G:110:ILE:HG13	2.11	0.50
1:H:33:LYS:HD3	1:H:215:GLU:CD	2.37	0.50
1:E:215:GLU:CG	1:E:219:TRP:NE1	2.37	0.50
1:A:16:TYR:CD2	1:A:209:THR:HG21	2.47	0.50
1:B:106:LEU:O	1:B:110:ILE:HG13	2.10	0.50
1:D:6:GLU:OE1	1:D:6:GLU:HA	2.12	0.50
1:E:4:LEU:HD21	1:E:119:MET:HE1	1.94	0.50
1:B:215:GLU:CD	1:B:219:TRP:HE1	2.17	0.49
1:E:199:THR:HG22	1:E:200:VAL:N	2.25	0.49
1:E:206:SER:HA	1:E:209:THR:HG22	1.93	0.49
1:C:199:THR:HG22	1:C:200:VAL:N	2.26	0.49
1:E:161:GLU:CD	1:E:161:GLU:H	2.19	0.49
1:A:34:VAL:HA	1:A:219:TRP:CZ3	2.41	0.49
1:F:206:SER:HA	1:F:209:THR:HG22	1.95	0.49
1:G:182:GLY:O	1:G:183:ALA:CB	2.59	0.49
1:B:199:THR:HG22	1:B:200:VAL:N	2.26	0.49
1:D:137:ASN:N	1:D:138:PRO:HD3	2.26	0.49
1:H:137:ASN:N	1:H:138:PRO:HD3	2.26	0.49
1:A:6:GLU:HB2	1:A:7:PRO:HD3	1.95	0.49
1:C:4:LEU:HD21	1:C:119:MET:HE1	1.93	0.49
1:E:16:TYR:CD2	1:E:209:THR:HG21	2.48	0.49
1:H:145:PRO:HD2	1:H:155:VAL:HG11	1.94	0.49
1:A:34:VAL:CG2	1:A:219:TRP:CD2	2.88	0.49
1:F:58:VAL:HG11	1:F:60:ILE:HG12	1.94	0.49
1:F:129:VAL:HG23	1:H:107:GLU:HG2	1.95	0.49
1:B:199:THR:HG21	1:B:201:GLY:O	2.13	0.49
1:D:6:GLU:HB2	1:D:7:PRO:HD3	1.95	0.49
1:D:159:LYS:N	1:D:160:PRO:CD	2.76	0.49
1:F:137:ASN:N	1:F:138:PRO:HD3	2.26	0.49
1:H:159:LYS:N	1:H:160:PRO:CD	2.76	0.49
1:H:204:LEU:C	2:H:1226:3PP:HCB1	2.38	0.49
1:B:4:LEU:HD21	1:B:119:MET:HE1	1.95	0.48
1:B:33:LYS:HD3	1:B:215:GLU:CD	2.38	0.48
1:D:16:TYR:CD2	1:D:209:THR:HG21	2.47	0.48
1:A:4:LEU:HD21	1:A:119:MET:HE1	1.95	0.48
1:C:16:TYR:CD2	1:C:209:THR:HG21	2.48	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:95:ASN:N	1:C:101:MET:HE2	2.27	0.48
1:E:34:VAL:CG2	1:E:219:TRP:CD2	2.88	0.48
1:E:107:GLU:O	1:E:111:ARG:HG2	2.13	0.48
1:D:4:LEU:HD21	1:D:119:MET:HE1	1.94	0.48
1:E:199:THR:HG21	1:E:201:GLY:O	2.12	0.48
1:G:75:HIS:CD2	1:H:15:THR:HG21	2.49	0.48
1:A:6:GLU:HA	1:A:6:GLU:OE1	2.14	0.48
1:E:6:GLU:OE1	1:E:6:GLU:HA	2.13	0.48
1:H:182:GLY:O	1:H:183:ALA:CB	2.61	0.48
1:A:33:LYS:HD3	1:A:215:GLU:CD	2.38	0.48
1:A:137:ASN:N	1:A:138:PRO:HD3	2.27	0.48
1:B:182:GLY:O	1:B:183:ALA:CB	2.62	0.48
1:D:33:LYS:HD3	1:D:215:GLU:CD	2.38	0.48
1:D:209:THR:CG2	1:D:210:LYS:HG2	2.42	0.48
1:E:38:THR:CG2	1:E:40:VAL:HG23	2.43	0.48
1:E:96:HIS:HE1	2:E:1226:3PP:OE1	1.95	0.48
1:F:33:LYS:HD3	1:F:215:GLU:CD	2.39	0.48
1:B:137:ASN:N	1:B:138:PRO:HD3	2.27	0.48
1:C:6:GLU:HB2	1:C:7:PRO:HD3	1.96	0.48
1:E:94:LEU:HB3	1:E:101:MET:HE1	1.96	0.48
1:E:205:ALA:N	2:E:1226:3PP:CB	2.73	0.48
1:G:199:THR:HG22	1:G:200:VAL:N	2.27	0.48
1:G:199:THR:HG21	1:G:201:GLY:O	2.13	0.48
1:H:106:LEU:O	1:H:110:ILE:HG13	2.14	0.48
1:H:199:THR:HG21	1:H:201:GLY:O	2.14	0.48
1:H:206:SER:HA	1:H:209:THR:HG22	1.96	0.48
1:E:6:GLU:HB2	1:E:7:PRO:HD3	1.96	0.48
1:E:33:LYS:HD3	1:E:215:GLU:CD	2.38	0.48
1:G:33:LYS:HD3	1:G:215:GLU:CD	2.39	0.48
1:H:209:THR:CG2	1:H:210:LYS:HG2	2.43	0.48
1:B:149:ILE:O	2:B:1226:3PP:O2	2.32	0.47
1:C:106:LEU:O	1:C:110:ILE:HG13	2.13	0.47
1:C:199:THR:HG21	1:C:201:GLY:O	2.13	0.47
1:F:6:GLU:HB2	1:F:7:PRO:HD3	1.96	0.47
1:F:106:LEU:O	1:F:110:ILE:HG13	2.13	0.47
1:A:129:VAL:HG23	1:C:107:GLU:HG2	1.97	0.47
1:H:6:GLU:HB2	1:H:7:PRO:HD3	1.96	0.47
1:F:150:GLY:N	3:F:2039:HOH:O	2.38	0.47
1:G:168:GLU:HG3	1:G:169:LEU:N	2.29	0.47
1:H:16:TYR:CD2	1:H:209:THR:HG21	2.48	0.47
1:A:32:GLU:HG3	1:A:59:GLU:HG2	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:149:ILE:O	2:A:1226:3PP:O2	2.33	0.47
1:C:182:GLY:O	1:C:183:ALA:CB	2.61	0.47
1:G:137:ASN:N	1:G:138:PRO:HD3	2.29	0.47
1:G:168:GLU:CG	1:G:169:LEU:N	2.78	0.47
1:A:159:LYS:N	1:A:160:PRO:CD	2.75	0.47
1:A:182:GLY:O	1:A:183:ALA:CB	2.62	0.47
1:C:6:GLU:HA	1:C:6:GLU:OE1	2.14	0.47
1:D:155:VAL:HG11	1:D:182:GLY:O	2.14	0.47
1:E:137:ASN:N	1:E:138:PRO:HD3	2.28	0.47
1:E:159:LYS:N	1:E:160:PRO:CD	2.75	0.47
1:F:6:GLU:HA	1:F:6:GLU:OE1	2.14	0.47
1:A:199:THR:HG22	1:A:200:VAL:N	2.28	0.47
1:B:94:LEU:HD22	1:B:101:MET:CE	2.44	0.47
1:D:58:VAL:HG11	1:D:60:ILE:HG12	1.91	0.47
1:F:159:LYS:N	1:F:160:PRO:CD	2.75	0.47
1:B:159:LYS:N	1:B:160:PRO:CD	2.76	0.47
1:C:3:LYS:HE2	3:C:2047:HOH:O	2.15	0.47
1:D:199:THR:HG21	1:D:201:GLY:O	2.14	0.47
1:E:12:ASN:HB3	2:E:1226:3PP:OE2	2.15	0.47
1:H:94:LEU:HD22	1:H:101:MET:CE	2.44	0.47
1:A:95:ASN:N	1:A:101:MET:HE2	2.30	0.46
1:C:137:ASN:N	1:C:138:PRO:HD3	2.30	0.46
1:H:52:ARG:O	1:H:56:GLU:HG3	2.15	0.46
1:G:96:HIS:CE1	2:G:1226:3PP:OE1	2.61	0.46
1:C:33:LYS:HD3	1:C:215:GLU:CD	2.40	0.46
1:C:76:THR:O	1:C:78:HIS:HD2	1.99	0.46
1:C:199:THR:CG2	1:C:201:GLY:H	2.14	0.46
1:D:14:LYS:HE3	2:D:1226:3PP:HCG1	1.97	0.46
1:D:164:THR:O	1:D:168:GLU:HG3	2.16	0.46
1:E:36:LYS:HE2	3:E:2011:HOH:O	2.15	0.46
1:E:149:ILE:O	2:E:1226:3PP:O2	2.34	0.46
1:F:157:LYS:HE3	1:F:190:ASP:OD1	2.15	0.46
1:G:38:THR:HG21	1:G:222:VAL:CG1	2.44	0.46
1:H:4:LEU:HD21	1:H:119:MET:HE1	1.96	0.46
1:A:76:THR:O	1:A:78:HIS:HD2	1.97	0.46
1:A:199:THR:HG21	1:A:201:GLY:O	2.15	0.46
1:C:38:THR:HG21	1:C:222:VAL:CG1	2.43	0.46
1:E:76:THR:O	1:E:78:HIS:HD2	1.98	0.46
1:E:106:LEU:O	1:E:110:ILE:HG13	2.14	0.46
1:A:209:THR:CG2	1:A:210:LYS:HG2	2.44	0.46
1:C:94:LEU:HB3	1:C:101:MET:HE1	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:199:THR:HG21	1:F:201:GLY:O	2.16	0.46
1:A:126:ASN:ND2	1:A:126:ASN:C	2.66	0.46
1:A:205:ALA:N	2:A:1226:3PP:CB	2.78	0.46
1:B:33:LYS:O	1:B:37:GLU:HG2	2.16	0.46
1:G:165:ASN:C	1:G:168:GLU:HG2	2.39	0.46
1:A:94:LEU:HD22	1:A:101:MET:CE	2.45	0.46
1:E:129:VAL:HG23	1:G:107:GLU:HG2	1.98	0.46
1:G:6:GLU:O	1:G:200:VAL:HG13	2.16	0.46
1:A:136:LEU:HD21	1:C:132:ALA:HB2	1.97	0.46
1:A:155:VAL:HG11	1:A:182:GLY:O	2.15	0.46
1:A:215:GLU:CD	1:A:219:TRP:HE1	2.17	0.45
1:B:94:LEU:HB3	1:B:101:MET:HE1	1.99	0.45
1:C:159:LYS:N	1:C:160:PRO:CD	2.78	0.45
1:E:94:LEU:HD22	1:E:101:MET:CE	2.46	0.45
1:G:165:ASN:HA	1:G:168:GLU:CD	2.41	0.45
1:A:122:VAL:O	1:A:141:VAL:HA	2.16	0.45
1:B:71:LYS:HA	1:B:112:ARG:HH21	1.82	0.45
1:D:182:GLY:O	1:D:183:ALA:CB	2.64	0.45
1:F:122:VAL:O	1:F:141:VAL:HA	2.16	0.45
1:A:144:GLU:HG2	1:A:182:GLY:CA	2.47	0.45
1:E:209:THR:CG2	1:E:210:LYS:HG2	2.45	0.45
1:G:159:LYS:N	1:G:160:PRO:CD	2.79	0.45
1:G:95:ASN:N	1:G:101:MET:HE2	2.30	0.45
1:C:45:ALA:HA	1:C:63:PHE:O	2.16	0.45
1:F:155:VAL:HG11	1:F:182:GLY:O	2.17	0.45
1:B:14:LYS:NZ	2:B:1226:3PP:HCG1	2.32	0.45
1:C:33:LYS:O	1:C:37:GLU:HG2	2.16	0.45
1:D:204:LEU:HA	2:D:1226:3PP:HCB1	1.99	0.45
1:E:122:VAL:O	1:E:141:VAL:HA	2.17	0.45
1:E:164:THR:O	1:E:168:GLU:HG3	2.17	0.45
1:F:71:LYS:HA	1:F:112:ARG:HH21	1.82	0.45
1:F:164:THR:O	1:F:168:GLU:HG3	2.16	0.45
1:G:149:ILE:O	2:G:1226:3PP:O2	2.34	0.45
1:H:155:VAL:CG1	1:H:183:ALA:HB3	2.47	0.45
1:G:4:LEU:HD21	1:G:119:MET:HE1	1.98	0.45
1:D:21:GLY:O	1:D:24:ALA:HB3	2.17	0.45
1:E:4:LEU:CD2	1:E:119:MET:HE1	2.47	0.45
1:E:107:GLU:HG2	1:G:129:VAL:CG2	2.44	0.45
1:G:94:LEU:HB3	1:G:101:MET:HE1	1.98	0.45
1:H:164:THR:O	1:H:168:GLU:HG3	2.17	0.45
1:D:106:LEU:O	1:D:110:ILE:HG13	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:33:LYS:O	1:G:37:GLU:HG2	2.17	0.44
1:G:76:THR:O	1:G:78:HIS:HD2	2.00	0.44
1:H:94:LEU:HB3	1:H:101:MET:HE1	1.98	0.44
1:E:215:GLU:CD	1:E:219:TRP:HE1	2.19	0.44
1:G:155:VAL:CG1	1:G:183:ALA:HB3	2.46	0.44
1:H:205:ALA:H	2:H:1226:3PP:HCB1	1.78	0.44
1:A:164:THR:O	1:A:168:GLU:HG3	2.17	0.44
1:B:8:ILE:H	1:B:8:ILE:HG13	1.42	0.44
1:D:218:ILE:O	1:D:222:VAL:HG23	2.17	0.44
1:F:182:GLY:O	1:F:183:ALA:CB	2.65	0.44
1:G:122:VAL:O	1:G:141:VAL:HA	2.16	0.44
1:B:96:HIS:HE1	2:B:1226:3PP:OE2	2.00	0.44
1:C:126:ASN:ND2	1:C:126:ASN:C	2.66	0.44
1:D:155:VAL:CG1	1:D:183:ALA:HB3	2.47	0.44
1:F:94:LEU:HB3	1:F:101:MET:HE1	2.00	0.44
1:G:199:THR:CG2	1:G:201:GLY:O	2.65	0.44
1:H:155:VAL:HG11	1:H:182:GLY:O	2.17	0.44
1:A:96:HIS:CE1	2:A:1226:3PP:OE1	2.59	0.44
1:A:107:GLU:HG2	1:C:129:VAL:CG2	2.45	0.44
1:D:159:LYS:HA	1:D:161:GLU:OE2	2.18	0.44
1:E:33:LYS:O	1:E:37:GLU:HG2	2.18	0.44
1:G:159:LYS:HA	1:G:161:GLU:OE2	2.18	0.44
3:G:2026:HOH:O	1:H:22:LYS:HD2	2.17	0.44
1:D:94:LEU:HD22	1:D:101:MET:CE	2.46	0.44
1:E:182:GLY:O	1:E:183:ALA:CB	2.66	0.44
1:E:199:THR:CG2	1:E:201:GLY:O	2.66	0.44
1:H:40:VAL:HG21	1:H:222:VAL:HG13	2.00	0.44
1:A:4:LEU:CD2	1:A:119:MET:HE1	2.48	0.44
1:B:155:VAL:CG1	1:B:183:ALA:HB3	2.47	0.44
1:E:218:ILE:O	1:E:222:VAL:HG23	2.16	0.44
1:G:165:ASN:HA	1:G:168:GLU:OE2	2.18	0.44
1:H:159:LYS:HA	1:H:161:GLU:OE2	2.18	0.44
1:C:73:GLY:C	1:D:15:THR:HG22	2.43	0.44
1:A:94:LEU:HB3	1:A:101:MET:HE1	2.00	0.43
1:E:45:ALA:HA	1:E:63:PHE:O	2.17	0.43
1:E:155:VAL:CG1	1:E:183:ALA:HB3	2.48	0.43
1:G:71:LYS:HA	1:G:112:ARG:HH21	1.82	0.43
1:B:155:VAL:HG11	1:B:182:GLY:O	2.18	0.43
1:F:45:ALA:HA	1:F:63:PHE:O	2.18	0.43
1:A:199:THR:CG2	1:A:201:GLY:O	2.67	0.43
1:D:94:LEU:HB3	1:D:101:MET:HE1	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:71:LYS:HA	1:E:112:ARG:HH21	1.82	0.43
1:A:121:MET:HE2	1:A:180:LEU:HD11	2.00	0.43
1:C:199:THR:CG2	1:C:201:GLY:O	2.66	0.43
1:D:4:LEU:CD2	1:D:119:MET:HE1	2.48	0.43
1:G:22:LYS:O	1:G:26:GLU:HG3	2.18	0.43
1:B:164:THR:O	1:B:168:GLU:HG3	2.18	0.43
1:C:155:VAL:CG1	1:C:183:ALA:HB3	2.47	0.43
1:D:38:THR:HG21	1:D:222:VAL:CG1	2.46	0.43
1:B:34:VAL:CA	1:B:219:TRP:CZ2	2.99	0.43
1:C:71:LYS:HA	1:C:112:ARG:HH21	1.83	0.43
1:E:102:ILE:O	1:E:103:LEU:C	2.61	0.43
1:G:144:GLU:HG2	1:G:182:GLY:CA	2.49	0.43
1:B:199:THR:CG2	1:B:201:GLY:O	2.67	0.43
1:C:94:LEU:HD22	1:C:101:MET:CE	2.47	0.43
1:D:40:VAL:HG21	1:D:222:VAL:HG13	2.01	0.43
1:E:47:GLN:HA	3:E:2014:HOH:O	2.19	0.43
1:F:159:LYS:HA	1:F:161:GLU:OE2	2.18	0.43
1:A:155:VAL:CG1	1:A:183:ALA:HB3	2.49	0.43
1:B:76:THR:O	1:B:78:HIS:HD2	2.01	0.43
1:B:159:LYS:HA	1:B:161:GLU:OE2	2.18	0.43
1:B:180:LEU:HD22	1:B:201:GLY:CA	2.49	0.43
1:D:144:GLU:HG2	1:D:182:GLY:CA	2.48	0.43
1:F:4:LEU:CD2	1:F:119:MET:HE1	2.48	0.43
1:B:121:MET:HE2	1:B:180:LEU:HD11	2.01	0.43
1:D:122:VAL:O	1:D:141:VAL:HA	2.19	0.43
1:F:94:LEU:HD22	1:F:101:MET:CE	2.48	0.43
1:G:45:ALA:HA	1:G:63:PHE:O	2.18	0.43
1:G:180:LEU:HD22	1:G:201:GLY:CA	2.49	0.43
1:H:199:THR:CG2	1:H:201:GLY:O	2.67	0.43
1:C:149:ILE:O	2:C:1226:3PP:O2	2.36	0.42
1:E:144:GLU:HG2	1:E:182:GLY:CA	2.49	0.42
1:E:155:VAL:HG11	1:E:182:GLY:O	2.19	0.42
1:F:218:ILE:O	1:F:222:VAL:HG23	2.19	0.42
1:G:94:LEU:HD22	1:G:101:MET:CE	2.48	0.42
1:H:33:LYS:O	1:H:37:GLU:HG2	2.18	0.42
1:H:45:ALA:HA	1:H:63:PHE:O	2.17	0.42
1:A:45:ALA:HA	1:A:63:PHE:O	2.19	0.42
1:C:122:VAL:O	1:C:141:VAL:HA	2.18	0.42
1:D:76:THR:O	1:D:78:HIS:HD2	2.01	0.42
1:D:187:THR:CG2	1:D:189:GLU:H	2.09	0.42
1:E:136:LEU:HD21	1:G:132:ALA:HB2	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:159:LYS:HA	1:E:161:GLU:OE2	2.19	0.42
1:F:155:VAL:CG1	1:F:183:ALA:HB3	2.49	0.42
1:F:204:LEU:C	2:F:1226:3PP:HCB1	2.45	0.42
1:H:4:LEU:CD2	1:H:119:MET:HE1	2.49	0.42
1:B:204:LEU:C	2:B:1226:3PP:HCB1	2.43	0.42
1:C:4:LEU:CD2	1:C:119:MET:HE1	2.49	0.42
1:D:126:ASN:HD21	1:D:129:VAL:HG23	1.84	0.42
1:A:22:LYS:O	1:A:26:GLU:HG3	2.19	0.42
1:A:106:LEU:O	1:A:110:ILE:HG13	2.18	0.42
1:E:121:MET:HE2	1:E:180:LEU:HD11	2.01	0.42
1:F:102:ILE:HA	1:H:102:ILE:HA	2.01	0.42
1:G:75:HIS:HD2	1:H:15:THR:HG21	1.83	0.42
1:A:136:LEU:CD2	1:C:132:ALA:HB2	2.50	0.42
1:B:122:VAL:O	1:B:141:VAL:HA	2.19	0.42
1:C:75:HIS:CD2	1:D:15:THR:HG21	2.54	0.42
1:C:144:GLU:HG2	1:C:182:GLY:CA	2.49	0.42
1:C:164:THR:O	1:C:168:GLU:HG3	2.20	0.42
1:F:144:GLU:HG2	1:F:182:GLY:CA	2.49	0.42
1:H:108:ALA:O	1:H:112:ARG:HB2	2.19	0.42
1:C:15:THR:HG21	1:D:75:HIS:HD2	1.84	0.42
1:C:218:ILE:O	1:C:222:VAL:HG23	2.20	0.42
1:H:38:THR:HG21	1:H:222:VAL:CG1	2.47	0.42
1:H:38:THR:HG22	1:H:40:VAL:N	2.21	0.42
1:B:218:ILE:O	1:B:222:VAL:HG23	2.20	0.42
1:D:45:ALA:HA	1:D:63:PHE:O	2.18	0.42
1:E:22:LYS:O	1:E:26:GLU:HG3	2.19	0.42
1:F:76:THR:O	1:F:78:HIS:HD2	2.02	0.42
1:C:15:THR:O	1:C:15:THR:CG2	2.68	0.42
1:B:21:GLY:O	1:B:24:ALA:HB3	2.20	0.42
1:E:204:LEU:C	1:E:204:LEU:HD12	2.45	0.42
1:G:157:LYS:HG3	3:G:2043:HOH:O	2.20	0.42
1:A:102:ILE:O	1:A:103:LEU:C	2.62	0.41
1:A:204:LEU:C	1:A:204:LEU:HD12	2.45	0.41
1:B:199:THR:HG22	1:B:201:GLY:N	2.17	0.41
1:C:159:LYS:HA	1:C:161:GLU:OE2	2.20	0.41
1:H:141:VAL:O	1:H:141:VAL:HG13	2.20	0.41
1:B:144:GLU:HG2	1:B:182:GLY:CA	2.50	0.41
1:G:205:ALA:CB	2:G:1226:3PP:HCB1	2.49	0.41
1:B:108:ALA:O	1:B:112:ARG:HB2	2.21	0.41
1:C:6:GLU:HG2	3:C:2001:HOH:O	2.20	0.41
1:C:22:LYS:O	1:C:26:GLU:HG3	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:204:LEU:C	1:C:204:LEU:HD12	2.46	0.41
1:D:199:THR:CG2	1:D:201:GLY:O	2.68	0.41
1:D:33:LYS:O	1:D:37:GLU:HG2	2.20	0.41
1:D:100:ARG:HD2	3:D:2025:HOH:O	2.20	0.41
1:H:143:VAL:HG11	3:H:2038:HOH:O	2.19	0.41
1:B:45:ALA:HA	1:B:63:PHE:O	2.21	0.41
1:C:141:VAL:O	1:C:141:VAL:HG13	2.20	0.41
1:D:204:LEU:HB2	3:D:2044:HOH:O	2.21	0.41
1:D:205:ALA:H	2:D:1226:3PP:HCB1	1.84	0.41
1:E:38:THR:HG22	1:E:40:VAL:N	2.21	0.41
1:A:71:LYS:HA	1:A:112:ARG:HH21	1.84	0.41
1:C:143:VAL:HG22	1:C:166:THR:HG21	2.02	0.41
1:D:71:LYS:HA	1:D:112:ARG:HH21	1.85	0.41
1:F:119:MET:HE2	3:F:2024:HOH:O	2.20	0.41
1:G:111:ARG:HA	1:G:111:ARG:HD3	1.84	0.41
1:G:149:ILE:O	1:G:149:ILE:CG2	2.68	0.41
1:G:155:VAL:HG11	1:G:182:GLY:O	2.20	0.41
1:B:4:LEU:CD2	1:B:119:MET:HE1	2.50	0.41
1:B:6:GLU:OE1	1:B:6:GLU:HA	2.20	0.41
1:B:14:LYS:HE3	2:B:1226:3PP:HCG1	2.03	0.41
1:G:149:ILE:O	1:G:149:ILE:HG22	2.20	0.41
1:C:102:ILE:O	1:C:103:LEU:C	2.63	0.41
1:H:31:ALA:HB1	1:H:42:ILE:HG21	2.03	0.41
1:B:126:ASN:ND2	1:B:126:ASN:C	2.64	0.41
1:B:143:VAL:CG1	1:B:166:THR:HG21	2.46	0.41
1:C:31:ALA:HB1	1:C:42:ILE:HG21	2.03	0.41
1:C:155:VAL:HG11	1:C:182:GLY:O	2.21	0.41
1:B:163:ILE:O	1:B:166:THR:OG1	2.34	0.41
1:F:33:LYS:O	1:F:37:GLU:HG2	2.21	0.41
1:F:40:VAL:HG21	1:F:222:VAL:HG13	2.02	0.41
1:G:21:GLY:O	1:G:24:ALA:HB3	2.21	0.41
1:B:209:THR:CG2	1:B:210:LYS:HG2	2.43	0.40
1:D:111:ARG:HA	1:D:111:ARG:HD3	1.82	0.40
1:E:34:VAL:CA	1:E:219:TRP:CZ2	2.98	0.40
1:F:94:LEU:HB3	1:F:101:MET:HE2	2.03	0.40
1:H:21:GLY:O	1:H:24:ALA:HB3	2.21	0.40
1:C:15:THR:HG21	1:D:75:HIS:CD2	2.57	0.40
1:D:141:VAL:O	1:D:141:VAL:HG13	2.20	0.40
1:F:141:VAL:HG13	1:F:141:VAL:O	2.21	0.40
1:G:209:THR:CG2	1:G:210:LYS:HG2	2.46	0.40
1:H:218:ILE:O	1:H:222:VAL:HG23	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:6:GLU:CB	1:A:7:PRO:HD3	2.51	0.40
1:C:40:VAL:HG21	1:C:222:VAL:HG13	2.02	0.40
1:F:6:GLU:CB	1:F:7:PRO:HD3	2.51	0.40
1:F:14:LYS:CE	2:F:1226:3PP:HCG1	2.49	0.40
1:G:218:ILE:O	1:G:222:VAL:HG23	2.22	0.40
1:H:162:VAL:CG1	3:H:2038:HOH:O	2.62	0.40
1:A:102:ILE:HA	1:C:102:ILE:HA	2.03	0.40
1:C:155:VAL:O	1:C:158:ALA:O	2.39	0.40
1:F:126:ASN:ND2	1:F:126:ASN:C	2.66	0.40
1:B:31:ALA:HB1	1:B:42:ILE:HG21	2.02	0.40
1:B:155:VAL:O	1:B:158:ALA:O	2.39	0.40
1:C:112:ARG:HD2	1:C:112:ARG:HA	1.86	0.40
1:H:14:LYS:HE3	2:H:1226:3PP:HCG1	2.02	0.40
1:H:71:LYS:HA	1:H:112:ARG:HH21	1.86	0.40

All (4) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:37:GLU:OE2	1:C:168:GLU:OE1[2_646]	1.73	0.47
1:E:37:GLU:CB	1:G:165:ASN:OD1[2_747]	1.92	0.28
1:F:137:ASN:OD1	1:G:212:LYS:NZ[2_647]	1.99	0.21
1:F:137:ASN:OD1	1:G:212:LYS:CE[2_647]	2.18	0.02

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	222/225 (99%)	205 (92%)	14 (6%)	3 (1%)	9	23
1	B	222/225 (99%)	206 (93%)	14 (6%)	2 (1%)	14	35
1	C	222/225 (99%)	206 (93%)	13 (6%)	3 (1%)	9	23

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	222/225 (99%)	204 (92%)	15 (7%)	3 (1%)	9	23
1	E	222/225 (99%)	206 (93%)	13 (6%)	3 (1%)	9	23
1	F	222/225 (99%)	206 (93%)	13 (6%)	3 (1%)	9	23
1	G	222/225 (99%)	207 (93%)	12 (5%)	3 (1%)	9	23
1	H	222/225 (99%)	206 (93%)	13 (6%)	3 (1%)	9	23
All	All	1776/1800 (99%)	1646 (93%)	107 (6%)	23 (1%)	9	25

All (23) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	6	GLU
1	C	6	GLU
1	D	6	GLU
1	E	6	GLU
1	F	6	GLU
1	H	6	GLU
1	A	149	ILE
1	B	149	ILE
1	C	149	ILE
1	D	149	ILE
1	E	149	ILE
1	F	149	ILE
1	G	149	ILE
1	H	149	ILE
1	A	183	ALA
1	B	183	ALA
1	C	183	ALA
1	D	183	ALA
1	E	183	ALA
1	F	183	ALA
1	G	183	ALA
1	H	183	ALA
1	G	6	GLU

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	179/180 (99%)	166 (93%)	13 (7%)	13	32
1	B	179/180 (99%)	166 (93%)	13 (7%)	13	32
1	C	179/180 (99%)	165 (92%)	14 (8%)	11	29
1	D	179/180 (99%)	165 (92%)	14 (8%)	11	29
1	E	179/180 (99%)	165 (92%)	14 (8%)	11	29
1	F	179/180 (99%)	165 (92%)	14 (8%)	11	29
1	G	179/180 (99%)	164 (92%)	15 (8%)	10	26
1	H	179/180 (99%)	165 (92%)	14 (8%)	11	29
All	All	1432/1440 (99%)	1321 (92%)	111 (8%)	11	29

All (111) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	15	THR
1	A	38	THR
1	A	44	VAL
1	A	58	VAL
1	A	82	GLU
1	A	106	LEU
1	A	107	GLU
1	A	116	VAL
1	A	126	ASN
1	A	143	VAL
1	A	147	GLU
1	A	200	VAL
1	A	209	THR
1	B	15	THR
1	B	38	THR
1	B	44	VAL
1	B	82	GLU
1	B	106	LEU
1	B	107	GLU
1	B	116	VAL
1	B	126	ASN
1	B	143	VAL
1	B	147	GLU
1	B	161	GLU

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Mol	Chain	Res	Type
1	B	200	VAL
1	B	209	THR
1	C	15	THR
1	C	38	THR
1	C	44	VAL
1	C	58	VAL
1	C	82	GLU
1	C	106	LEU
1	C	107	GLU
1	C	116	VAL
1	C	126	ASN
1	C	143	VAL
1	C	147	GLU
1	C	161	GLU
1	C	200	VAL
1	C	209	THR
1	D	15	THR
1	D	38	THR
1	D	44	VAL
1	D	58	VAL
1	D	82	GLU
1	D	106	LEU
1	D	107	GLU
1	D	116	VAL
1	D	126	ASN
1	D	143	VAL
1	D	147	GLU
1	D	161	GLU
1	D	200	VAL
1	D	209	THR
1	E	15	THR
1	E	38	THR
1	E	44	VAL
1	E	58	VAL
1	E	82	GLU
1	E	106	LEU
1	E	107	GLU
1	E	116	VAL
1	E	126	ASN
1	E	143	VAL
1	E	147	GLU
1	E	161	GLU

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Mol	Chain	Res	Type
1	E	200	VAL
1	E	209	THR
1	F	15	THR
1	F	38	THR
1	F	44	VAL
1	F	58	VAL
1	F	82	GLU
1	F	106	LEU
1	F	107	GLU
1	F	116	VAL
1	F	126	ASN
1	F	143	VAL
1	F	147	GLU
1	F	161	GLU
1	F	200	VAL
1	F	209	THR
1	G	6	GLU
1	G	15	THR
1	G	38	THR
1	G	44	VAL
1	G	58	VAL
1	G	82	GLU
1	G	106	LEU
1	G	107	GLU
1	G	116	VAL
1	G	126	ASN
1	G	143	VAL
1	G	147	GLU
1	G	161	GLU
1	G	200	VAL
1	G	209	THR
1	H	15	THR
1	H	38	THR
1	H	44	VAL
1	H	58	VAL
1	H	82	GLU
1	H	106	LEU
1	H	107	GLU
1	H	116	VAL
1	H	126	ASN
1	H	143	VAL
1	H	147	GLU

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Mol	Chain	Res	Type
1	H	161	GLU
1	H	200	VAL
1	H	209	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	47	GLN
1	A	96	HIS
1	A	126	ASN
1	A	165	ASN
1	B	47	GLN
1	B	96	HIS
1	B	126	ASN
1	C	47	GLN
1	C	78	HIS
1	C	96	HIS
1	C	126	ASN
1	D	47	GLN
1	D	96	HIS
1	D	126	ASN
1	E	47	GLN
1	E	96	HIS
1	E	126	ASN
1	F	47	GLN
1	F	96	HIS
1	F	126	ASN
1	G	47	GLN
1	G	96	HIS
1	G	126	ASN
1	H	47	GLN
1	H	96	HIS
1	H	126	ASN

5.3.3 RNA ⓘ

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](#)

8 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
2	3PP	H	1226	-	8,8,8	2.49	4 (50%)	11,11,11	2.25	3 (27%)
2	3PP	E	1226	-	8,8,8	2.18	4 (50%)	11,11,11	2.39	3 (27%)
2	3PP	F	1226	-	8,8,8	2.49	3 (37%)	11,11,11	2.29	3 (27%)
2	3PP	G	1226	-	8,8,8	2.03	3 (37%)	11,11,11	2.37	4 (36%)
2	3PP	A	1226	-	8,8,8	2.19	3 (37%)	11,11,11	2.39	3 (27%)
2	3PP	B	1226	-	8,8,8	2.25	3 (37%)	11,11,11	2.17	4 (36%)
2	3PP	D	1226	-	8,8,8	2.66	4 (50%)	11,11,11	2.23	3 (27%)
2	3PP	C	1226	-	8,8,8	1.99	3 (37%)	11,11,11	2.32	3 (27%)

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	3PP	H	1226	-	-	2/6/6/6	-
2	3PP	E	1226	-	-	2/6/6/6	-
2	3PP	F	1226	-	-	2/6/6/6	-
2	3PP	G	1226	-	-	2/6/6/6	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	3PP	A	1226	-	-	2/6/6/6	-
2	3PP	B	1226	-	-	2/6/6/6	-
2	3PP	D	1226	-	-	5/6/6/6	-
2	3PP	C	1226	-	-	1/6/6/6	-

All (27) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	1226	3PP	P-CB	6.20	1.85	1.78
2	H	1226	3PP	P-CB	5.37	1.84	1.78
2	F	1226	3PP	P-CB	5.24	1.84	1.78
2	A	1226	3PP	P-CB	4.64	1.83	1.78
2	B	1226	3PP	P-CB	4.64	1.83	1.78
2	E	1226	3PP	P-CB	4.59	1.83	1.78
2	G	1226	3PP	P-CB	3.68	1.82	1.78
2	C	1226	3PP	P-CB	3.28	1.82	1.78
2	C	1226	3PP	P-O3	-2.87	1.48	1.55
2	G	1226	3PP	P-O3	-2.81	1.48	1.55
2	C	1226	3PP	P-O1	-2.81	1.48	1.55
2	F	1226	3PP	P-O3	-2.67	1.49	1.55
2	G	1226	3PP	P-O1	-2.52	1.49	1.55
2	F	1226	3PP	P-O1	-2.50	1.49	1.55
2	E	1226	3PP	P-O3	-2.40	1.49	1.55
2	A	1226	3PP	P-O3	-2.40	1.49	1.55
2	H	1226	3PP	P-O1	-2.35	1.49	1.55
2	B	1226	3PP	P-O3	-2.33	1.49	1.55
2	H	1226	3PP	CG-CD	2.30	1.55	1.50
2	E	1226	3PP	P-O1	-2.29	1.49	1.55
2	B	1226	3PP	P-O1	-2.17	1.50	1.55
2	D	1226	3PP	CG-CD	2.14	1.55	1.50
2	H	1226	3PP	P-O3	-2.14	1.50	1.55
2	E	1226	3PP	OE2-CD	-2.10	1.23	1.30
2	D	1226	3PP	P-O3	-2.08	1.50	1.55
2	A	1226	3PP	P-O1	-2.05	1.50	1.55
2	D	1226	3PP	P-O1	-2.03	1.50	1.55

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	1226	3PP	O2-P-CB	-5.78	100.56	111.45
2	A	1226	3PP	O2-P-CB	-5.69	100.74	111.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	1226	3PP	O2-P-CB	-5.40	101.27	111.45
2	H	1226	3PP	O2-P-CB	-5.36	101.35	111.45
2	G	1226	3PP	O2-P-CB	-5.35	101.36	111.45
2	D	1226	3PP	O2-P-CB	-5.18	101.69	111.45
2	B	1226	3PP	O2-P-CB	-4.72	102.56	111.45
2	C	1226	3PP	O2-P-CB	-4.63	102.73	111.45
2	C	1226	3PP	O3-P-CB	-3.75	97.96	106.91
2	D	1226	3PP	O3-P-CB	-3.19	99.29	106.91
2	G	1226	3PP	O3-P-O1	2.96	116.41	107.96
2	C	1226	3PP	O3-P-O1	2.96	116.39	107.96
2	A	1226	3PP	O3-P-O1	2.92	116.28	107.96
2	A	1226	3PP	O3-P-CB	-2.89	100.02	106.91
2	E	1226	3PP	O3-P-CB	-2.89	100.02	106.91
2	E	1226	3PP	O3-P-O1	2.88	116.17	107.96
2	F	1226	3PP	O3-P-CB	-2.83	100.15	106.91
2	G	1226	3PP	O3-P-CB	-2.82	100.17	106.91
2	F	1226	3PP	O3-P-O1	2.81	115.97	107.96
2	B	1226	3PP	O3-P-O1	2.80	115.94	107.96
2	H	1226	3PP	O3-P-O1	2.76	115.82	107.96
2	B	1226	3PP	O3-P-CB	-2.73	100.38	106.91
2	D	1226	3PP	O3-P-O1	2.70	115.67	107.96
2	H	1226	3PP	O3-P-CB	-2.70	100.47	106.91
2	B	1226	3PP	O1-P-CB	-2.24	101.56	106.91
2	G	1226	3PP	O1-P-CB	-2.22	101.61	106.91

There are no chirality outliers.

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	1226	3PP	CG-CB-P-O1
2	D	1226	3PP	CG-CB-P-O2
2	D	1226	3PP	CG-CB-P-O3
2	F	1226	3PP	OE1-CD-CG-CB
2	D	1226	3PP	OE2-CD-CG-CB
2	F	1226	3PP	OE2-CD-CG-CB
2	H	1226	3PP	OE2-CD-CG-CB
2	D	1226	3PP	OE1-CD-CG-CB
2	E	1226	3PP	OE2-CD-CG-CB
2	H	1226	3PP	OE1-CD-CG-CB
2	A	1226	3PP	OE2-CD-CG-CB
2	B	1226	3PP	OE2-CD-CG-CB
2	C	1226	3PP	OE2-CD-CG-CB

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Mol	Chain	Res	Type	Atoms
2	E	1226	3PP	OE1-CD-CG-CB
2	G	1226	3PP	OE2-CD-CG-CB
2	B	1226	3PP	OE1-CD-CG-CB
2	G	1226	3PP	OE1-CD-CG-CB
2	A	1226	3PP	OE1-CD-CG-CB

There are no ring outliers.

8 monomers are involved in 46 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	H	1226	3PP	4	0
2	E	1226	3PP	7	0
2	F	1226	3PP	5	0
2	G	1226	3PP	6	0
2	A	1226	3PP	8	0
2	B	1226	3PP	6	0
2	D	1226	3PP	4	0
2	C	1226	3PP	6	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.