Full wwPDB X-ray Structure Validation Report (i)

Jun 15, 2024 – 01:16 PM EDT

PDB ID	:	1T72
Title	:	Crystal structure of phosphate transport system protein phoU from Aquifex aeolicus
Authors	:	Oganesyan, V.; Kim, S.-H.; Oganesyan, N.; Jancarik, J.; Adams, P.D.; Kim, R.; Berkeley Structural Genomics Center (BSGC)
Deposited on	:	2004-05-07
Resolution	:	2.90 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 (i) 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 (i)) were used in the production of this report:

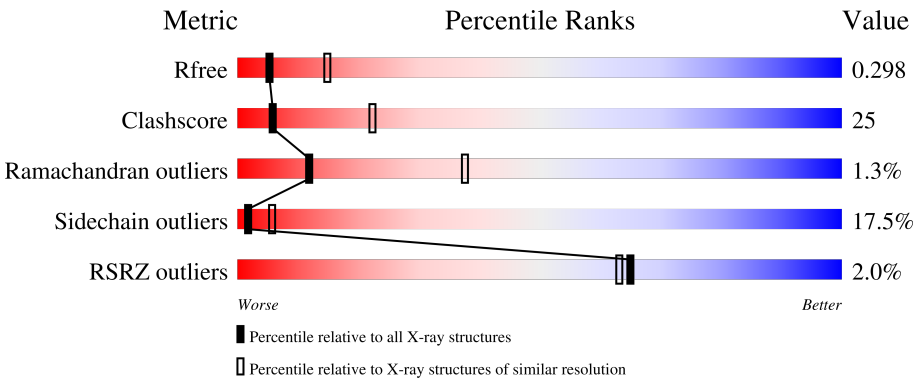
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (2.90-2.90)

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 <=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	227	3%; 47%, 37%, 11%, 5%
1	B	227	4%; 51%, 33%, 11%, 6%
1	D	227	2%; 51%, 37%, 7%, 5%
1	E	227	%; 47%, 37%, 10%, 6%
1	F	227	43%, 46%, 5%, 6%

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Mol	Chain	Length	Quality of chain
1	G	227	% 49% 34% 9% 7%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 10391 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 Phosphate transport system protein phoU homolog.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	A	215	1712	1076	286	339	1	10	0	0	0
1	B	214	1708	1074	285	338	1	10	0	0	0
1	D	215	1712	1076	286	339	1	10	0	0	0
1	E	213	1699	1069	284	335	1	10	0	0	0
1	F	214	1708	1074	285	338	1	10	0	0	0
1	G	211	1691	1065	282	333	1	10	0	0	0

There are 96 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	GLY	-	CLONING ARTIFACT	UNP O67053
A	2	GLY	-	CLONING ARTIFACT	UNP O67053
A	3	GLY	-	CLONING ARTIFACT	UNP O67053
A	4	GLY	-	CLONING ARTIFACT	UNP O67053
A	5	GLY	-	CLONING ARTIFACT	UNP O67053
A	6	GLY	-	CLONING ARTIFACT	UNP O67053
A	7	MSE	MET	MODIFIED RESIDUE	UNP O67053
A	23	MSE	MET	MODIFIED RESIDUE	UNP O67053
A	70	MSE	MET	MODIFIED RESIDUE	UNP O67053
A	83	MSE	MET	MODIFIED RESIDUE	UNP O67053
A	85	MSE	MET	MODIFIED RESIDUE	UNP O67053
A	97	MSE	MET	MODIFIED RESIDUE	UNP O67053
A	126	MSE	MET	MODIFIED RESIDUE	UNP O67053
A	133	MSE	MET	MODIFIED RESIDUE	UNP O67053
A	171	MSE	MET	MODIFIED RESIDUE	UNP O67053
A	185	MSE	MET	MODIFIED RESIDUE	UNP O67053
B	1	GLY	-	CLONING ARTIFACT	UNP O67053

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Chain	Residue	Modelled	Actual	Comment	Reference
B	2	GLY	-	CLONING ARTIFACT	UNP O67053
B	3	GLY	-	CLONING ARTIFACT	UNP O67053
B	4	GLY	-	CLONING ARTIFACT	UNP O67053
B	5	GLY	-	CLONING ARTIFACT	UNP O67053
B	6	GLY	-	CLONING ARTIFACT	UNP O67053
B	7	MSE	MET	MODIFIED RESIDUE	UNP O67053
B	23	MSE	MET	MODIFIED RESIDUE	UNP O67053
B	70	MSE	MET	MODIFIED RESIDUE	UNP O67053
B	83	MSE	MET	MODIFIED RESIDUE	UNP O67053
B	85	MSE	MET	MODIFIED RESIDUE	UNP O67053
B	97	MSE	MET	MODIFIED RESIDUE	UNP O67053
B	126	MSE	MET	MODIFIED RESIDUE	UNP O67053
B	133	MSE	MET	MODIFIED RESIDUE	UNP O67053
B	171	MSE	MET	MODIFIED RESIDUE	UNP O67053
B	185	MSE	MET	MODIFIED RESIDUE	UNP O67053
D	1	GLY	-	CLONING ARTIFACT	UNP O67053
D	2	GLY	-	CLONING ARTIFACT	UNP O67053
D	3	GLY	-	CLONING ARTIFACT	UNP O67053
D	4	GLY	-	CLONING ARTIFACT	UNP O67053
D	5	GLY	-	CLONING ARTIFACT	UNP O67053
D	6	GLY	-	CLONING ARTIFACT	UNP O67053
D	7	MSE	MET	MODIFIED RESIDUE	UNP O67053
D	23	MSE	MET	MODIFIED RESIDUE	UNP O67053
D	70	MSE	MET	MODIFIED RESIDUE	UNP O67053
D	83	MSE	MET	MODIFIED RESIDUE	UNP O67053
D	85	MSE	MET	MODIFIED RESIDUE	UNP O67053
D	97	MSE	MET	MODIFIED RESIDUE	UNP O67053
D	126	MSE	MET	MODIFIED RESIDUE	UNP O67053
D	133	MSE	MET	MODIFIED RESIDUE	UNP O67053
D	171	MSE	MET	MODIFIED RESIDUE	UNP O67053
D	185	MSE	MET	MODIFIED RESIDUE	UNP O67053
E	1	GLY	-	CLONING ARTIFACT	UNP O67053
E	2	GLY	-	CLONING ARTIFACT	UNP O67053
E	3	GLY	-	CLONING ARTIFACT	UNP O67053
E	4	GLY	-	CLONING ARTIFACT	UNP O67053
E	5	GLY	-	CLONING ARTIFACT	UNP O67053
E	6	GLY	-	CLONING ARTIFACT	UNP O67053
E	7	MSE	MET	MODIFIED RESIDUE	UNP O67053
E	23	MSE	MET	MODIFIED RESIDUE	UNP O67053
E	70	MSE	MET	MODIFIED RESIDUE	UNP O67053
E	83	MSE	MET	MODIFIED RESIDUE	UNP O67053
E	85	MSE	MET	MODIFIED RESIDUE	UNP O67053

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Chain	Residue	Modelled	Actual	Comment	Reference
E	97	MSE	MET	MODIFIED RESIDUE	UNP O67053
E	126	MSE	MET	MODIFIED RESIDUE	UNP O67053
E	133	MSE	MET	MODIFIED RESIDUE	UNP O67053
E	171	MSE	MET	MODIFIED RESIDUE	UNP O67053
E	185	MSE	MET	MODIFIED RESIDUE	UNP O67053
F	1	GLY	-	CLONING ARTIFACT	UNP O67053
F	2	GLY	-	CLONING ARTIFACT	UNP O67053
F	3	GLY	-	CLONING ARTIFACT	UNP O67053
F	4	GLY	-	CLONING ARTIFACT	UNP O67053
F	5	GLY	-	CLONING ARTIFACT	UNP O67053
F	6	GLY	-	CLONING ARTIFACT	UNP O67053
F	7	MSE	MET	MODIFIED RESIDUE	UNP O67053
F	23	MSE	MET	MODIFIED RESIDUE	UNP O67053
F	70	MSE	MET	MODIFIED RESIDUE	UNP O67053
F	83	MSE	MET	MODIFIED RESIDUE	UNP O67053
F	85	MSE	MET	MODIFIED RESIDUE	UNP O67053
F	97	MSE	MET	MODIFIED RESIDUE	UNP O67053
F	126	MSE	MET	MODIFIED RESIDUE	UNP O67053
F	133	MSE	MET	MODIFIED RESIDUE	UNP O67053
F	171	MSE	MET	MODIFIED RESIDUE	UNP O67053
F	185	MSE	MET	MODIFIED RESIDUE	UNP O67053
G	1	GLY	-	CLONING ARTIFACT	UNP O67053
G	2	GLY	-	CLONING ARTIFACT	UNP O67053
G	3	GLY	-	CLONING ARTIFACT	UNP O67053
G	4	GLY	-	CLONING ARTIFACT	UNP O67053
G	5	GLY	-	CLONING ARTIFACT	UNP O67053
G	6	GLY	-	CLONING ARTIFACT	UNP O67053
G	7	MSE	MET	MODIFIED RESIDUE	UNP O67053
G	23	MSE	MET	MODIFIED RESIDUE	UNP O67053
G	70	MSE	MET	MODIFIED RESIDUE	UNP O67053
G	83	MSE	MET	MODIFIED RESIDUE	UNP O67053
G	85	MSE	MET	MODIFIED RESIDUE	UNP O67053
G	97	MSE	MET	MODIFIED RESIDUE	UNP O67053
G	126	MSE	MET	MODIFIED RESIDUE	UNP O67053
G	133	MSE	MET	MODIFIED RESIDUE	UNP O67053
G	171	MSE	MET	MODIFIED RESIDUE	UNP O67053
G	185	MSE	MET	MODIFIED RESIDUE	UNP O67053

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	33	33	33	0	0

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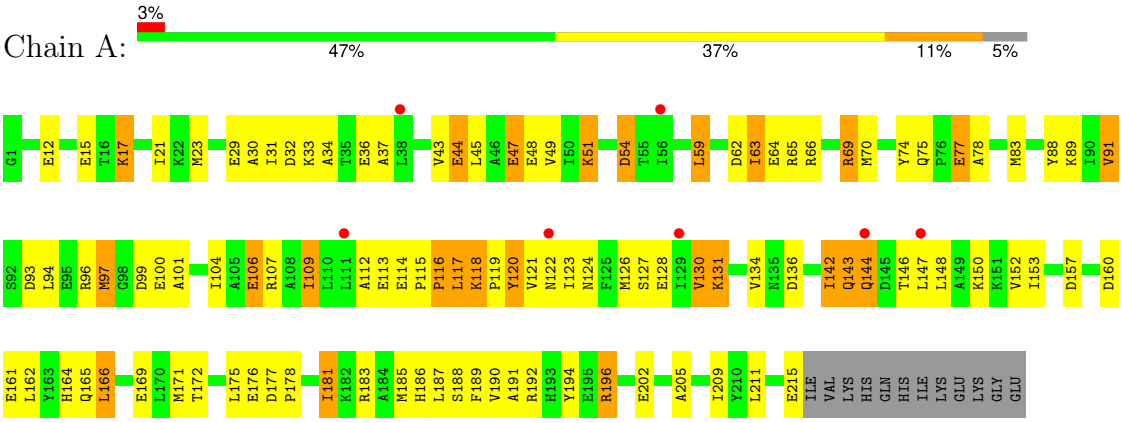
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	B	26	26	26	0	0
			Total	O		
2	D	30	30	30	0	0
			Total	O		
2	E	18	18	18	0	0
			Total	O		
2	F	27	27	27	0	0
			Total	O		
2	G	27	27	27	0	0

3 Residue-property plots (i)

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

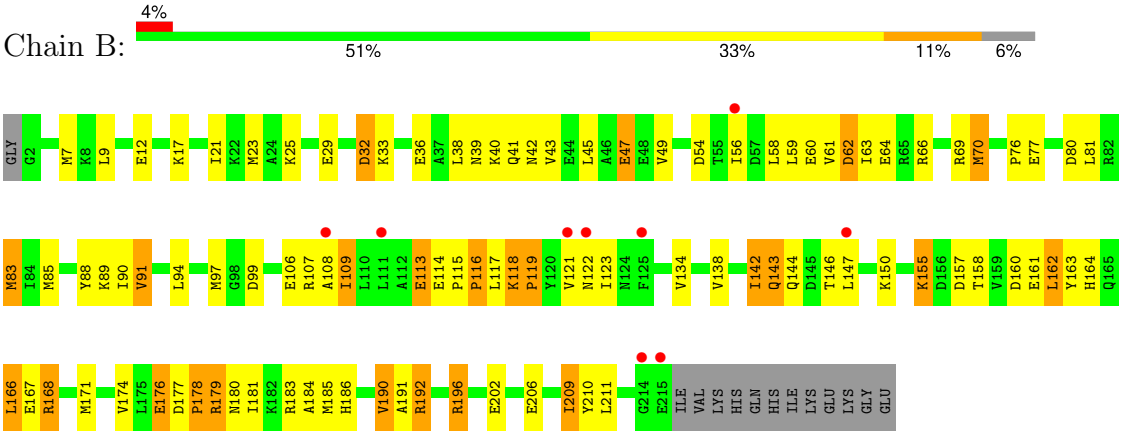
• Molecule 1: Phosphate transport system protein phoU homolog

Chain A:



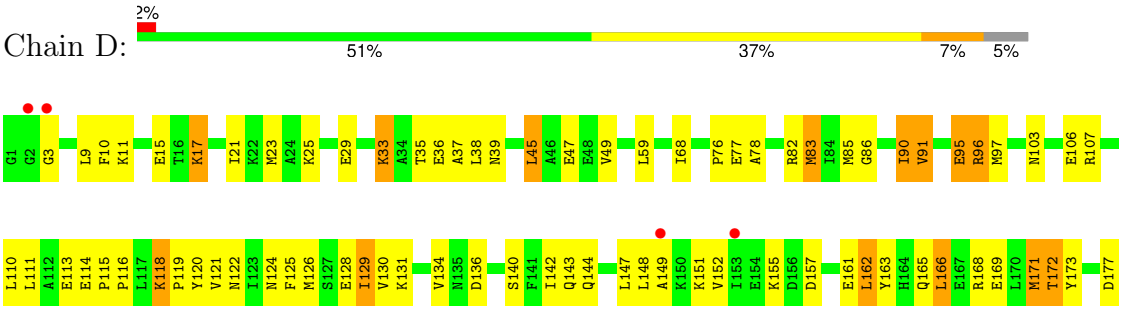
• Molecule 1: Phosphate transport system protein phoU homolog

Chain B:



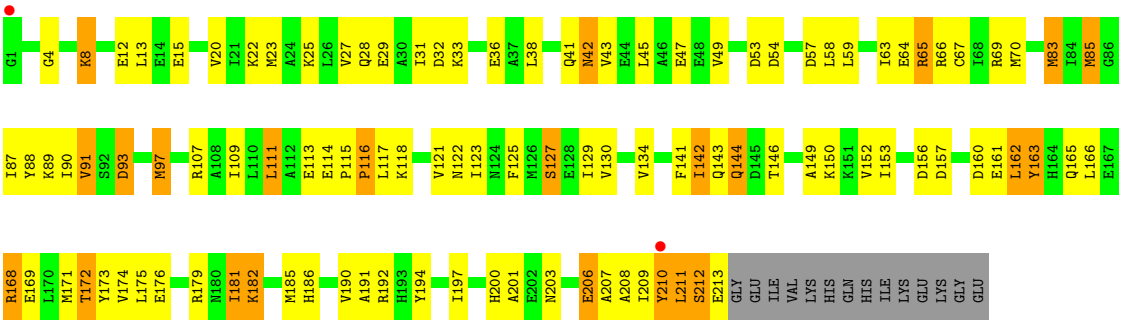
• Molecule 1: Phosphate transport system protein phoU homolog

Chain D:



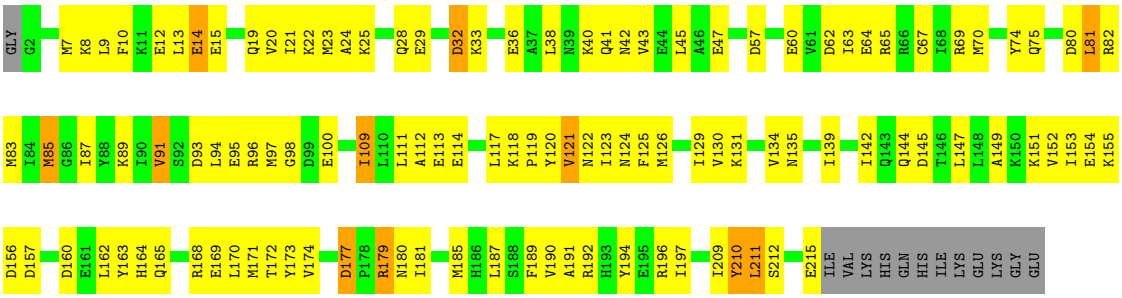
• Molecule 1: Phosphate transport system protein phoU homolog

Chain E:



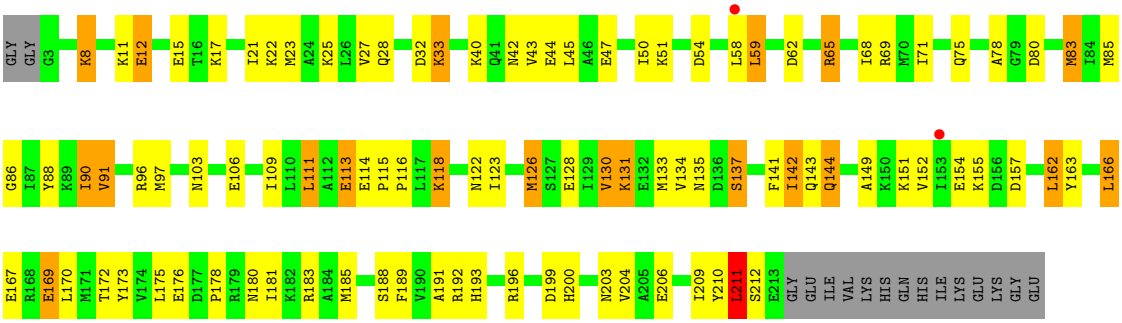
• Molecule 1: Phosphate transport system protein phoU homolog

Chain F:



• Molecule 1: Phosphate transport system protein phoU homolog

Chain G:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 43	Depositor
Cell constants a, b, c, α, β, γ	113.50Å 113.50Å 155.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.00 – 2.90 27.91 – 2.80	Depositor EDS
% Data completeness (in resolution range)	84.5 (15.00-2.90) 84.5 (27.91-2.80)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.09	Depositor
$< I/\sigma(I) >$ [1]	2.74 (at 2.80Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.201 , 0.293 0.240 , 0.298	Depositor DCC
R_{free} test set	2050 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å^2)	52.1	Xtriage
Anisotropy	0.132	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 1.0	EDS
L-test for twinning[2]	$< \|L\| > = 0.47$, $< L^2 > = 0.30$	Xtriage
Estimated twinning fraction	0.488 for h,-k,-l	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	10391	wwPDB-VP
Average B, all atoms (Å^2)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 66.64 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 5.8534e-06. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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)

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	0.65	1/1721 (0.1%)	0.98	11/2298 (0.5%)
1	B	0.63	0/1717	1.07	9/2293 (0.4%)
1	D	0.63	0/1721	0.92	6/2298 (0.3%)
1	E	0.67	1/1708 (0.1%)	1.09	11/2281 (0.5%)
1	F	0.67	0/1717	0.90	7/2293 (0.3%)
1	G	0.63	0/1700	0.96	7/2271 (0.3%)
All	All	0.65	2/10284 (0.0%)	0.99	51/13734 (0.4%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	97	MSE	SE-CE	-5.21	1.64	1.95
1	E	97	MSE	SE-CE	-5.17	1.65	1.95

All (51) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	E	115	PRO	CA-N-CD	-22.31	80.26	111.50
1	B	119	PRO	CA-N-CD	-19.37	84.38	111.50
1	B	116	PRO	CA-N-CD	-17.55	86.93	111.50
1	A	116	PRO	CA-N-CD	-17.50	87.00	111.50
1	D	76	PRO	CA-N-CD	-17.13	87.52	111.50
1	B	115	PRO	CA-N-CD	-15.26	90.14	111.50
1	E	116	PRO	CA-N-CD	-15.08	90.38	111.50
1	G	115	PRO	CA-N-CD	-14.84	90.72	111.50
1	G	116	PRO	CA-N-CD	-13.95	91.98	111.50
1	A	119	PRO	CA-N-CD	-13.38	92.76	111.50
1	B	178	PRO	CA-N-CD	-11.22	95.80	111.50
1	A	115	PRO	CA-N-CD	-10.41	96.92	111.50
1	D	76	PRO	C-N-CA	-8.12	101.41	121.70
1	E	160	ASP	CB-CG-OD2	7.31	124.88	118.30
1	F	57	ASP	CB-CG-OD2	7.26	124.83	118.30
1	F	93	ASP	CB-CG-OD2	7.01	124.61	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	E	54	ASP	CB-CG-OD2	6.66	124.29	118.30
1	E	157	ASP	CB-CG-OD2	6.56	124.20	118.30
1	E	115	PRO	N-CA-CB	6.50	111.10	103.30
1	F	157	ASP	CB-CG-OD2	6.45	124.11	118.30
1	A	99	ASP	CB-CG-OD2	6.35	124.02	118.30
1	E	156	ASP	CB-CG-OD2	6.26	123.94	118.30
1	B	54	ASP	CB-CG-OD2	6.16	123.84	118.30
1	B	99	ASP	CB-CG-OD2	6.10	123.79	118.30
1	B	160	ASP	CB-CG-OD2	5.98	123.69	118.30
1	E	57	ASP	CB-CG-OD2	5.80	123.52	118.30
1	G	199	ASP	CB-CG-OD2	5.70	123.43	118.30
1	F	80	ASP	CB-CG-OD2	5.63	123.37	118.30
1	A	177	ASP	CB-CG-OD2	5.62	123.36	118.30
1	G	111	LEU	CA-CB-CG	-5.61	102.41	115.30
1	E	93	ASP	CB-CG-OD2	5.59	123.33	118.30
1	D	157	ASP	CB-CG-OD2	5.55	123.30	118.30
1	G	157	ASP	CB-CG-OD2	5.54	123.28	118.30
1	F	160	ASP	CB-CG-OD2	5.52	123.27	118.30
1	A	54	ASP	CB-CG-OD2	5.50	123.25	118.30
1	E	53	ASP	CB-CG-OD2	5.47	123.22	118.30
1	D	77	GLU	CB-CA-C	5.38	121.15	110.40
1	B	157	ASP	CB-CG-OD2	5.34	123.11	118.30
1	A	32	ASP	CB-CG-OD2	5.33	123.10	118.30
1	A	160	ASP	CB-CG-OD2	5.30	123.08	118.30
1	D	136	ASP	CB-CG-OD2	5.29	123.06	118.30
1	G	80	ASP	CB-CG-OD2	5.25	123.02	118.30
1	A	93	ASP	CB-CG-OD2	5.24	123.02	118.30
1	G	54	ASP	CB-CG-OD2	5.22	123.00	118.30
1	A	136	ASP	CB-CG-OD2	5.21	122.99	118.30
1	E	144	GLN	CB-CA-C	-5.17	100.06	110.40
1	F	62	ASP	CB-CG-OD2	5.15	122.94	118.30
1	B	32	ASP	CB-CG-OD2	5.10	122.89	118.30
1	A	157	ASP	CB-CG-OD2	5.05	122.85	118.30
1	D	199	ASP	CB-CG-OD2	5.05	122.85	118.30
1	F	177	ASP	CB-CG-OD2	5.02	122.82	118.30

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	1712	0	1727	96	0
1	B	1708	0	1721	108	1
1	D	1712	0	1727	79	1
1	E	1699	0	1718	117	0
1	F	1708	0	1721	106	0
1	G	1691	0	1709	80	0
2	A	33	0	0	2	0
2	B	26	0	0	3	0
2	D	30	0	0	3	0
2	E	18	0	0	0	0
2	F	27	0	0	0	0
2	G	27	0	0	2	0
All	All	10391	0	10323	517	1

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 25.

All (517) 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:F:85:MSE:SE	1:G:85:MSE:HE1	1.82	1.29
1:G:97:MSE:HE3	1:G:134:VAL:HG22	1.25	1.16
1:B:97:MSE:CE	1:B:134:VAL:HG22	1.75	1.16
1:D:97:MSE:CE	1:D:134:VAL:HG22	1.78	1.13
1:F:13:LEU:HA	1:F:70:MSE:HE1	1.34	1.10
1:B:168:ARG:HD2	1:D:172:THR:HG21	1.26	1.08
1:E:172:THR:HG21	1:F:168:ARG:HG2	1.11	1.06
1:E:97:MSE:CE	1:E:134:VAL:HG22	1.86	1.05
1:B:107:ARG:HH11	1:B:186:HIS:CD2	1.76	1.02
1:D:97:MSE:HE2	1:D:134:VAL:HG22	1.38	1.02
1:G:65:ARG:HH11	1:G:65:ARG:HG2	0.86	1.01
1:A:59:LEU:O	1:A:63:ILE:HG13	1.61	0.99
1:E:172:THR:HG22	1:F:168:ARG:HH11	1.23	0.98
1:E:174:VAL:HG22	1:E:181:ILE:HD12	1.43	0.98
1:A:122:ASN:HB2	2:A:233:HOH:O	1.63	0.97

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:13:LEU:CA	1:F:70:MSE:HE1	1.94	0.97
1:G:200:HIS:O	1:G:203:ASN:HB2	1.65	0.96
1:G:65:ARG:HH11	1:G:65:ARG:CG	1.78	0.96
1:A:83:MSE:SE	1:A:211:LEU:HD21	2.15	0.95
1:G:17:LYS:O	1:G:21:ILE:HD12	1.66	0.95
1:E:172:THR:CG2	1:F:168:ARG:HH11	1.79	0.95
1:E:172:THR:HG22	1:F:168:ARG:NH1	1.80	0.94
1:B:107:ARG:HH11	1:B:186:HIS:HD2	0.95	0.94
1:G:189:PHE:HA	1:G:192:ARG:HD2	1.49	0.94
1:E:97:MSE:HE3	1:E:134:VAL:HG22	1.48	0.94
1:F:97:MSE:CE	1:F:134:VAL:HG22	1.97	0.94
1:A:97:MSE:CE	1:A:134:VAL:HG22	1.98	0.94
1:A:189:PHE:HZ	1:B:163:TYR:OH	1.49	0.93
1:A:106:GLU:OE2	1:A:106:GLU:HA	1.69	0.93
1:A:196:ARG:HE	1:B:196:ARG:HH21	1.14	0.93
1:G:97:MSE:CE	1:G:134:VAL:HG22	1.98	0.92
1:B:23:MSE:SE	1:B:91:VAL:HG22	2.20	0.92
1:A:189:PHE:HZ	1:B:163:TYR:HH	0.93	0.92
1:G:65:ARG:HG2	1:G:65:ARG:NH1	1.66	0.91
1:A:107:ARG:HH11	1:A:186:HIS:CD2	1.86	0.91
1:A:107:ARG:HH11	1:A:186:HIS:HD2	1.13	0.91
1:B:106:GLU:OE2	1:B:106:GLU:HA	1.70	0.90
1:B:168:ARG:HD2	1:D:172:THR:CG2	2.02	0.90
1:F:100:GLU:HA	1:F:100:GLU:OE1	1.69	0.89
1:E:172:THR:CG2	1:F:168:ARG:NH1	2.38	0.87
1:A:83:MSE:SE	1:A:211:LEU:CD2	2.74	0.86
1:E:97:MSE:HG2	1:E:197:ILE:HG23	1.58	0.85
1:A:43:VAL:O	1:A:47:GLU:HG3	1.77	0.85
1:B:97:MSE:HE1	1:B:134:VAL:HG22	1.57	0.85
1:B:178:PRO:O	1:B:181:ILE:HG13	1.77	0.84
1:G:86:GLY:O	1:G:90:ILE:HG23	1.77	0.84
1:A:120:TYR:O	1:A:124:ASN:ND2	2.11	0.84
1:G:114:GLU:HB3	1:G:183:ARG:HD3	1.59	0.84
1:B:166:LEU:HD22	1:B:191:ALA:HB2	1.60	0.83
1:B:97:MSE:HE3	1:B:134:VAL:HG22	1.61	0.82
1:D:95:GLU:OE2	1:D:96:ARG:NH1	2.12	0.82
1:D:177:ASP:O	1:D:180:ASN:HB2	1.80	0.82
1:A:166:LEU:HD22	1:A:191:ALA:HB2	1.61	0.82
1:E:172:THR:CG2	1:F:168:ARG:HG2	2.04	0.82
1:B:117:LEU:HD13	1:B:183:ARG:CZ	2.10	0.81
1:F:163:TYR:OH	1:G:192:ARG:NH2	2.13	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:23:MSE:SE	1:E:91:VAL:HG22	2.32	0.79
1:F:97:MSE:HE1	1:F:134:VAL:HG13	1.65	0.79
1:E:142:ILE:HG22	1:E:143:GLN:CG	2.14	0.78
1:B:113:GLU:O	1:B:114:GLU:HG3	1.82	0.78
1:A:43:VAL:HG13	1:A:109:ILE:HD13	1.65	0.76
1:D:86:GLY:O	1:D:90:ILE:HG23	1.86	0.76
1:A:23:MSE:HG3	1:A:59:LEU:HD23	1.67	0.76
1:B:107:ARG:NH1	1:B:186:HIS:CD2	2.52	0.76
1:D:171:MSE:SE	1:E:181:ILE:HG13	2.36	0.76
1:E:142:ILE:HG22	1:E:143:GLN:HG3	1.69	0.75
1:E:122:ASN:OD1	1:F:165:GLN:NE2	2.19	0.75
1:D:118:LYS:HD2	1:D:169:GLU:OE2	1.86	0.75
1:A:146:THR:HG22	1:A:150:LYS:HE3	1.69	0.74
1:B:168:ARG:NH2	1:D:173:TYR:CZ	2.55	0.74
1:E:13:LEU:HA	1:E:70:MSE:HE1	1.67	0.74
1:E:161:GLU:HG2	1:F:119:PRO:HG2	1.68	0.74
1:B:83:MSE:HE2	1:B:144:GLN:NE2	2.02	0.74
1:B:146:THR:CG2	1:B:150:LYS:HE3	2.18	0.74
1:F:85:MSE:SE	1:G:85:MSE:CE	2.76	0.73
1:E:200:HIS:O	1:E:203:ASN:HB2	1.86	0.73
1:E:174:VAL:HG22	1:E:181:ILE:CD1	2.18	0.73
1:B:36:GLU:HB3	1:B:45:LEU:CD2	2.18	0.73
1:A:178:PRO:O	1:A:181:ILE:HB	1.88	0.73
1:G:17:LYS:O	1:G:21:ILE:CD1	2.37	0.73
1:A:65:ARG:HE	1:A:69:ARG:NH2	1.87	0.73
1:B:142:ILE:HG22	1:B:143:GLN:HG2	1.71	0.72
1:G:118:LYS:HB3	1:G:173:TYR:OH	1.89	0.72
1:E:210:TYR:C	1:E:210:TYR:CD2	2.63	0.72
1:B:174:VAL:O	1:B:178:PRO:HD3	1.89	0.72
1:F:122:ASN:ND2	1:F:169:GLU:OE2	2.21	0.72
1:E:97:MSE:HE1	1:E:134:VAL:HG13	1.71	0.72
1:F:97:MSE:HE3	1:F:134:VAL:HG22	1.70	0.72
1:A:127:SER:HB3	1:A:194:TYR:CE2	2.25	0.71
1:B:121:VAL:HG21	1:D:161:GLU:HB3	1.72	0.71
1:E:13:LEU:CA	1:E:70:MSE:HE1	2.20	0.71
1:A:166:LEU:HD22	1:A:191:ALA:CB	2.21	0.70
1:F:43:VAL:HG22	1:F:109:ILE:HG13	1.72	0.70
1:B:117:LEU:HD13	1:B:183:ARG:NH2	2.05	0.70
1:E:210:TYR:C	1:E:210:TYR:HD2	1.95	0.69
1:A:153:ILE:HD12	1:A:202:GLU:HG2	1.74	0.69
1:D:118:LYS:HB3	1:D:173:TYR:OH	1.93	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:196:ARG:NE	1:B:196:ARG:HH21	1.88	0.69
1:B:146:THR:HG22	1:B:150:LYS:HE3	1.73	0.69
1:D:97:MSE:HE2	1:D:134:VAL:CG2	2.21	0.69
1:F:13:LEU:CA	1:F:70:MSE:CE	2.70	0.69
1:B:62:ASP:O	1:B:66:ARG:HG3	1.92	0.69
1:B:83:MSE:SE	1:B:211:LEU:HD21	2.44	0.68
1:F:97:MSE:HE1	1:F:134:VAL:HG22	1.75	0.68
1:B:121:VAL:HG11	1:D:162:LEU:HD13	1.76	0.68
1:A:17:LYS:HE3	1:A:83:MSE:HE2	1.74	0.68
1:D:68:ILE:HG12	1:E:85:MSE:CE	2.23	0.67
1:B:117:LEU:O	1:B:118:LYS:HB3	1.93	0.67
1:G:113:GLU:O	1:G:114:GLU:HG2	1.95	0.67
1:F:210:TYR:C	1:F:210:TYR:CD2	2.67	0.66
1:B:59:LEU:O	1:B:63:ILE:HD12	1.95	0.66
1:E:117:LEU:HD11	1:E:173:TYR:CD2	2.31	0.66
1:F:85:MSE:HE3	1:G:68:ILE:HG12	1.77	0.66
1:F:111:LEU:HD11	1:F:187:LEU:HD23	1.77	0.66
1:A:189:PHE:CZ	1:B:163:TYR:OH	2.40	0.66
1:F:144:GLN:OE1	1:F:215:GLU:OE1	2.13	0.65
1:A:196:ARG:HE	1:B:196:ARG:NH2	1.91	0.65
1:E:97:MSE:HE1	1:E:134:VAL:HG22	1.76	0.65
1:A:36:GLU:HB3	1:A:45:LEU:HD23	1.79	0.65
1:A:107:ARG:NH1	1:A:186:HIS:CD2	2.63	0.65
1:E:122:ASN:ND2	1:E:169:GLU:OE2	2.29	0.65
1:F:33:LYS:HD3	1:F:45:LEU:HD11	1.79	0.65
1:F:117:LEU:HD11	1:F:173:TYR:CD2	2.32	0.64
1:E:111:LEU:HD22	1:E:186:HIS:CB	2.27	0.64
1:D:140:SER:HB3	1:D:148:LEU:HD23	1.80	0.64
1:G:118:LYS:NZ	1:G:169:GLU:OE2	2.19	0.64
1:A:118:LYS:O	1:A:120:TYR:CD2	2.51	0.64
1:B:17:LYS:HE3	1:B:83:MSE:HE3	1.79	0.64
1:D:144:GLN:OE1	1:D:211:LEU:HD21	1.98	0.63
1:A:36:GLU:HB3	1:A:45:LEU:CD2	2.27	0.63
1:B:161:GLU:HG2	1:D:119:PRO:HG3	1.79	0.63
1:F:85:MSE:CE	1:G:68:ILE:HG12	2.28	0.63
1:F:75:GLN:NE2	1:G:75:GLN:HG2	2.14	0.63
1:F:171:MSE:SE	1:G:181:ILE:HG23	2.49	0.63
1:B:162:LEU:HD13	1:D:121:VAL:HG11	1.81	0.62
1:F:163:TYR:HD2	1:F:164:HIS:HD2	1.45	0.62
1:D:90:ILE:HB	1:D:204:VAL:HG22	1.79	0.62
1:E:122:ASN:OD1	1:F:165:GLN:CD	2.38	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:23:MSE:CB	1:B:59:LEU:HD23	2.29	0.62
1:D:166:LEU:HD22	1:D:191:ALA:HB2	1.82	0.62
1:E:13:LEU:CB	1:E:70:MSE:HE1	2.29	0.62
1:F:100:GLU:OE2	1:F:196:ARG:HD2	2.00	0.62
1:A:106:GLU:OE2	1:A:106:GLU:CA	2.44	0.61
1:B:43:VAL:HG13	1:B:109:ILE:HD13	1.82	0.61
1:B:142:ILE:HG22	1:B:143:GLN:HE21	1.65	0.61
1:E:97:MSE:CE	1:E:134:VAL:CG2	2.73	0.61
1:B:186:HIS:O	1:B:190:VAL:HG23	2.01	0.61
1:A:118:LYS:O	1:A:120:TYR:CE2	2.54	0.61
1:D:192:ARG:HE	1:E:192:ARG:CZ	2.14	0.61
1:G:90:ILE:HB	1:G:204:VAL:HG22	1.83	0.61
1:B:180:ASN:OD1	1:B:183:ARG:NH2	2.34	0.61
1:A:63:ILE:HB	1:A:91:VAL:HG21	1.83	0.60
1:D:106:GLU:HG3	2:D:245:HOH:O	2.01	0.60
1:E:142:ILE:HG22	1:E:143:GLN:HG2	1.82	0.60
1:E:144:GLN:HE22	1:E:211:LEU:HD12	1.66	0.60
1:G:43:VAL:HG22	1:G:109:ILE:HG12	1.84	0.60
1:A:83:MSE:SE	1:A:211:LEU:HD23	2.51	0.60
1:A:117:LEU:O	1:A:117:LEU:HG	2.02	0.59
1:D:45:LEU:O	1:D:49:VAL:HG23	2.03	0.59
1:E:20:VAL:HG21	1:E:87:ILE:HG23	1.85	0.59
1:F:85:MSE:HG2	1:G:88:TYR:CE2	2.37	0.59
1:D:23:MSE:SE	1:D:91:VAL:HG22	2.52	0.59
1:E:142:ILE:CG2	1:E:143:GLN:HG2	2.32	0.59
1:E:144:GLN:HE22	1:E:211:LEU:CD1	2.16	0.59
1:A:97:MSE:HE2	1:A:134:VAL:HG22	1.81	0.59
1:D:147:LEU:H	1:D:147:LEU:HD12	1.66	0.59
1:F:13:LEU:CB	1:F:70:MSE:HE1	2.32	0.59
1:A:196:ARG:HH21	1:B:196:ARG:HE	1.50	0.59
1:B:161:GLU:HG2	1:D:119:PRO:CG	2.33	0.58
1:E:169:GLU:OE2	1:F:165:GLN:HG3	2.03	0.58
1:G:97:MSE:HE1	1:G:134:VAL:HG13	1.85	0.58
1:G:188:SER:O	1:G:192:ARG:HG3	2.04	0.58
1:A:12:GLU:OE2	1:A:74:TYR:OH	2.14	0.58
1:B:36:GLU:HB3	1:B:45:LEU:HD23	1.83	0.58
1:D:144:GLN:OE1	1:D:211:LEU:CD2	2.51	0.58
1:E:41:GLN:NE2	1:E:116:PRO:HD3	2.19	0.58
1:B:17:LYS:O	1:B:21:ILE:HD12	2.04	0.58
1:E:161:GLU:HG2	1:F:119:PRO:CG	2.34	0.58
1:E:146:THR:HG22	1:E:150:LYS:HE3	1.86	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:122:ASN:ND2	1:D:166:LEU:HD12	2.19	0.57
1:A:75:GLN:HG2	1:A:75:GLN:O	2.04	0.57
1:A:101:ALA:HA	1:A:104:ILE:HD12	1.86	0.57
1:G:181:ILE:O	1:G:185:MSE:HG3	2.05	0.57
1:A:65:ARG:O	1:A:69:ARG:HD2	2.05	0.57
1:E:111:LEU:HD22	1:E:186:HIS:HB3	1.87	0.57
1:D:130:VAL:HG12	1:D:131:LYS:N	2.20	0.57
1:E:42:ASN:OD1	1:E:42:ASN:C	2.42	0.56
1:A:192:ARG:CZ	1:B:192:ARG:NH1	2.68	0.56
1:A:192:ARG:CZ	1:B:192:ARG:CZ	2.82	0.56
1:E:161:GLU:HB3	1:F:121:VAL:HG21	1.87	0.56
1:A:130:VAL:HG12	1:A:131:LYS:N	2.20	0.56
1:G:83:MSE:SE	1:G:211:LEU:HD11	2.56	0.56
1:B:168:ARG:HB3	1:D:172:THR:HG21	1.87	0.56
1:G:65:ARG:CG	1:G:65:ARG:NH1	2.46	0.56
1:D:210:TYR:O	1:D:211:LEU:C	2.45	0.56
1:A:127:SER:CB	1:A:194:TYR:CE2	2.89	0.55
1:F:36:GLU:HG2	1:F:40:LYS:HD2	1.88	0.55
1:F:63:ILE:HG21	1:F:91:VAL:HG21	1.87	0.55
1:F:97:MSE:HE1	1:F:134:VAL:CG1	2.36	0.55
1:G:96:ARG:HA	1:G:96:ARG:NE	2.21	0.55
1:D:85:MSE:HG3	1:E:88:TYR:CE2	2.42	0.55
1:F:95:GLU:O	1:F:98:GLY:N	2.40	0.55
1:F:97:MSE:CE	1:F:134:VAL:CG2	2.79	0.55
1:A:116:PRO:O	1:A:117:LEU:HB3	2.07	0.55
1:E:41:GLN:NE2	1:E:116:PRO:CD	2.70	0.55
1:E:83:MSE:SE	1:E:141:PHE:CE2	3.10	0.55
1:D:118:LYS:NZ	1:D:169:GLU:OE2	2.37	0.55
1:E:141:PHE:C	1:E:141:PHE:CD2	2.80	0.55
1:B:107:ARG:NH1	1:B:186:HIS:HD2	1.80	0.55
1:A:122:ASN:CB	1:A:166:LEU:HD12	2.36	0.55
1:D:9:LEU:HD12	1:D:9:LEU:O	2.07	0.54
1:F:111:LEU:HD11	1:F:187:LEU:CD2	2.36	0.54
1:E:111:LEU:CD2	1:E:186:HIS:CB	2.84	0.54
1:G:133:MSE:O	1:G:137:SER:HB3	2.07	0.54
1:E:83:MSE:HE2	1:E:144:GLN:NE2	2.22	0.54
1:E:168:ARG:O	1:E:171:MSE:HB3	2.07	0.54
1:F:210:TYR:C	1:F:210:TYR:HD2	2.10	0.54
1:G:200:HIS:O	1:G:203:ASN:CB	2.50	0.54
1:D:148:LEU:O	1:D:152:VAL:HG23	2.07	0.53
1:A:62:ASP:O	1:A:66:ARG:HG3	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:12:GLU:OE1	1:G:69:ARG:HD3	2.09	0.53
1:G:28:GLN:OE1	1:G:135:ASN:HB2	2.08	0.53
1:F:97:MSE:CE	1:F:134:VAL:HG13	2.38	0.53
1:G:126:MSE:HG3	1:G:162:LEU:HD23	1.91	0.53
1:A:23:MSE:HE2	1:A:94:LEU:HB3	1.89	0.53
1:D:78:ALA:O	1:D:82:ARG:HD3	2.09	0.53
1:A:122:ASN:HB3	1:A:166:LEU:CD1	2.39	0.53
1:A:196:ARG:HG3	1:B:196:ARG:NH2	2.23	0.53
1:F:163:TYR:HD2	1:F:164:HIS:CD2	2.26	0.53
1:G:103:ASN:HB3	1:G:193:HIS:NE2	2.23	0.53
1:A:175:LEU:HG	1:B:181:ILE:HD12	1.91	0.53
1:D:36:GLU:O	1:D:37:ALA:C	2.47	0.53
1:D:192:ARG:HE	1:E:192:ARG:NH1	2.06	0.53
1:B:64:GLU:HG2	1:B:88:TYR:CE1	2.44	0.53
1:F:130:VAL:O	1:F:134:VAL:HG23	2.09	0.53
1:A:196:ARG:HH21	1:B:196:ARG:HG3	1.74	0.52
1:F:13:LEU:N	1:F:70:MSE:HE3	2.24	0.52
1:A:142:ILE:HG22	1:A:143:GLN:HE21	1.74	0.52
1:B:178:PRO:O	1:B:181:ILE:CG1	2.54	0.52
1:D:118:LYS:HB3	1:D:173:TYR:CZ	2.45	0.52
1:E:172:THR:HG21	1:F:168:ARG:HH11	1.71	0.52
1:G:71:ILE:HD11	1:G:85:MSE:HE3	1.90	0.52
1:E:13:LEU:HB2	1:E:70:MSE:HE1	1.92	0.52
1:F:120:TYR:O	1:F:124:ASN:ND2	2.42	0.52
1:A:122:ASN:HB3	1:A:166:LEU:HD12	1.92	0.52
1:E:111:LEU:CD2	1:E:186:HIS:HB2	2.40	0.52
1:E:130:VAL:O	1:E:134:VAL:HG23	2.10	0.52
1:F:96:ARG:HA	1:F:96:ARG:NE	2.25	0.52
1:F:174:VAL:CG1	1:G:175:LEU:HD21	2.40	0.52
1:A:185:MSE:SE	1:B:167:GLU:OE2	2.78	0.52
1:E:111:LEU:CD2	1:E:186:HIS:HB3	2.40	0.52
1:G:23:MSE:SE	1:G:91:VAL:HG22	2.60	0.52
1:B:122:ASN:HB2	1:B:166:LEU:CD1	2.39	0.52
1:F:42:ASN:ND2	1:F:45:LEU:HB2	2.25	0.52
1:F:179:ARG:C	1:F:181:ILE:H	2.14	0.52
1:B:83:MSE:HE2	1:B:144:GLN:HE22	1.74	0.51
2:B:249:HOH:O	1:E:182:LYS:HD3	2.10	0.51
1:E:146:THR:HG23	1:E:209:ILE:HD12	1.92	0.51
1:D:97:MSE:HE3	1:D:134:VAL:HG22	1.84	0.51
1:B:36:GLU:HB3	1:B:45:LEU:HD21	1.90	0.51
1:D:182:LYS:HE3	1:D:186:HIS:HE2	1.75	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:13:LEU:N	1:F:70:MSE:CE	2.73	0.51
1:A:113:GLU:O	1:A:114:GLU:HG3	2.11	0.51
1:A:143:GLN:C	1:A:144:GLN:HG2	2.31	0.51
1:E:4:GLY:O	1:E:8:LYS:HB2	2.11	0.51
1:G:85:MSE:HE2	1:G:85:MSE:HA	1.93	0.51
1:G:210:TYR:O	1:G:211:LEU:C	2.49	0.51
1:B:161:GLU:HB3	1:D:121:VAL:HG21	1.93	0.51
1:B:142:ILE:CG2	1:B:143:GLN:HE21	2.24	0.50
1:D:103:ASN:HB3	1:D:193:HIS:NE2	2.26	0.50
1:E:168:ARG:HG2	1:E:168:ARG:HH11	1.76	0.50
1:E:107:ARG:O	1:E:111:LEU:HB2	2.11	0.50
1:F:12:GLU:OE2	1:F:74:TYR:OH	2.24	0.50
1:G:149:ALA:HB3	1:G:209:ILE:HD11	1.94	0.50
1:A:65:ARG:HE	1:A:69:ARG:HH21	1.58	0.50
1:F:192:ARG:NH1	1:G:163:TYR:OH	2.45	0.50
1:D:38:LEU:HD11	1:D:111:LEU:HD13	1.93	0.50
1:E:43:VAL:HG22	1:E:109:ILE:HG12	1.93	0.50
1:A:43:VAL:HG13	1:A:109:ILE:CD1	2.39	0.49
1:B:21:ILE:HD11	1:B:142:ILE:HG12	1.92	0.49
1:B:23:MSE:HE1	1:B:91:VAL:O	2.12	0.49
1:D:181:ILE:HG12	1:E:171:MSE:SE	2.62	0.49
1:B:23:MSE:SE	1:B:91:VAL:CG2	3.04	0.49
1:D:188:SER:O	1:D:192:ARG:HG3	2.11	0.49
1:E:168:ARG:CG	1:F:172:THR:HG21	2.41	0.49
1:E:27:VAL:O	1:E:31:ILE:HG13	2.13	0.49
1:E:149:ALA:O	1:E:152:VAL:HB	2.13	0.49
1:D:29:GLU:HB3	1:D:33:LYS:NZ	2.27	0.49
1:A:97:MSE:HE1	1:A:134:VAL:HG22	1.88	0.49
1:B:164:HIS:HB3	2:B:233:HOH:O	2.12	0.49
1:B:97:MSE:CE	1:B:134:VAL:CG2	2.68	0.49
1:D:205:ALA:O	1:D:208:ALA:HB3	2.13	0.49
1:E:83:MSE:HE2	1:E:211:LEU:HD12	1.95	0.49
1:B:33:LYS:HB3	1:B:49:VAL:CG2	2.43	0.48
1:B:43:VAL:O	1:B:47:GLU:HG2	2.13	0.48
1:F:179:ARG:O	1:F:181:ILE:N	2.46	0.48
1:F:210:TYR:O	1:F:211:LEU:C	2.49	0.48
1:A:164:HIS:HB3	2:A:240:HOH:O	2.14	0.48
1:A:171:MSE:O	1:A:175:LEU:HB2	2.12	0.48
1:B:23:MSE:HA	1:B:59:LEU:HD23	1.95	0.48
1:E:28:GLN:O	1:E:29:GLU:C	2.49	0.48
1:E:163:TYR:CD2	1:E:163:TYR:C	2.85	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:103:ASN:HB3	1:G:193:HIS:CD2	2.49	0.48
1:E:168:ARG:HG2	1:F:172:THR:HG21	1.95	0.48
1:A:186:HIS:O	1:A:190:VAL:HG23	2.14	0.48
1:A:117:LEU:HB2	1:A:183:ARG:HD2	1.95	0.48
1:D:129:ILE:HG23	1:D:155:LYS:CD	2.44	0.48
1:G:97:MSE:CE	1:G:134:VAL:CG2	2.84	0.48
1:B:23:MSE:CA	1:B:59:LEU:HD23	2.43	0.48
1:D:120:TYR:O	1:D:124:ASN:ND2	2.47	0.48
1:F:60:GLU:HG3	1:F:91:VAL:HG13	1.96	0.48
1:A:37:ALA:HB2	1:A:45:LEU:HB3	1.96	0.47
1:B:38:LEU:HA	1:B:108:ALA:HB1	1.96	0.47
1:D:23:MSE:SE	1:D:91:VAL:CG2	3.13	0.47
1:A:120:TYR:CE2	1:A:187:LEU:HD11	2.49	0.47
1:B:183:ARG:O	1:B:184:ALA:C	2.52	0.47
1:A:17:LYS:CE	1:A:83:MSE:HE2	2.42	0.47
1:E:165:GLN:HG3	1:F:169:GLU:OE2	2.15	0.47
1:F:28:GLN:O	1:F:29:GLU:C	2.53	0.47
1:B:40:LYS:O	1:B:42:ASN:N	2.48	0.47
1:E:64:GLU:HG2	1:E:88:TYR:CE1	2.49	0.47
1:E:93:ASP:OD2	1:E:203:ASN:HB3	2.14	0.47
1:A:196:ARG:HG3	1:B:196:ARG:HH21	1.78	0.47
1:E:38:LEU:HD12	1:E:38:LEU:HA	1.71	0.47
1:F:13:LEU:HA	1:F:70:MSE:CE	2.23	0.47
1:B:209:ILE:O	1:B:210:TYR:C	2.53	0.46
1:A:196:ARG:HH21	1:B:196:ARG:NE	2.12	0.46
1:D:10:PHE:HE1	2:D:234:HOH:O	1.97	0.46
1:B:38:LEU:HD21	1:B:123:ILE:HG21	1.97	0.46
1:D:192:ARG:HH21	1:E:192:ARG:NH2	2.13	0.46
1:E:43:VAL:HG22	1:E:109:ILE:CG1	2.46	0.46
1:E:165:GLN:CD	1:F:122:ASN:OD1	2.54	0.46
1:D:17:LYS:O	1:D:21:ILE:HD12	2.15	0.46
1:A:118:LYS:NZ	1:A:169:GLU:CD	2.69	0.46
1:D:163:TYR:OH	1:D:192:ARG:HD3	2.15	0.46
1:F:64:GLU:O	1:F:67:CYS:HB2	2.15	0.46
1:G:163:TYR:HD1	1:G:191:ALA:HB1	1.81	0.46
1:G:178:PRO:HA	1:G:181:ILE:HG13	1.98	0.46
1:E:144:GLN:CA	1:E:144:GLN:OE1	2.63	0.46
1:F:32:ASP:OD2	1:F:131:LYS:CE	2.64	0.46
1:B:119:PRO:CG	1:D:161:GLU:HG2	2.45	0.46
1:D:83:MSE:SE	1:D:211:LEU:HD21	2.66	0.46
1:F:85:MSE:HE3	1:G:68:ILE:CG1	2.45	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:45:LEU:HA	1:E:45:LEU:HD12	1.74	0.46
1:E:83:MSE:O	1:E:87:ILE:HG13	2.16	0.46
1:D:163:TYR:OH	1:D:192:ARG:HG2	2.16	0.45
1:F:97:MSE:HE1	1:F:134:VAL:CG2	2.43	0.45
1:A:96:ARG:O	1:A:100:GLU:HG2	2.17	0.45
1:B:155:LYS:O	1:B:158:THR:HB	2.16	0.45
1:D:83:MSE:HA	1:D:211:LEU:HD11	1.97	0.45
1:E:65:ARG:O	1:E:66:ARG:C	2.55	0.45
1:F:38:LEU:HD11	1:F:111:LEU:HD12	1.98	0.45
1:F:94:LEU:HD23	1:F:94:LEU:HA	1.69	0.45
1:F:177:ASP:OD1	1:F:177:ASP:C	2.53	0.45
1:B:76:PRO:HB2	1:B:80:ASP:HB2	1.98	0.45
1:F:32:ASP:OD2	1:F:131:LYS:HE2	2.15	0.45
1:F:94:LEU:HD23	1:F:97:MSE:SE	2.67	0.45
1:G:141:PHE:CD2	1:G:141:PHE:C	2.89	0.45
1:G:176:GLU:HA	2:G:235:HOH:O	2.15	0.45
1:G:210:TYR:O	1:G:212:SER:N	2.50	0.45
1:B:155:LYS:HD3	1:B:155:LYS:HA	1.80	0.45
1:D:149:ALA:O	1:D:152:VAL:HB	2.16	0.45
1:F:60:GLU:HG3	1:F:91:VAL:CG1	2.47	0.45
1:G:118:LYS:HB3	1:G:173:TYR:CZ	2.51	0.45
1:A:153:ILE:HD11	1:A:205:ALA:HB3	1.97	0.45
1:D:129:ILE:HG23	1:D:155:LYS:HD3	1.97	0.45
1:F:12:GLU:C	1:F:70:MSE:HE3	2.37	0.45
1:G:42:ASN:OD1	1:G:44:GLU:HB3	2.16	0.45
1:G:83:MSE:HE2	1:G:144:GLN:HE21	1.80	0.45
1:E:59:LEU:HD23	1:E:59:LEU:HA	1.57	0.45
1:F:45:LEU:HD12	1:F:45:LEU:HA	1.66	0.45
1:D:103:ASN:HB3	1:D:193:HIS:CE1	2.52	0.45
1:D:114:GLU:HB3	1:D:183:ARG:HD3	1.99	0.45
1:E:212:SER:O	1:E:213:GLU:C	2.55	0.45
1:G:21:ILE:HD11	1:G:142:ILE:HG12	1.98	0.45
1:G:111:LEU:HA	1:G:111:LEU:HD23	1.65	0.45
1:B:23:MSE:HG3	1:B:59:LEU:HD23	1.98	0.45
1:B:23:MSE:HB2	1:B:59:LEU:HD23	1.97	0.44
1:B:171:MSE:HE2	1:B:171:MSE:HB3	1.97	0.44
1:E:143:GLN:C	1:E:144:GLN:HG2	2.37	0.44
1:E:58:LEU:O	1:E:58:LEU:HG	2.16	0.44
1:E:165:GLN:NE2	1:F:122:ASN:OD1	2.50	0.44
1:F:24:ALA:O	1:F:28:GLN:HG3	2.17	0.44
1:A:44:GLU:O	1:A:48:GLU:HG3	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:81:LEU:O	1:B:85:MSE:HG2	2.17	0.44
1:A:169:GLU:O	1:A:172:THR:HB	2.17	0.44
1:E:111:LEU:HD22	1:E:186:HIS:HB2	1.98	0.44
1:G:71:ILE:CD1	1:G:85:MSE:HE3	2.47	0.44
1:G:47:GLU:HA	1:G:50:ILE:HD12	2.00	0.44
1:D:111:LEU:HD22	2:D:230:HOH:O	2.18	0.44
1:B:63:ILE:HD13	1:B:91:VAL:HG21	1.99	0.44
1:D:126:MSE:HG3	1:D:162:LEU:HD23	2.00	0.44
1:F:163:TYR:CD2	1:F:163:TYR:C	2.91	0.44
1:G:167:GLU:HA	1:G:167:GLU:OE1	2.18	0.44
1:G:211:LEU:HD22	1:G:211:LEU:HA	1.83	0.44
1:A:64:GLU:HG2	1:A:88:TYR:CD1	2.53	0.44
1:A:142:ILE:HG22	1:A:143:GLN:HG2	2.00	0.44
1:A:142:ILE:CG2	1:A:143:GLN:HE21	2.31	0.44
1:D:118:LYS:CD	1:D:169:GLU:OE2	2.60	0.44
1:G:33:LYS:HD3	1:G:45:LEU:HD11	1.98	0.44
1:G:142:ILE:O	1:G:142:ILE:HG23	2.17	0.44
1:A:196:ARG:NH2	1:B:196:ARG:HG3	2.32	0.43
1:D:115:PRO:HA	1:D:116:PRO:HD3	1.88	0.43
1:D:206:GLU:O	1:D:207:ALA:C	2.56	0.43
1:E:200:HIS:O	1:E:201:ALA:C	2.57	0.43
1:G:142:ILE:HG22	1:G:143:GLN:HG2	2.00	0.43
1:A:45:LEU:HA	1:A:45:LEU:HD12	1.73	0.43
1:D:107:ARG:NH1	1:D:186:HIS:ND1	2.65	0.43
1:E:33:LYS:HB2	1:E:49:VAL:CG2	2.49	0.43
1:F:19:GLN:O	1:F:20:VAL:C	2.57	0.43
1:G:45:LEU:HD12	1:G:45:LEU:HA	1.80	0.43
1:A:123:ILE:HG23	1:A:194:TYR:CE1	2.52	0.43
1:E:127:SER:HB2	1:E:194:TYR:CE2	2.53	0.43
1:G:166:LEU:HD22	1:G:191:ALA:HB2	1.99	0.43
1:F:126:MSE:HE2	1:F:194:TYR:CB	2.49	0.43
1:D:33:LYS:HD3	1:D:45:LEU:HD11	2.01	0.43
1:B:9:LEU:HD12	1:B:70:MSE:HG2	2.01	0.43
1:D:95:GLU:OE2	1:D:96:ARG:HD2	2.18	0.43
1:D:96:ARG:HB3	1:D:200:HIS:NE2	2.33	0.43
1:E:162:LEU:HD12	1:E:162:LEU:HA	1.88	0.43
1:G:97:MSE:CE	1:G:134:VAL:HG13	2.49	0.43
1:B:60:GLU:HG3	1:B:91:VAL:HG13	2.01	0.43
1:D:35:THR:HG22	1:D:39:ASN:ND2	2.34	0.43
1:E:190:VAL:O	1:E:191:ALA:C	2.56	0.43
1:F:41:GLN:HE21	1:F:112:ALA:HA	1.84	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:154:GLU:C	1:F:156:ASP:H	2.23	0.43
1:B:117:LEU:HB2	1:B:183:ARG:HD2	2.00	0.42
1:E:211:LEU:HD22	1:E:211:LEU:HA	1.73	0.42
1:F:23:MSE:SE	1:F:91:VAL:HG22	2.69	0.42
1:A:47:GLU:O	1:A:51:LYS:HG2	2.20	0.42
1:E:111:LEU:HA	1:E:111:LEU:HD13	1.60	0.42
1:E:122:ASN:HB2	1:E:166:LEU:HD12	2.01	0.42
1:F:149:ALA:O	1:F:152:VAL:HB	2.19	0.42
1:A:148:LEU:O	1:A:152:VAL:HG23	2.19	0.42
1:E:63:ILE:HD12	1:E:91:VAL:HG21	2.00	0.42
1:G:106:GLU:HG3	2:G:246:HOH:O	2.20	0.42
1:G:122:ASN:CG	1:G:166:LEU:HD12	2.40	0.42
1:A:23:MSE:HE1	1:A:91:VAL:O	2.20	0.42
1:G:59:LEU:HD12	1:G:59:LEU:HA	1.79	0.42
1:F:118:LYS:HZ3	1:F:169:GLU:CD	2.22	0.42
1:G:170:LEU:HD23	1:G:170:LEU:HA	1.75	0.42
1:G:192:ARG:O	1:G:196:ARG:HG3	2.20	0.42
1:A:33:LYS:HB3	1:A:49:VAL:HG22	2.02	0.42
1:B:94:LEU:HD23	1:B:97:MSE:SE	2.70	0.42
1:D:96:ARG:HD2	1:D:96:ARG:HA	1.64	0.42
1:F:97:MSE:HG2	1:F:197:ILE:HG23	2.02	0.42
1:F:185:MSE:HE2	1:F:185:MSE:HB2	1.98	0.42
1:G:27:VAL:HG11	1:G:97:MSE:HE2	2.02	0.42
1:G:200:HIS:C	1:G:203:ASN:HB2	2.37	0.42
1:D:125:PHE:O	1:D:129:ILE:HG13	2.19	0.42
1:E:142:ILE:HA	1:E:142:ILE:HD12	1.74	0.42
1:F:13:LEU:HB2	1:F:70:MSE:CE	2.49	0.42
1:F:190:VAL:O	1:F:191:ALA:C	2.56	0.42
1:B:94:LEU:HD11	1:B:138:VAL:HG21	2.02	0.42
1:E:97:MSE:HE1	1:E:134:VAL:CG1	2.45	0.42
1:E:117:LEU:HD11	1:E:173:TYR:CG	2.55	0.42
1:E:174:VAL:HG12	1:E:175:LEU:N	2.34	0.42
1:F:10:PHE:CE1	1:F:14:GLU:OE2	2.73	0.42
1:B:58:LEU:HD12	1:B:61:VAL:HB	2.02	0.41
1:B:179:ARG:HG2	1:B:180:ASN:N	2.34	0.41
1:F:20:VAL:HG21	1:F:87:ILE:HG23	2.01	0.41
1:G:23:MSE:SE	1:G:91:VAL:CG2	3.18	0.41
1:A:116:PRO:O	1:A:117:LEU:CB	2.68	0.41
1:B:83:MSE:HE2	1:B:144:GLN:HE21	1.84	0.41
1:B:176:GLU:O	1:B:177:ASP:HB2	2.20	0.41
1:E:111:LEU:HD21	1:E:186:HIS:HB2	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:8:LYS:O	1:G:12:GLU:HB2	2.20	0.41
1:G:151:LYS:O	1:G:152:VAL:C	2.59	0.41
1:A:17:LYS:O	1:A:21:ILE:HD12	2.21	0.41
1:A:118:LYS:NZ	1:A:169:GLU:OE2	2.50	0.41
1:A:192:ARG:NH1	1:B:192:ARG:CZ	2.83	0.41
1:D:68:ILE:CG1	1:E:85:MSE:CE	2.97	0.41
1:E:206:GLU:O	1:E:207:ALA:C	2.59	0.41
1:E:210:TYR:HD2	1:E:211:LEU:N	2.19	0.41
1:F:170:LEU:HD23	1:F:170:LEU:HA	1.75	0.41
1:E:125:PHE:O	1:E:129:ILE:HG13	2.20	0.41
1:F:21:ILE:O	1:F:22:LYS:C	2.59	0.41
1:F:81:LEU:HD13	1:F:81:LEU:HA	1.92	0.41
1:E:13:LEU:HB2	1:E:70:MSE:CE	2.50	0.41
1:G:118:LYS:HD2	1:G:169:GLU:OE1	2.20	0.41
1:A:185:MSE:O	1:A:188:SER:HB3	2.20	0.41
1:B:176:GLU:OE2	1:D:168:ARG:NH2	2.53	0.41
1:F:163:TYR:HH	1:G:192:ARG:HH21	1.59	0.41
1:B:90:ILE:HG13	1:B:91:VAL:N	2.35	0.41
1:B:94:LEU:HD23	1:B:94:LEU:HA	1.64	0.41
1:E:207:ALA:O	1:E:208:ALA:C	2.58	0.41
1:F:13:LEU:CB	1:F:70:MSE:CE	2.98	0.41
1:F:125:PHE:O	1:F:129:ILE:HG13	2.20	0.41
1:A:34:ALA:HB1	1:A:104:ILE:HG22	2.03	0.41
1:A:196:ARG:NH2	1:B:196:ARG:HE	2.17	0.41
1:E:168:ARG:HH11	1:E:168:ARG:CG	2.34	0.41
1:F:9:LEU:HD12	1:F:9:LEU:O	2.21	0.41
1:F:209:ILE:HD13	1:F:209:ILE:HA	1.87	0.41
1:G:130:VAL:HG12	1:G:131:LYS:N	2.35	0.41
1:B:70:MSE:HE2	1:B:70:MSE:HB2	1.97	0.41
1:B:97:MSE:HE1	1:B:134:VAL:CG2	2.40	0.41
1:B:117:LEU:O	1:B:118:LYS:CB	2.60	0.41
1:E:83:MSE:HE2	1:E:211:LEU:CD1	2.51	0.41
1:E:150:LYS:HA	1:E:153:ILE:HD12	2.03	0.41
1:G:62:ASP:C	1:G:62:ASP:OD1	2.58	0.41
1:D:68:ILE:CG1	1:E:85:MSE:HE3	2.51	0.41
1:A:63:ILE:CB	1:A:91:VAL:HG21	2.49	0.40
1:B:143:GLN:C	1:B:144:GLN:HG2	2.40	0.40
1:E:64:GLU:O	1:E:67:CYS:HB2	2.21	0.40
1:E:144:GLN:OE1	1:E:144:GLN:HA	2.21	0.40
1:B:23:MSE:HG3	1:B:59:LEU:HB3	2.03	0.40
1:A:77:GLU:H	1:A:77:GLU:HG2	1.55	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:23:MSE:CG	1:B:59:LEU:HD23	2.51	0.40
1:E:36:GLU:HB3	1:E:45:LEU:HD23	2.03	0.40
1:F:189:PHE:O	1:F:190:VAL:C	2.58	0.40
1:E:13:LEU:CA	1:E:70:MSE:CE	2.97	0.40
1:E:83:MSE:SE	1:E:141:PHE:HE2	2.55	0.40
1:F:135:ASN:O	1:F:139:ILE:HG13	2.21	0.40
1:A:30:ALA:O	1:A:31:ILE:C	2.57	0.40
1:A:112:ALA:C	1:A:114:GLU:H	2.24	0.40
1:B:171:MSE:SE	2:B:252:HOH:O	2.90	0.40
1:F:36:GLU:HB3	1:F:45:LEU:HD23	2.04	0.40

All (1) 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:B:179:ARG:NH1	1:D:179:ARG:NH1[4_565]	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	213/227 (94%)	189 (89%)	21 (10%)	3 (1%)	11	36
1	B	212/227 (93%)	185 (87%)	24 (11%)	3 (1%)	11	36
1	D	213/227 (94%)	186 (87%)	24 (11%)	3 (1%)	11	36
1	E	211/227 (93%)	191 (90%)	18 (8%)	2 (1%)	17	48
1	F	212/227 (93%)	191 (90%)	19 (9%)	2 (1%)	17	48
1	G	209/227 (92%)	184 (88%)	21 (10%)	4 (2%)	8	28
All	All	1270/1362 (93%)	1126 (89%)	127 (10%)	17 (1%)	12	37

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	78	ALA
1	A	117	LEU
1	B	41	GLN
1	D	180	ASN
1	E	176	GLU
1	F	180	ASN
1	G	144	GLN
1	G	180	ASN
1	A	120	TYR
1	G	78	ALA
1	B	39	ASN
1	D	214	GLY
1	B	190	VAL
1	E	212	SER
1	G	211	LEU
1	D	3	GLY
1	F	114	GLU

5.3.2 Protein sidechains (i)

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	185/186 (100%)	150 (81%)	35 (19%)	1 4
1	B	185/186 (100%)	152 (82%)	33 (18%)	2 5
1	D	185/186 (100%)	156 (84%)	29 (16%)	2 8
1	E	184/186 (99%)	150 (82%)	34 (18%)	1 5
1	F	185/186 (100%)	155 (84%)	30 (16%)	2 7
1	G	184/186 (99%)	151 (82%)	33 (18%)	2 5
All	All	1108/1116 (99%)	914 (82%)	194 (18%)	2 6

All (194) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	15	GLU
1	A	17	LYS

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Mol	Chain	Res	Type
1	A	29	GLU
1	A	44	GLU
1	A	47	GLU
1	A	51	LYS
1	A	54	ASP
1	A	59	LEU
1	A	63	ILE
1	A	69	ARG
1	A	70	MSE
1	A	77	GLU
1	A	89	LYS
1	A	91	VAL
1	A	106	GLU
1	A	109	ILE
1	A	118	LYS
1	A	121	VAL
1	A	126	MSE
1	A	128	GLU
1	A	130	VAL
1	A	131	LYS
1	A	142	ILE
1	A	143	GLN
1	A	144	GLN
1	A	147	LEU
1	A	161	GLU
1	A	162	LEU
1	A	165	GLN
1	A	166	LEU
1	A	176	GLU
1	A	181	ILE
1	A	196	ARG
1	A	209	ILE
1	A	215	GLU
1	B	7	MSE
1	B	12	GLU
1	B	25	LYS
1	B	29	GLU
1	B	32	ASP
1	B	47	GLU
1	B	56	ILE
1	B	62	ASP
1	B	69	ARG

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Mol	Chain	Res	Type
1	B	70	MSE
1	B	77	GLU
1	B	83	MSE
1	B	89	LYS
1	B	91	VAL
1	B	109	ILE
1	B	113	GLU
1	B	116	PRO
1	B	118	LYS
1	B	142	ILE
1	B	143	GLN
1	B	147	LEU
1	B	155	LYS
1	B	162	LEU
1	B	166	LEU
1	B	168	ARG
1	B	176	GLU
1	B	179	ARG
1	B	185	MSE
1	B	192	ARG
1	B	196	ARG
1	B	202	GLU
1	B	206	GLU
1	B	209	ILE
1	D	11	LYS
1	D	15	GLU
1	D	17	LYS
1	D	25	LYS
1	D	33	LYS
1	D	45	LEU
1	D	47	GLU
1	D	59	LEU
1	D	83	MSE
1	D	90	ILE
1	D	91	VAL
1	D	95	GLU
1	D	96	ARG
1	D	110	LEU
1	D	113	GLU
1	D	118	LYS
1	D	128	GLU
1	D	129	ILE

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Mol	Chain	Res	Type
1	D	142	ILE
1	D	143	GLN
1	D	151	LYS
1	D	162	LEU
1	D	165	GLN
1	D	166	LEU
1	D	171	MSE
1	D	172	THR
1	D	202	GLU
1	D	206	GLU
1	D	212	SER
1	E	8	LYS
1	E	12	GLU
1	E	15	GLU
1	E	22	LYS
1	E	25	LYS
1	E	32	ASP
1	E	42	ASN
1	E	47	GLU
1	E	65	ARG
1	E	69	ARG
1	E	83	MSE
1	E	85	MSE
1	E	89	LYS
1	E	90	ILE
1	E	91	VAL
1	E	111	LEU
1	E	113	GLU
1	E	114	GLU
1	E	118	LYS
1	E	121	VAL
1	E	123	ILE
1	E	127	SER
1	E	142	ILE
1	E	162	LEU
1	E	163	TYR
1	E	168	ARG
1	E	172	THR
1	E	179	ARG
1	E	181	ILE
1	E	182	LYS
1	E	185	MSE

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Mol	Chain	Res	Type
1	E	206	GLU
1	E	210	TYR
1	E	211	LEU
1	F	7	MSE
1	F	8	LYS
1	F	14	GLU
1	F	15	GLU
1	F	25	LYS
1	F	32	ASP
1	F	47	GLU
1	F	65	ARG
1	F	69	ARG
1	F	81	LEU
1	F	82	ARG
1	F	83	MSE
1	F	85	MSE
1	F	89	LYS
1	F	91	VAL
1	F	109	ILE
1	F	113	GLU
1	F	121	VAL
1	F	123	ILE
1	F	142	ILE
1	F	145	ASP
1	F	147	LEU
1	F	151	LYS
1	F	153	ILE
1	F	155	LYS
1	F	162	LEU
1	F	179	ARG
1	F	210	TYR
1	F	211	LEU
1	F	212	SER
1	G	8	LYS
1	G	11	LYS
1	G	12	GLU
1	G	15	GLU
1	G	22	LYS
1	G	25	LYS
1	G	32	ASP
1	G	33	LYS
1	G	40	LYS

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Mol	Chain	Res	Type
1	G	51	LYS
1	G	58	LEU
1	G	59	LEU
1	G	65	ARG
1	G	83	MSE
1	G	90	ILE
1	G	91	VAL
1	G	113	GLU
1	G	118	LYS
1	G	123	ILE
1	G	126	MSE
1	G	128	GLU
1	G	130	VAL
1	G	131	LYS
1	G	137	SER
1	G	142	ILE
1	G	154	GLU
1	G	155	LYS
1	G	162	LEU
1	G	166	LEU
1	G	169	GLU
1	G	172	THR
1	G	206	GLU
1	G	211	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	143	GLN
1	A	144	GLN
1	A	186	HIS
1	A	200	HIS
1	B	39	ASN
1	B	41	GLN
1	B	124	ASN
1	B	143	GLN
1	B	186	HIS
1	D	75	GLN
1	E	41	GLN
1	E	164	HIS
1	E	165	GLN
1	F	39	ASN

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Mol	Chain	Res	Type
1	F	41	GLN
1	F	124	ASN
1	F	164	HIS
1	F	165	GLN
1	G	41	GLN
1	G	144	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

There are no ligands in this entry.

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	205/227 (90%)	0.31	7 (3%)	45	40	9, 24, 38, 53	0
1	B	204/227 (89%)	0.28	9 (4%)	34	30	11, 24, 38, 61	0
1	D	205/227 (90%)	0.26	4 (1%)	65	63	6, 24, 41, 63	0
1	E	203/227 (89%)	0.25	2 (0%)	82	82	10, 23, 36, 56	0
1	F	204/227 (89%)	0.27	0	100	100	10, 23, 37, 55	0
1	G	201/227 (88%)	0.21	2 (0%)	82	82	5, 25, 35, 49	0
All	All	1222/1362 (89%)	0.26	24 (1%)	65	63	5, 24, 38, 63	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	122	ASN	4.2
1	B	121	VAL	3.3
1	B	111	LEU	3.0
1	A	147	LEU	2.9
1	G	58	LEU	2.9
1	D	3	GLY	2.8
1	B	56	ILE	2.8
1	E	1	GLY	2.7
1	D	149	ALA	2.6
1	D	153	ILE	2.5
1	A	38	LEU	2.5
1	A	129	ILE	2.5
1	B	108	ALA	2.4
1	D	2	GLY	2.3
1	A	122	ASN	2.3
1	B	214	GLY	2.2
1	G	153	ILE	2.2
1	A	144	GLN	2.1
1	A	111	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	147	LEU	2.1
1	B	215	GLU	2.1
1	A	56	ILE	2.1
1	B	125	PHE	2.1
1	E	210	TYR	2.0

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

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

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

6.4 Ligands (i)

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.