



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 2, 2026 – 02:58 AM UTC

PDB ID : 3TUZ / pdb_00003tuz
Title : Inward facing conformations of the MetNI methionine ABC transporter: CY5
SeMet soak crystal form
Authors : Johnson, E.; Nguyen, P.; Rees, D.C.
Deposited on : 2011-09-19
Resolution : 3.40 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

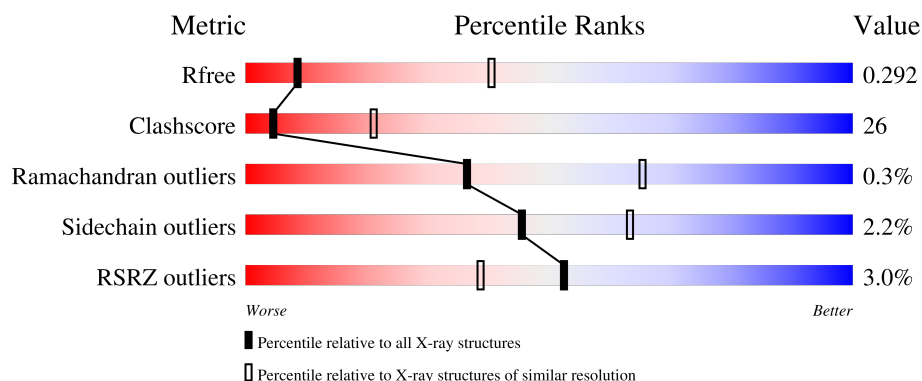
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.40 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1001 (3.44-3.36)
Clashscore	190562	1022 (3.44-3.36)
Ramachandran outliers	187476	1012 (3.44-3.36)
Sidechain outliers	187428	1012 (3.44-3.36)
RSRZ outliers	180081	1001 (3.44-3.36)

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

Mol	Chain	Length	Quality of chain
1	A	217	
1	B	217	
1	E	217	
1	F	217	
2	C	366	

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Mol	Chain	Length	Quality of chain
2	D	366	
2	G	366	
2	H	366	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	MSE	H	1000	-	-	X	-
4	ADP	C	2000	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 17234 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 D-methionine transport system permease protein metI.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	216	Total	C	N	O	S	0	0	0
			1627	1079	266	272	10			
1	B	216	Total	C	N	O	S	0	0	0
			1627	1079	266	272	10			
1	E	215	Total	C	N	O	S	0	0	0
			1616	1073	262	271	10			
1	F	216	Total	C	N	O	S	0	0	0
			1627	1079	266	272	10			

- Molecule 2 is a protein called Methionine import ATP-binding protein MetN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	344	Total	C	N	O	S	0	0	0
			2661	1670	471	508	12			
2	D	343	Total	C	N	O	S	0	0	0
			2651	1664	468	507	12			
2	G	343	Total	C	N	O	S	0	0	0
			2651	1664	468	507	12			
2	H	340	Total	C	N	O	S	0	0	0
			2630	1648	467	503	12			

There are 96 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	-22	MET	-	expression tag	UNP P30750
C	-21	GLY	-	expression tag	UNP P30750
C	-20	HIS	-	expression tag	UNP P30750
C	-19	HIS	-	expression tag	UNP P30750
C	-18	HIS	-	expression tag	UNP P30750
C	-17	HIS	-	expression tag	UNP P30750
C	-16	HIS	-	expression tag	UNP P30750
C	-15	HIS	-	expression tag	UNP P30750

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-14	HIS	-	expression tag	UNP P30750
C	-13	HIS	-	expression tag	UNP P30750
C	-12	HIS	-	expression tag	UNP P30750
C	-11	HIS	-	expression tag	UNP P30750
C	-10	SER	-	expression tag	UNP P30750
C	-9	SER	-	expression tag	UNP P30750
C	-8	GLY	-	expression tag	UNP P30750
C	-7	HIS	-	expression tag	UNP P30750
C	-6	ILE	-	expression tag	UNP P30750
C	-5	ASP	-	expression tag	UNP P30750
C	-4	ASP	-	expression tag	UNP P30750
C	-3	ASP	-	expression tag	UNP P30750
C	-2	ASP	-	expression tag	UNP P30750
C	-1	LYS	-	expression tag	UNP P30750
C	0	HIS	-	expression tag	UNP P30750
C	166	GLN	GLU	engineered mutation	UNP P30750
D	-22	MET	-	expression tag	UNP P30750
D	-21	GLY	-	expression tag	UNP P30750
D	-20	HIS	-	expression tag	UNP P30750
D	-19	HIS	-	expression tag	UNP P30750
D	-18	HIS	-	expression tag	UNP P30750
D	-17	HIS	-	expression tag	UNP P30750
D	-16	HIS	-	expression tag	UNP P30750
D	-15	HIS	-	expression tag	UNP P30750
D	-14	HIS	-	expression tag	UNP P30750
D	-13	HIS	-	expression tag	UNP P30750
D	-12	HIS	-	expression tag	UNP P30750
D	-11	HIS	-	expression tag	UNP P30750
D	-10	SER	-	expression tag	UNP P30750
D	-9	SER	-	expression tag	UNP P30750
D	-8	GLY	-	expression tag	UNP P30750
D	-7	HIS	-	expression tag	UNP P30750
D	-6	ILE	-	expression tag	UNP P30750
D	-5	ASP	-	expression tag	UNP P30750
D	-4	ASP	-	expression tag	UNP P30750
D	-3	ASP	-	expression tag	UNP P30750
D	-2	ASP	-	expression tag	UNP P30750
D	-1	LYS	-	expression tag	UNP P30750
D	0	HIS	-	expression tag	UNP P30750
D	166	GLN	GLU	engineered mutation	UNP P30750
G	-22	MET	-	expression tag	UNP P30750
G	-21	GLY	-	expression tag	UNP P30750

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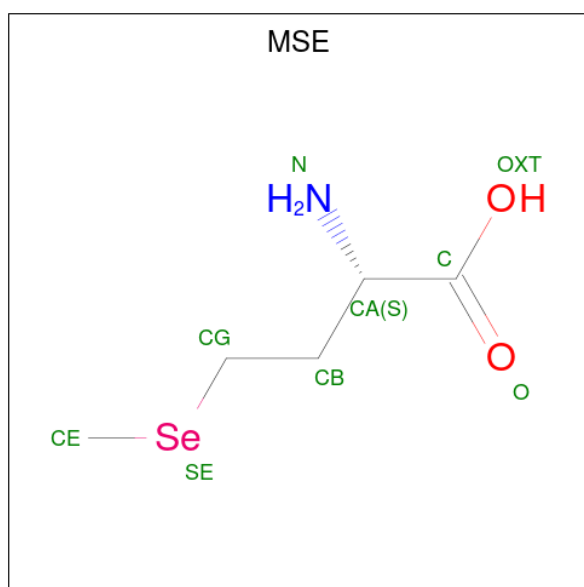
Chain	Residue	Modelled	Actual	Comment	Reference
G	-20	HIS	-	expression tag	UNP P30750
G	-19	HIS	-	expression tag	UNP P30750
G	-18	HIS	-	expression tag	UNP P30750
G	-17	HIS	-	expression tag	UNP P30750
G	-16	HIS	-	expression tag	UNP P30750
G	-15	HIS	-	expression tag	UNP P30750
G	-14	HIS	-	expression tag	UNP P30750
G	-13	HIS	-	expression tag	UNP P30750
G	-12	HIS	-	expression tag	UNP P30750
G	-11	HIS	-	expression tag	UNP P30750
G	-10	SER	-	expression tag	UNP P30750
G	-9	SER	-	expression tag	UNP P30750
G	-8	GLY	-	expression tag	UNP P30750
G	-7	HIS	-	expression tag	UNP P30750
G	-6	ILE	-	expression tag	UNP P30750
G	-5	ASP	-	expression tag	UNP P30750
G	-4	ASP	-	expression tag	UNP P30750
G	-3	ASP	-	expression tag	UNP P30750
G	-2	ASP	-	expression tag	UNP P30750
G	-1	LYS	-	expression tag	UNP P30750
G	0	HIS	-	expression tag	UNP P30750
G	166	GLN	GLU	engineered mutation	UNP P30750
H	-22	MET	-	expression tag	UNP P30750
H	-21	GLY	-	expression tag	UNP P30750
H	-20	HIS	-	expression tag	UNP P30750
H	-19	HIS	-	expression tag	UNP P30750
H	-18	HIS	-	expression tag	UNP P30750
H	-17	HIS	-	expression tag	UNP P30750
H	-16	HIS	-	expression tag	UNP P30750
H	-15	HIS	-	expression tag	UNP P30750
H	-14	HIS	-	expression tag	UNP P30750
H	-13	HIS	-	expression tag	UNP P30750
H	-12	HIS	-	expression tag	UNP P30750
H	-11	HIS	-	expression tag	UNP P30750
H	-10	SER	-	expression tag	UNP P30750
H	-9	SER	-	expression tag	UNP P30750
H	-8	GLY	-	expression tag	UNP P30750
H	-7	HIS	-	expression tag	UNP P30750
H	-6	ILE	-	expression tag	UNP P30750
H	-5	ASP	-	expression tag	UNP P30750
H	-4	ASP	-	expression tag	UNP P30750
H	-3	ASP	-	expression tag	UNP P30750

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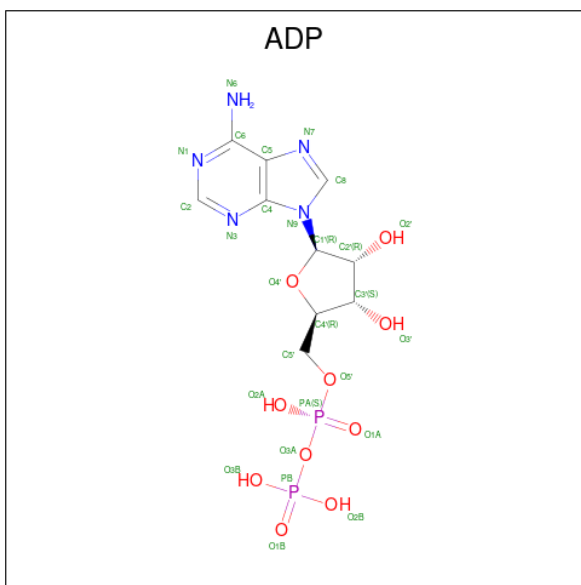
Chain	Residue	Modelled	Actual	Comment	Reference
H	-2	ASP	-	expression tag	UNP P30750
H	-1	LYS	-	expression tag	UNP P30750
H	0	HIS	-	expression tag	UNP P30750
H	166	GLN	GLU	engineered mutation	UNP P30750

- Molecule 3 is SELENOMETHIONINE (CCD ID: MSE) (formula: $C_5H_{11}NO_2Se$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	C	1	Total	C	N	O	Se	0	0
			9	5	1	2	1		
3	D	1	Total	C	N	O	Se	0	0
			9	5	1	2	1		
3	G	1	Total	C	N	O	Se	0	0
			9	5	1	2	1		
3	H	1	Total	C	N	O	Se	0	0
			9	5	1	2	1		

- Molecule 4 is ADENOSINE-5'-DIPHOSPHATE (CCD ID: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).

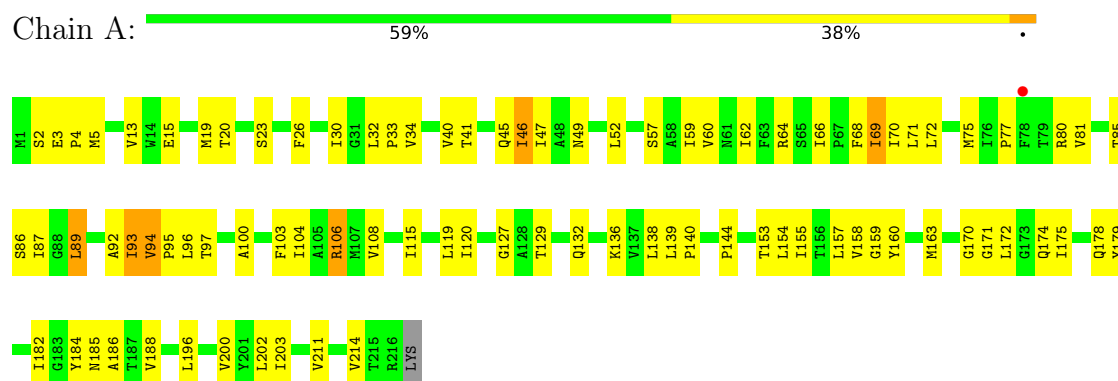


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	C	1	Total 27	C 10	N 5	O 10	P 2	0	0
4	D	1	Total 27	C 10	N 5	O 10	P 2	0	0
4	G	1	Total 27	C 10	N 5	O 10	P 2	0	0
4	H	1	Total 27	C 10	N 5	O 10	P 2	0	0

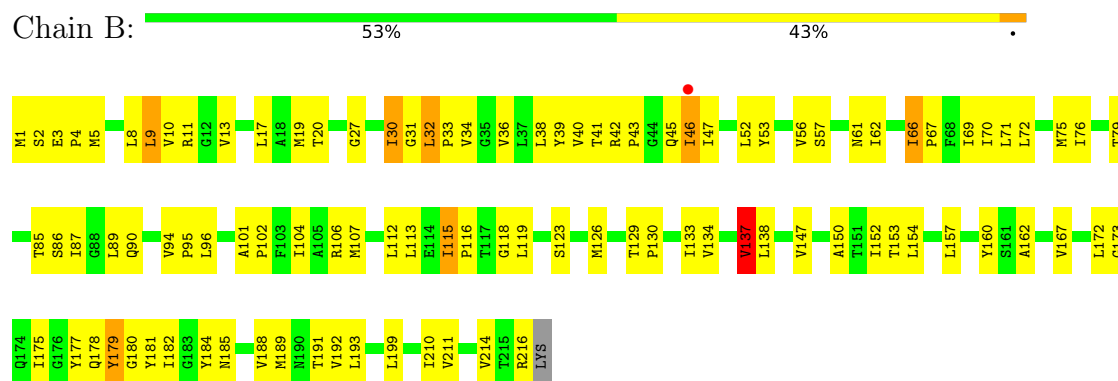
3 Residue-property plots

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

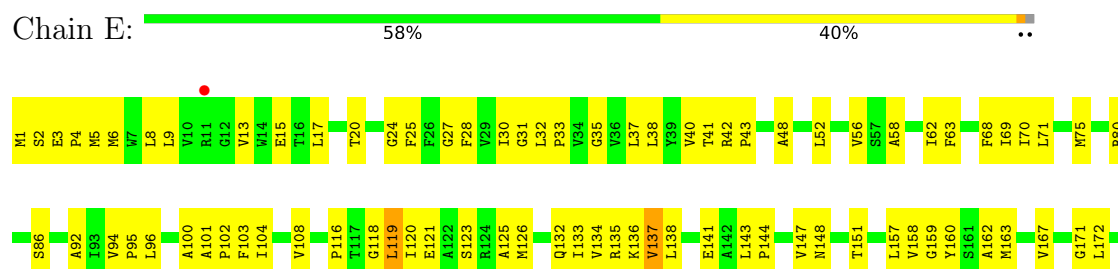
- Molecule 1: D-methionine transport system permease protein metI



- Molecule 1: D-methionine transport system permease protein metI

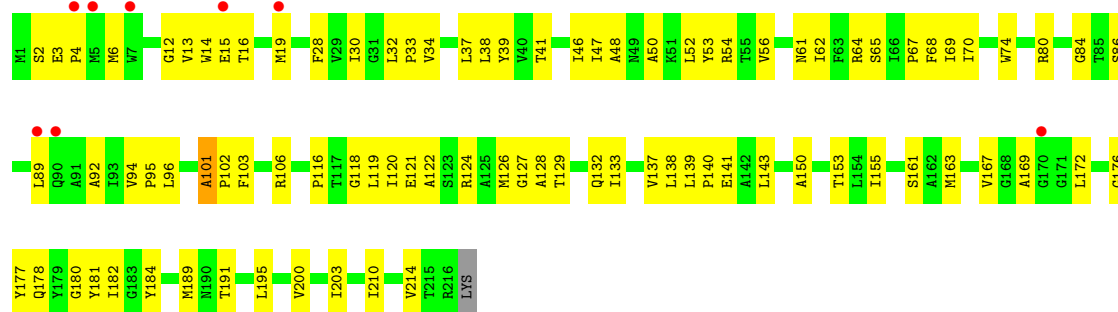


- Molecule 1: D-methionine transport system permease protein metI

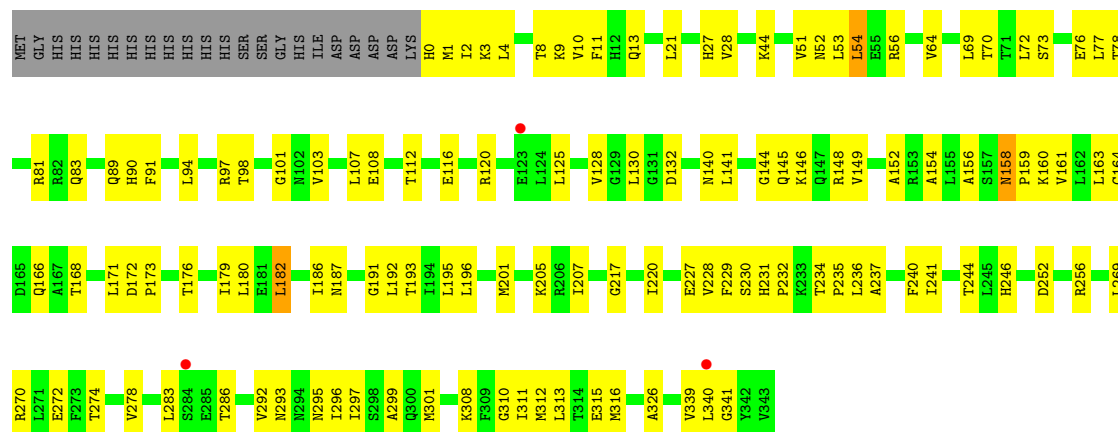




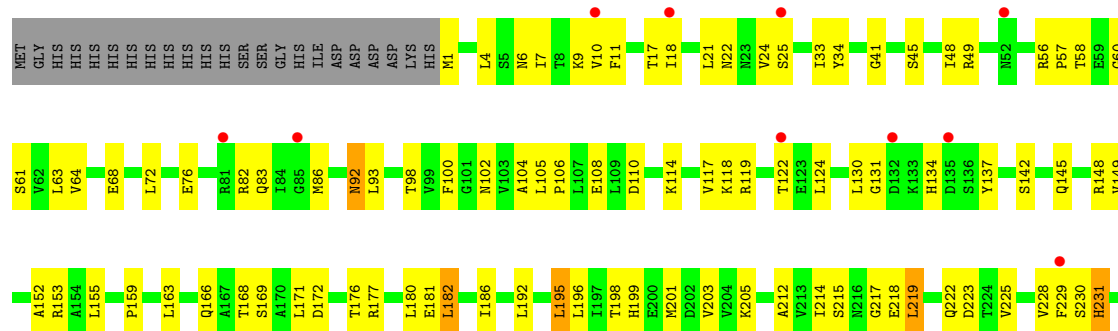
- Molecule 1: D-methionine transport system permease protein metI

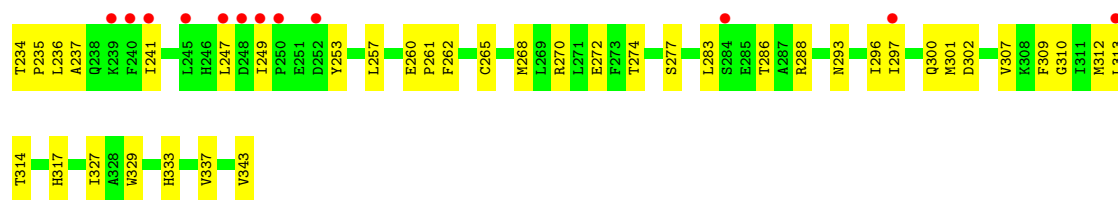


- Molecule 2: Methionine import ATP-binding protein MetN

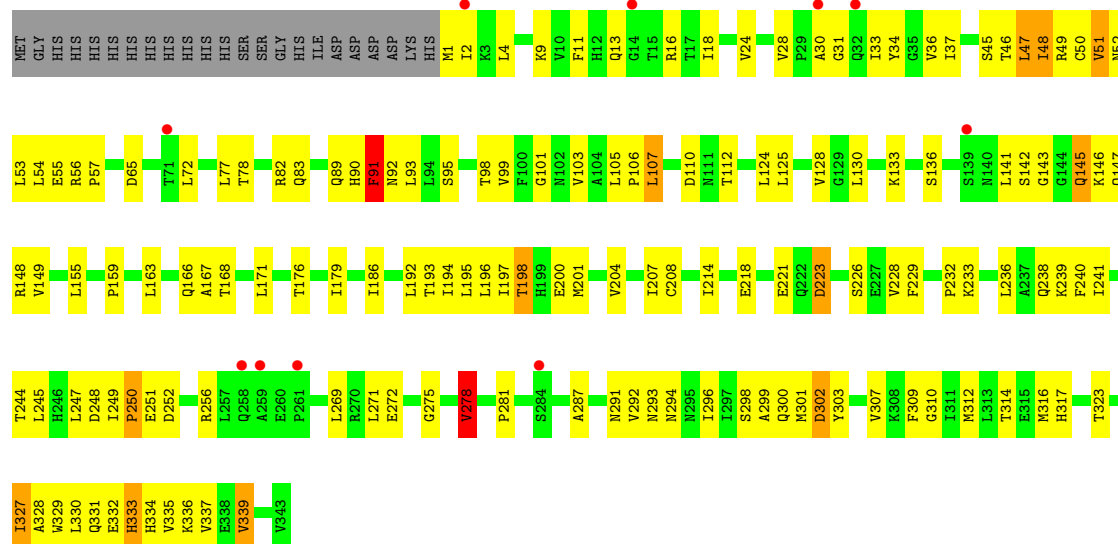


- Molecule 2: Methionine import ATP-binding protein MetN

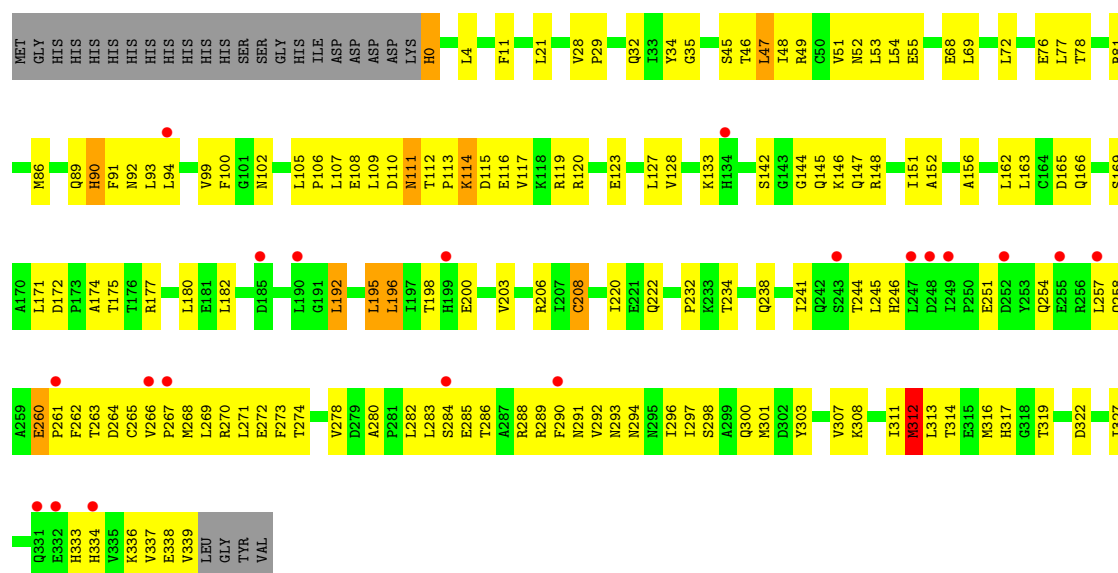




• Molecule 2: Methionine import ATP-binding protein MetN



• Molecule 2: Methionine import ATP-binding protein MetN



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	83.31Å 138.89Å 147.51Å 90.00° 95.72° 90.00°	Depositor
Resolution (Å)	38.80 – 3.40 38.80 – 3.40	Depositor EDS
% Data completeness (in resolution range)	93.4 (38.80-3.40) 93.3 (38.80-3.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.00 (at 3.18Å)	Xtriage
Refinement program	PHENIX 1.7.1 _743	Depositor
R, R_{free}	0.260 , 0.305 0.250 , 0.292	Depositor DCC
R_{free} test set	2337 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	58.8	Xtriage
Anisotropy	0.033	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 28.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.87	EDS
Total number of atoms	17234	wwPDB-VP
Average B, all atoms (Å ²)	80.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.25% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

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

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	0.78	0/1660	0.99	0/2267
1	B	0.81	1/1660 (0.1%)	1.13	8/2267 (0.4%)
1	E	0.73	1/1649 (0.1%)	1.01	4/2253 (0.2%)
1	F	0.76	0/1660	1.01	2/2267 (0.1%)
2	C	0.69	0/2701	0.97	4/3662 (0.1%)
2	D	0.69	0/2690	1.03	6/3647 (0.2%)
2	G	0.78	0/2690	1.08	11/3647 (0.3%)
2	H	0.74	0/2669	1.09	5/3618 (0.1%)
All	All	0.74	2/17379 (0.0%)	1.04	40/23628 (0.2%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	1	MET	SD-CE	-5.52	1.65	1.79
1	E	1	MET	CG-SD	-5.46	1.67	1.80

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	115	ILE	N-CA-C	8.55	115.98	108.63
1	E	119	LEU	N-CA-C	-7.50	103.08	111.71
2	G	333	HIS	CA-C-N	-7.41	111.25	122.21
2	G	333	HIS	C-N-CA	-7.41	111.25	122.21
2	H	90	HIS	N-CA-C	-7.39	98.90	110.36
1	B	46	ILE	N-CA-C	7.02	117.13	110.53
2	D	231	HIS	CA-C-N	6.60	126.96	119.83
2	D	231	HIS	C-N-CA	6.60	126.96	119.83
2	G	278	VAL	N-CA-C	6.46	116.61	110.53
2	D	92	ASN	N-CA-C	6.08	119.63	111.24
2	C	90	HIS	N-CA-C	-6.04	102.14	110.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	208	CYS	N-CA-C	6.04	119.36	109.76
2	C	54	LEU	N-CA-C	-5.96	105.27	112.54
2	G	167	ALA	N-CA-C	5.92	117.73	111.28
2	G	302	ASP	N-CA-C	5.91	117.25	107.73
2	G	223	ASP	N-CA-C	5.87	115.99	108.24
1	F	101	ALA	CA-C-N	5.85	125.22	118.85
1	F	101	ALA	C-N-CA	5.85	125.22	118.85
2	G	51	VAL	CB-CA-C	-5.72	104.42	112.14
1	B	137	VAL	N-CA-C	5.66	116.43	111.56
2	G	207	ILE	N-CA-C	5.65	118.62	113.20
2	D	131	GLY	N-CA-C	-5.54	108.57	114.67
2	G	91	PHE	N-CA-CB	-5.51	102.03	110.12
2	G	90	HIS	N-CA-C	-5.47	100.26	109.07
1	E	137	VAL	N-CA-C	5.42	115.90	111.62
1	B	61	ASN	N-CA-C	5.42	116.99	111.14
1	B	66	ILE	CA-C-N	5.41	125.35	119.78
1	B	66	ILE	C-N-CA	5.41	125.35	119.78
2	H	114	LYS	CB-CA-C	-5.38	110.39	116.63
1	E	48	ALA	N-CA-C	5.29	118.88	111.90
2	H	312	MET	N-CA-C	5.29	117.22	109.24
1	B	32	LEU	CA-C-N	5.28	124.60	118.85
1	B	32	LEU	C-N-CA	5.28	124.60	118.85
2	G	91	PHE	N-CA-C	5.25	117.00	111.28
1	E	24	GLY	N-CA-C	-5.18	106.22	112.49
2	D	33	ILE	N-CA-C	5.14	115.74	107.98
2	C	158	ASN	CA-C-N	5.01	124.92	119.85
2	C	158	ASN	C-N-CA	5.01	124.92	119.85
2	D	199	HIS	N-CA-C	-5.00	107.56	113.97
2	H	307	VAL	N-CA-C	5.00	115.68	108.48

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1627	0	1746	93	0
1	B	1627	0	1746	126	0
1	E	1616	0	1733	84	0
1	F	1627	0	1746	94	0
2	C	2661	0	2714	131	0
2	D	2651	0	2707	121	0
2	G	2651	0	2707	138	0
2	H	2630	0	2682	184	0
3	C	9	0	8	2	0
3	D	9	0	8	1	0
3	G	9	0	8	2	0
3	H	9	0	8	6	0
4	C	27	0	12	12	0
4	D	27	0	12	5	0
4	G	27	0	12	1	0
4	H	27	0	12	5	0
All	All	17234	0	17861	917	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 26.

All (917) 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:13:VAL:HA	1:F:172:LEU:HD11	1.16	1.16
2:G:293:ASN:OD1	2:H:284:SER:HB2	1.48	1.13
2:G:31:GLY:H	2:G:193:THR:HG22	1.03	1.06
1:B:115:ILE:HD11	1:B:138:LEU:HD23	1.38	1.05
1:B:8:LEU:O	1:B:11:ARG:N	1.91	1.04
2:H:91:PHE:CD2	2:H:146:LYS:HG3	1.93	1.04
1:A:185:ASN:HB3	1:A:188:VAL:CG1	1.88	1.02
2:H:312:MET:HB3	3:H:1000:MSE:CE	1.91	1.01
1:A:185:ASN:HB3	1:A:188:VAL:HG12	1.46	0.97
1:E:123:SER:HA	1:E:126:MET:HE3	1.45	0.97
2:D:307:VAL:HG13	2:D:309:PHE:HE1	1.27	0.97
1:A:2:SER:OG	1:A:5:MET:HB2	1.66	0.96
1:A:23:SER:HA	1:A:97:THR:HG22	1.44	0.95
1:F:19:MET:HE3	1:F:161:SER:OG	1.67	0.94
2:C:241:ILE:O	2:C:244:THR:HG22	1.68	0.94
2:D:9:LYS:HA	2:D:58:THR:HG22	1.47	0.94
2:G:49:ARG:HB2	2:G:54:LEU:HB2	1.48	0.93
2:G:31:GLY:N	2:G:193:THR:HG22	1.83	0.93

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:182:LEU:HD11	2:D:186:ILE:HD11	1.50	0.92
1:B:2:SER:C	1:B:4:PRO:HD2	1.94	0.92
1:A:5:MET:HE2	1:A:188:VAL:HG23	1.49	0.92
1:A:170:GLY:HA2	1:A:174:GLN:HE21	1.33	0.92
2:D:307:VAL:HG13	2:D:309:PHE:CE1	2.03	0.92
1:F:13:VAL:CA	1:F:172:LEU:HD11	2.00	0.91
2:H:266:VAL:HG13	2:H:316:MET:O	1.69	0.91
1:E:41:THR:HG21	1:E:52:LEU:HD11	1.49	0.90
1:F:155:ILE:HB	1:F:203:ILE:HD11	1.55	0.88
1:F:176:GLY:HA2	1:F:189:MET:HE1	1.55	0.87
2:C:11:PHE:HE2	4:C:2000:ADP:C4	1.93	0.87
1:E:147:VAL:HG23	1:E:211:VAL:HG22	1.57	0.87
2:H:128:VAL:HG11	2:H:152:ALA:HB2	1.54	0.87
2:H:260:GLU:OE1	2:H:261:PRO:HD2	1.75	0.86
2:H:48:ILE:HD12	2:H:195:LEU:HD12	1.58	0.86
1:B:42:ARG:HH11	1:B:113:LEU:HD11	1.39	0.85
2:H:292:VAL:HG21	2:H:316:MET:HE2	1.59	0.84
2:D:182:LEU:CD1	2:D:186:ILE:HD11	2.07	0.84
1:F:129:THR:H	1:F:132:GLN:HE21	1.26	0.84
1:E:118:GLY:C	2:G:92:ASN:ND2	2.35	0.83
2:H:114:LYS:HG2	2:H:115:ASP:H	1.42	0.82
2:C:11:PHE:HE2	4:C:2000:ADP:C5	1.97	0.82
1:A:5:MET:CE	1:A:188:VAL:HG23	2.09	0.82
1:A:69:ILE:HD12	1:A:163:MET:HE2	1.59	0.82
2:C:2:ILE:HD11	2:C:28:VAL:HG23	1.62	0.82
1:E:118:GLY:O	2:G:92:ASN:ND2	2.11	0.82
1:B:20:THR:HG22	1:B:154:LEU:HD12	1.62	0.82
2:H:110:ASP:O	2:H:111:ASN:HB2	1.78	0.82
1:B:72:LEU:HD21	1:B:167:VAL:HG11	1.62	0.82
1:E:28:PHE:HA	1:E:104:ILE:HD12	1.62	0.82
2:D:72:LEU:HD12	2:D:76:GLU:HB3	1.63	0.81
1:E:143:LEU:HD23	1:E:214:VAL:HG21	1.62	0.81
1:E:172:LEU:HD11	1:E:195:LEU:HD22	1.61	0.81
2:D:182:LEU:HD13	2:D:182:LEU:C	2.06	0.81
1:B:179:TYR:HD1	1:B:188:VAL:HG21	1.46	0.81
1:F:80:ARG:HB2	1:F:86:SER:HB3	1.64	0.80
2:G:248:ASP:HB3	2:H:300:GLN:HE22	1.45	0.80
1:A:185:ASN:HB3	1:A:188:VAL:HG11	1.64	0.80
1:A:196:LEU:HD13	1:B:70:ILE:HG13	1.63	0.80
2:C:2:ILE:HG22	2:C:64:VAL:HG13	1.63	0.80
1:A:13:VAL:HG22	1:A:172:LEU:HD11	1.64	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:G:9:LYS:HE2	2:G:56:ARG:O	1.82	0.79
2:H:312:MET:HB3	3:H:1000:MSE:HE2	1.64	0.79
2:C:11:PHE:CE2	4:C:2000:ADP:C4	2.70	0.79
1:E:134:VAL:HG12	1:E:138:LEU:HD12	1.65	0.79
1:F:92:ALA:HB2	1:F:169:ALA:HB1	1.65	0.79
2:G:329:TRP:CE2	2:G:333:HIS:NE2	2.51	0.79
2:D:98:THR:HG22	2:D:137:TYR:CE2	2.18	0.79
2:G:310:GLY:HA3	3:G:1000:MSE:CE	2.13	0.78
2:C:246:HIS:O	2:D:300:GLN:NE2	2.16	0.78
1:F:118:GLY:C	2:H:92:ASN:ND2	2.42	0.78
2:D:72:LEU:CD1	2:D:76:GLU:HB3	2.14	0.77
1:B:19:MET:HE2	1:B:96:LEU:HD11	1.66	0.77
1:E:20:THR:HG21	1:E:158:VAL:HG22	1.66	0.77
2:H:11:PHE:HB3	4:H:2000:ADP:C2	2.18	0.77
1:B:211:VAL:HA	1:B:214:VAL:HG12	1.65	0.77
2:C:0:HIS:HA	2:C:27:HIS:NE2	2.00	0.76
1:E:42:ARG:HB3	1:E:43:PRO:HD2	1.65	0.76
1:E:133:ILE:O	1:E:137:VAL:HG22	1.85	0.76
2:H:114:LYS:HG2	2:H:115:ASP:N	2.01	0.76
1:A:185:ASN:CB	1:A:188:VAL:HG12	2.15	0.76
2:D:296:ILE:HG23	2:D:312:MET:CE	2.16	0.76
2:D:100:PHE:HD2	2:D:134:HIS:CE1	2.04	0.75
2:G:300:GLN:HE21	2:H:246:HIS:CD2	2.03	0.75
2:H:45:SER:OG	4:H:2000:ADP:O1B	2.02	0.75
2:C:2:ILE:HD11	2:C:28:VAL:CG2	2.16	0.75
2:C:1:MET:HE3	2:C:161:VAL:HG23	1.66	0.75
1:E:32:LEU:HB3	1:E:33:PRO:HD3	1.69	0.75
2:G:99:VAL:HG11	2:G:141:LEU:HD11	1.68	0.75
2:H:174:ALA:HA	2:H:177:ARG:HH11	1.51	0.75
1:A:2:SER:OG	1:A:5:MET:CB	2.35	0.75
1:B:42:ARG:NH1	1:B:113:LEU:HD11	2.01	0.75
2:H:91:PHE:HD2	2:H:146:LYS:HG3	1.52	0.75
2:H:220:ILE:HG22	2:H:234:THR:HG21	1.67	0.75
2:D:296:ILE:HA	2:D:314:THR:HG22	1.67	0.74
2:G:300:GLN:HE21	2:H:246:HIS:HD2	1.35	0.74
2:H:91:PHE:CE2	2:H:146:LYS:HG3	2.22	0.74
1:E:13:VAL:HG22	1:E:195:LEU:HD21	1.68	0.74
2:G:1:MET:HB2	2:G:65:ASP:H	1.53	0.74
2:G:13:GLN:HB2	2:G:16:ARG:HB3	1.68	0.74
1:F:41:THR:HA	1:F:48:ALA:HA	1.70	0.74
1:F:137:VAL:O	1:F:141:GLU:HB2	1.87	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:118:GLY:C	2:G:92:ASN:HD21	1.96	0.74
2:G:99:VAL:CG1	2:G:141:LEU:HD11	2.18	0.73
2:H:48:ILE:CD1	2:H:195:LEU:HD12	2.18	0.73
2:D:6:ASN:H	2:D:25:SER:HB3	1.53	0.73
2:G:223:ASP:HB3	2:G:228:VAL:HG23	1.71	0.73
2:C:220:ILE:HD12	2:C:234:THR:HG21	1.70	0.73
1:F:46:ILE:HG22	1:F:124:ARG:HD3	1.69	0.73
1:F:143:LEU:HD23	1:F:214:VAL:HG21	1.70	0.73
2:D:4:LEU:HB3	2:D:7:ILE:HD12	1.68	0.72
1:E:172:LEU:CD1	1:E:195:LEU:HD22	2.19	0.72
1:B:178:GLN:OE1	1:B:182:ILE:HD12	1.89	0.72
1:B:72:LEU:CD2	1:B:167:VAL:HG11	2.18	0.72
2:D:182:LEU:HD13	2:D:182:LEU:O	1.90	0.72
1:B:188:VAL:HA	1:B:191:THR:HG22	1.72	0.72
2:D:201:MET:HE3	2:D:229:PHE:CZ	2.25	0.72
1:F:118:GLY:CA	2:H:92:ASN:HD21	2.02	0.72
1:B:32:LEU:HB3	1:B:33:PRO:HD3	1.72	0.71
2:D:307:VAL:CG1	2:D:309:PHE:HE1	2.02	0.71
1:A:69:ILE:CD1	1:A:163:MET:HE2	2.19	0.71
2:D:230:SER:HG	2:D:231:HIS:CE1	2.08	0.71
2:D:142:SER:HB2	2:D:145:GLN:HG3	1.71	0.71
2:H:319:THR:HG23	2:H:322:ASP:HB2	1.73	0.71
1:E:20:THR:HG21	1:E:158:VAL:CG2	2.21	0.71
2:C:98:THR:HG23	2:C:101:GLY:H	1.56	0.70
1:B:85:THR:HG22	1:B:87:ILE:H	1.56	0.70
2:G:45:SER:HA	2:G:48:ILE:HG22	1.73	0.70
2:H:53:LEU:HD12	2:H:69:LEU:HB3	1.72	0.70
1:A:2:SER:HG	1:A:5:MET:HB2	1.55	0.70
2:D:124:LEU:HD21	2:D:155:LEU:HB2	1.74	0.70
2:H:260:GLU:CD	2:H:261:PRO:HD2	2.16	0.70
2:C:283:LEU:HD11	2:C:312:MET:SD	2.32	0.70
1:A:68:PHE:HZ	1:A:96:LEU:HD12	1.56	0.69
2:D:296:ILE:HG23	2:D:312:MET:HE1	1.73	0.69
1:B:40:VAL:HG12	1:B:47:ILE:HD12	1.73	0.69
2:C:1:MET:HE1	2:C:160:LYS:HB3	1.74	0.69
1:F:210:ILE:O	1:F:214:VAL:HG12	1.91	0.69
2:G:248:ASP:HB3	2:H:300:GLN:NE2	2.06	0.69
2:D:312:MET:HE2	2:D:314:THR:HG23	1.74	0.69
2:G:9:LYS:HE3	2:G:55:GLU:HG3	1.74	0.69
2:D:168:THR:HA	2:D:171:LEU:HD12	1.75	0.69
2:D:180:LEU:HD11	2:D:203:VAL:HG12	1.75	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:312:MET:HE2	2:D:314:THR:CG2	2.23	0.69
1:E:163:MET:HG2	1:F:163:MET:HE2	1.75	0.69
1:B:177:TYR:O	1:B:181:TYR:HB3	1.92	0.69
2:G:91:PHE:O	2:G:92:ASN:HB2	1.92	0.69
2:H:35:GLY:HA2	2:H:196:LEU:HD13	1.74	0.69
1:A:154:LEU:HB3	1:A:203:ILE:HD11	1.74	0.68
1:A:200:VAL:O	1:A:203:ILE:HG22	1.91	0.68
2:H:269:LEU:HD12	2:H:337:VAL:HG13	1.75	0.68
1:F:3:GLU:N	1:F:4:PRO:HD2	2.08	0.68
2:C:310:GLY:HA3	3:C:1000:MSE:HE1	1.75	0.68
1:B:72:LEU:HD21	1:B:167:VAL:CG1	2.23	0.68
2:C:310:GLY:HA3	3:C:1000:MSE:CE	2.22	0.68
2:D:124:LEU:CD2	2:D:155:LEU:HB2	2.23	0.68
1:B:41:THR:HG22	1:B:41:THR:O	1.94	0.68
2:H:266:VAL:CG1	2:H:316:MET:O	2.40	0.68
1:B:211:VAL:HA	1:B:214:VAL:CG1	2.22	0.68
2:D:230:SER:HB3	2:D:309:PHE:CE2	2.29	0.68
1:B:41:THR:HG22	1:B:53:TYR:HB2	1.76	0.67
2:D:48:ILE:HD11	2:D:163:LEU:HB3	1.76	0.67
1:E:69:ILE:HG13	1:E:163:MET:HE2	1.74	0.67
2:H:99:VAL:HG21	2:H:133:LYS:O	1.94	0.67
1:A:64:ARG:NH1	1:A:103:PHE:CD1	2.62	0.67
2:D:237:ALA:O	2:D:241:ILE:HG12	1.94	0.67
1:E:162:ALA:O	1:F:69:ILE:HD11	1.95	0.67
2:H:196:LEU:HD11	2:H:208:CYS:SG	2.34	0.67
2:H:107:LEU:HD11	2:H:156:ALA:O	1.93	0.67
1:B:118:GLY:O	2:D:92:ASN:ND2	2.27	0.67
2:G:125:LEU:O	2:G:130:LEU:N	2.22	0.67
2:H:112:THR:HG22	2:H:117:VAL:HG23	1.77	0.67
1:B:216:ARG:N	1:B:216:ARG:HD2	2.09	0.67
2:C:144:GLY:O	2:C:148:ARG:HG3	1.95	0.67
2:C:283:LEU:O	2:C:286:THR:HG22	1.95	0.67
1:E:80:ARG:HH12	1:F:184:TYR:HE2	1.41	0.67
2:C:274:THR:O	2:C:278:VAL:HG23	1.95	0.66
2:G:46:THR:OG1	4:G:2000:ADP:O1A	2.11	0.66
1:E:172:LEU:HD11	1:E:195:LEU:CD2	2.25	0.66
2:D:215:SER:CB	2:D:236:LEU:HD11	2.26	0.66
2:H:91:PHE:CE1	2:H:147:GLN:NE2	2.62	0.66
2:H:294:ASN:HD21	2:H:316:MET:HE3	1.61	0.66
2:H:303:TYR:OH	2:H:308:LYS:HD2	1.96	0.66
2:C:187:ASN:ND2	2:C:192:LEU:O	2.27	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:105:LEU:N	2:D:106:PRO:HD2	2.11	0.66
1:E:27:GLY:C	1:E:104:ILE:HD11	2.20	0.66
2:H:268:MET:HE1	2:H:297:ILE:HD13	1.76	0.66
1:B:2:SER:CB	1:B:4:PRO:HD2	2.25	0.66
2:C:316:MET:HE1	2:C:326:ALA:HB3	1.77	0.66
1:E:27:GLY:O	1:E:104:ILE:HD11	1.96	0.66
2:G:329:TRP:CZ2	2:G:333:HIS:NE2	2.64	0.66
2:H:290:PHE:HB3	2:H:322:ASP:HB3	1.78	0.65
2:H:172:ASP:HB2	2:H:175:THR:OG1	1.96	0.65
2:G:45:SER:O	2:G:49:ARG:HG2	1.97	0.65
1:F:16:THR:HB	1:F:172:LEU:HD13	1.78	0.65
2:C:301:MET:SD	2:C:308:LYS:HE2	2.37	0.65
2:D:182:LEU:CD1	2:D:186:ILE:CD1	2.74	0.65
1:F:189:MET:HE2	1:F:189:MET:HA	1.79	0.65
1:B:172:LEU:CD2	1:B:192:VAL:HG13	2.27	0.65
1:F:16:THR:HB	1:F:172:LEU:CD1	2.27	0.65
2:G:46:THR:O	2:G:50:CYS:HB2	1.97	0.65
1:B:94:VAL:HB	1:B:95:PRO:HD3	1.77	0.65
2:C:130:LEU:HD11	2:C:148:ARG:HB2	1.77	0.65
2:C:130:LEU:HD13	2:C:149:VAL:HG22	1.79	0.65
1:E:167:VAL:HG12	1:E:167:VAL:O	1.97	0.65
2:H:327:ILE:HG23	2:H:337:VAL:HG21	1.79	0.65
1:B:3:GLU:N	1:B:4:PRO:CD	2.60	0.64
2:C:1:MET:CE	2:C:160:LYS:HB3	2.27	0.64
2:C:231:HIS:N	2:C:232:PRO:HD3	2.13	0.64
2:G:328:ALA:O	2:G:332:GLU:HG3	1.96	0.64
1:A:23:SER:HA	1:A:97:THR:CG2	2.23	0.64
2:C:13:GLN:NE2	4:C:2000:ADP:C2	2.65	0.64
1:F:118:GLY:HA3	2:H:92:ASN:HD21	1.62	0.64
1:A:3:GLU:N	1:A:4:PRO:HD2	2.13	0.64
1:B:179:TYR:CD1	1:B:188:VAL:HG21	2.30	0.64
1:E:3:GLU:N	1:E:4:PRO:HD2	2.12	0.64
2:G:91:PHE:HD1	2:G:147:GLN:HE21	1.46	0.64
2:H:142:SER:N	2:H:145:GLN:OE1	2.28	0.64
1:F:38:LEU:HD21	1:F:106:ARG:NH1	2.13	0.64
1:B:185:ASN:HB3	1:B:188:VAL:HG22	1.80	0.63
1:B:119:LEU:O	1:B:123:SER:HB2	1.97	0.63
2:H:272:GLU:HB2	2:H:336:LYS:HB3	1.80	0.63
2:C:168:THR:HG22	2:C:171:LEU:HD12	1.80	0.63
1:F:129:THR:H	1:F:132:GLN:NE2	1.94	0.63
2:H:286:THR:HG22	2:H:292:VAL:CG2	2.28	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:46:ILE:HG13	1:F:46:ILE:O	1.98	0.63
1:A:103:PHE:HE2	1:A:160:TYR:HH	1.44	0.63
1:A:127:GLY:O	2:C:78:THR:HG23	1.99	0.63
1:F:118:GLY:C	2:H:92:ASN:HD22	2.07	0.63
2:C:2:ILE:HG13	2:C:28:VAL:HB	1.81	0.63
2:D:45:SER:HB3	4:D:2000:ADP:PA	2.38	0.63
1:E:13:VAL:HG22	1:E:195:LEU:CD2	2.28	0.63
2:D:145:GLN:O	2:D:149:VAL:HG23	1.99	0.62
1:E:71:LEU:O	1:E:75:MET:HG2	1.99	0.62
2:H:91:PHE:CD1	2:H:147:GLN:NE2	2.67	0.62
2:H:280:ALA:O	3:H:1000:MSE:N	2.32	0.62
1:A:26:PHE:O	1:A:30:ILE:HG22	1.99	0.62
1:B:3:GLU:N	1:B:4:PRO:HD2	2.14	0.62
2:D:41:GLY:HA2	4:D:2000:ADP:H3'	1.79	0.62
2:G:249:ILE:HG13	2:G:249:ILE:O	1.99	0.62
2:D:45:SER:O	2:D:49:ARG:HG2	1.98	0.62
2:G:198:THR:HB	2:G:200:GLU:H	1.65	0.62
2:C:53:LEU:CD1	2:C:69:LEU:HB3	2.29	0.62
1:F:47:ILE:HG12	1:F:47:ILE:O	1.98	0.62
1:B:40:VAL:HG12	1:B:47:ILE:CD1	2.30	0.62
2:D:166:GLN:HB3	2:D:169:SER:HB2	1.81	0.62
2:D:214:ILE:HA	2:D:219:LEU:HA	1.82	0.62
2:H:257:LEU:HG	2:H:258:GLN:O	2.00	0.62
1:A:93:ILE:HG22	1:A:93:ILE:O	1.99	0.62
1:B:45:GLN:HG3	1:B:46:ILE:H	1.64	0.62
1:B:36:VAL:HG12	1:B:112:LEU:HD11	1.81	0.61
2:H:220:ILE:CG2	2:H:234:THR:HG21	2.29	0.61
2:C:0:HIS:CA	2:C:27:HIS:NE2	2.64	0.61
2:D:230:SER:HB3	2:D:309:PHE:CZ	2.35	0.61
2:D:307:VAL:CG1	2:D:309:PHE:CE1	2.81	0.61
1:F:167:VAL:HG12	1:F:167:VAL:O	2.00	0.61
2:D:106:PRO:HG3	2:D:153:ARG:HG3	1.81	0.61
2:C:94:LEU:HD12	2:C:97:ARG:HD3	1.82	0.61
1:B:42:ARG:HG3	1:B:43:PRO:HD2	1.82	0.61
2:C:0:HIS:N	2:C:27:HIS:NE2	2.48	0.61
2:H:262:PHE:CE1	2:H:265:CYS:HB2	2.35	0.61
2:H:91:PHE:HZ	2:H:171:LEU:HD21	1.65	0.61
2:G:293:ASN:O	2:G:317:HIS:HB2	2.01	0.60
1:F:139:LEU:HB2	1:F:140:PRO:HD3	1.82	0.60
1:F:64:ARG:HD2	1:F:103:PHE:HA	1.83	0.60
2:H:105:LEU:HB3	2:H:106:PRO:HD3	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:155:ILE:CG1	1:A:203:ILE:CG2	2.79	0.60
1:B:45:GLN:HG3	1:B:46:ILE:N	2.17	0.59
2:C:272:GLU:HG2	2:C:311:ILE:HG12	1.84	0.59
1:A:32:LEU:HB3	1:A:33:PRO:HD3	1.83	0.59
1:B:42:ARG:CG	1:B:43:PRO:HD2	2.32	0.59
2:C:13:GLN:NE2	4:C:2000:ADP:H2	2.00	0.59
2:H:300:GLN:HE21	2:H:301:MET:H	1.50	0.59
1:B:167:VAL:HG22	1:B:167:VAL:O	2.02	0.59
2:H:269:LEU:HD13	2:H:338:GLU:O	2.02	0.59
2:C:164:CYS:HB2	2:C:196:LEU:CD1	2.32	0.59
2:D:21:LEU:HA	2:D:217:GLY:HA2	1.83	0.59
2:D:118:LYS:O	2:D:122:THR:HG23	2.02	0.59
2:D:182:LEU:CD1	2:D:182:LEU:C	2.75	0.59
1:F:176:GLY:CA	1:F:189:MET:HE1	2.30	0.59
1:B:8:LEU:O	1:B:9:LEU:C	2.44	0.59
2:D:177:ARG:O	2:D:181:GLU:HB2	2.01	0.59
2:G:248:ASP:O	2:G:249:ILE:HG22	2.03	0.59
2:C:297:ILE:HA	2:D:301:MET:HE3	1.84	0.59
1:E:159:GLY:HA2	1:F:67:PRO:HB3	1.84	0.59
2:C:168:THR:HB	2:C:176:THR:HG22	1.85	0.59
1:E:71:LEU:HG	1:E:75:MET:HE3	1.85	0.59
2:H:112:THR:HB	2:H:117:VAL:CG2	2.33	0.59
1:B:115:ILE:HG13	1:B:115:ILE:O	2.02	0.59
2:H:274:THR:HG23	2:H:334:HIS:HD2	1.67	0.59
2:G:310:GLY:HA3	3:G:1000:MSE:HE1	1.84	0.58
2:H:47:LEU:C	2:H:47:LEU:HD12	2.26	0.58
2:D:260:GLU:HB2	2:D:261:PRO:HD2	1.84	0.58
1:A:30:ILE:O	1:A:34:VAL:HG23	2.03	0.58
2:G:293:ASN:OD1	2:H:284:SER:CB	2.40	0.58
1:B:189:MET:HE2	1:B:189:MET:HA	1.85	0.58
2:D:215:SER:HB2	2:D:236:LEU:HD11	1.84	0.58
2:D:270:ARG:HH22	2:D:272:GLU:CD	2.12	0.58
2:G:98:THR:HG23	2:G:101:GLY:H	1.68	0.58
2:C:72:LEU:HD23	2:C:77:LEU:HD12	1.85	0.58
2:C:141:LEU:HB3	2:C:145:GLN:HB2	1.86	0.58
2:D:329:TRP:O	2:D:333:HIS:HD2	1.87	0.58
1:B:2:SER:HB2	1:B:5:MET:HB2	1.86	0.58
2:C:107:LEU:HB3	2:C:112:THR:HG21	1.86	0.58
2:G:2:ILE:HG12	2:G:2:ILE:O	2.04	0.58
1:E:17:LEU:HD21	1:E:199:LEU:HD22	1.85	0.57
2:H:45:SER:N	4:H:2000:ADP:O1B	2.37	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:G:2:ILE:HG23	2:G:28:VAL:HB	1.84	0.57
2:G:91:PHE:HD1	2:G:147:GLN:NE2	2.02	0.57
1:A:155:ILE:HG12	1:A:203:ILE:CG2	2.34	0.57
2:G:294:ASN:HB3	2:G:316:MET:HA	1.87	0.57
2:C:201:MET:HE2	2:C:240:PHE:HB3	1.85	0.57
2:G:327:ILE:O	2:G:331:GLN:HG3	2.05	0.57
1:B:172:LEU:HD23	1:B:192:VAL:HG13	1.86	0.57
2:C:220:ILE:CD1	2:C:236:LEU:HD23	2.35	0.57
2:H:271:LEU:HD11	2:H:283:LEU:HD21	1.87	0.57
2:G:287:ALA:O	2:G:291:ASN:HA	2.05	0.57
1:A:20:THR:HG22	1:A:157:LEU:O	2.05	0.57
2:G:333:HIS:O	2:G:335:VAL:N	2.38	0.57
2:H:89:GLN:NE2	2:H:169:SER:HB3	2.19	0.57
1:A:23:SER:CA	1:A:97:THR:HG22	2.27	0.56
1:A:157:LEU:HA	1:A:160:TYR:HB2	1.87	0.56
2:G:323:THR:O	2:G:327:ILE:HD12	2.05	0.56
1:F:118:GLY:CA	2:H:92:ASN:ND2	2.66	0.56
1:A:155:ILE:CG1	1:A:203:ILE:HG23	2.36	0.56
2:H:266:VAL:HG13	2:H:267:PRO:HD2	1.88	0.56
2:D:215:SER:OG	2:D:236:LEU:HD11	2.06	0.56
1:B:179:TYR:N	1:B:179:TYR:HD2	2.04	0.56
2:D:198:THR:HG21	2:D:203:VAL:CG2	2.36	0.56
1:E:31:GLY:HA3	1:E:104:ILE:HG13	1.86	0.56
2:G:51:VAL:HB	2:G:163:LEU:HD21	1.87	0.56
2:C:241:ILE:O	2:C:244:THR:CG2	2.50	0.56
1:B:179:TYR:N	1:B:179:TYR:CD2	2.73	0.56
1:A:57:SER:OG	1:A:106:ARG:NH2	2.34	0.56
2:C:91:PHE:CD2	2:C:91:PHE:N	2.74	0.56
1:E:214:VAL:O	1:E:214:VAL:HG13	2.04	0.55
2:H:286:THR:HG22	2:H:292:VAL:HG21	1.86	0.55
2:H:319:THR:CG2	2:H:322:ASP:HB2	2.36	0.55
2:C:103:VAL:HG21	2:C:125:LEU:HD11	1.88	0.55
2:D:274:THR:O	2:D:277:SER:HB2	2.07	0.55
2:G:232:PRO:HG2	2:G:238:GLN:HG2	1.87	0.55
2:D:223:ASP:HB3	2:D:228:VAL:HG23	1.89	0.55
2:G:28:VAL:HG22	2:G:34:TYR:CD2	2.41	0.55
1:A:144:PRO:HB3	1:A:214:VAL:HG21	1.88	0.55
1:F:128:ALA:HA	2:H:109:LEU:HD21	1.88	0.55
2:G:245:LEU:HD21	2:G:309:PHE:CZ	2.41	0.55
2:G:241:ILE:O	2:G:244:THR:HG22	2.05	0.55
2:C:182:LEU:CD1	2:C:186:ILE:HD11	2.37	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:125:ALA:HB1	2:G:52:ASN:HD21	1.72	0.55
2:G:201:MET:HE3	2:G:229:PHE:CZ	2.42	0.55
1:A:196:LEU:CD1	1:B:70:ILE:HG13	2.33	0.55
2:H:112:THR:CG2	2:H:117:VAL:HG23	2.36	0.55
1:B:2:SER:HB3	1:B:4:PRO:HD2	1.89	0.55
2:C:141:LEU:HB2	2:C:146:LYS:HG3	1.88	0.55
2:D:172:ASP:O	2:D:176:THR:N	2.39	0.55
2:H:268:MET:CE	2:H:313:LEU:HD13	2.36	0.55
2:H:303:TYR:CE1	2:H:308:LYS:HB2	2.42	0.54
2:G:99:VAL:HG13	2:G:136:SER:O	2.07	0.54
2:H:4:LEU:HD21	2:H:51:VAL:HG22	1.90	0.54
1:F:92:ALA:CB	1:F:169:ALA:HB1	2.36	0.54
1:F:180:GLY:HA3	1:F:189:MET:HE3	1.89	0.54
2:G:146:LYS:O	2:G:149:VAL:HG12	2.06	0.54
2:H:91:PHE:HE1	2:H:147:GLN:HE21	1.55	0.54
1:A:136:LYS:NZ	2:C:108:GLU:OE1	2.41	0.54
2:C:283:LEU:CD1	2:C:312:MET:SD	2.95	0.54
2:G:247:LEU:HD22	2:G:298:SER:HB3	1.89	0.54
2:C:299:ALA:CB	2:C:312:MET:HG2	2.38	0.54
1:F:32:LEU:HB3	1:F:33:PRO:HD3	1.88	0.54
2:H:336:LYS:HE2	2:H:338:GLU:OE2	2.07	0.54
1:B:133:ILE:O	1:B:137:VAL:HG13	2.07	0.54
2:D:205:LYS:HG2	2:D:229:PHE:CZ	2.43	0.54
2:D:105:LEU:HD23	2:D:105:LEU:O	2.07	0.54
1:F:41:THR:OG1	1:F:53:TYR:HB2	2.08	0.54
2:G:110:ASP:O	2:G:112:THR:HG23	2.08	0.54
2:H:270:ARG:HH21	2:H:311:ILE:HD11	1.72	0.54
1:A:93:ILE:O	1:A:93:ILE:CG2	2.56	0.53
1:A:100:ALA:O	1:A:104:ILE:HG12	2.07	0.53
1:B:150:ALA:O	1:B:154:LEU:HD23	2.09	0.53
2:G:272:GLU:HB2	2:G:336:LYS:HB3	1.91	0.53
1:F:191:THR:O	1:F:195:LEU:HB2	2.09	0.53
2:D:262:PHE:HZ	2:D:343:VAL:C	2.15	0.53
1:F:101:ALA:HB3	1:F:102:PRO:HD3	1.91	0.53
2:G:31:GLY:H	2:G:193:THR:CG2	1.97	0.53
2:H:47:LEU:HD12	2:H:47:LEU:O	2.08	0.53
2:D:286:THR:HG22	2:D:329:TRP:CE3	2.44	0.53
2:G:103:VAL:HG21	2:G:125:LEU:HD11	1.91	0.53
1:A:120:ILE:HD12	1:A:138:LEU:HD11	1.91	0.53
1:B:40:VAL:HA	1:B:45:GLN:HG2	1.90	0.53
2:G:252:ASP:O	2:G:256:ARG:HG2	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:40:VAL:HA	1:A:45:GLN:HG2	1.90	0.53
1:B:20:THR:CG2	1:B:154:LEU:HD12	2.38	0.53
1:B:185:ASN:HB3	1:B:188:VAL:CG2	2.40	0.52
2:C:11:PHE:CE2	4:C:2000:ADP:C5	2.89	0.52
2:H:312:MET:HB3	3:H:1000:MSE:HE3	1.87	0.52
2:H:333:HIS:C	2:H:334:HIS:ND1	2.67	0.52
1:F:2:SER:C	1:F:4:PRO:HD2	2.34	0.52
1:E:13:VAL:HG21	1:E:195:LEU:HD11	1.92	0.52
1:A:120:ILE:CD1	1:A:138:LEU:HD11	2.38	0.52
2:G:72:LEU:HD23	2:G:77:LEU:HD23	1.90	0.52
2:H:28:VAL:HG22	2:H:34:TYR:CD2	2.45	0.52
2:H:333:HIS:O	2:H:334:HIS:CG	2.62	0.52
1:F:13:VAL:HA	1:F:172:LEU:CD1	2.11	0.52
2:G:226:SER:HB2	2:G:307:VAL:CG1	2.40	0.52
2:C:103:VAL:HG21	2:C:125:LEU:CD1	2.40	0.52
2:G:245:LEU:HD11	2:G:302:ASP:HB2	1.92	0.52
2:H:112:THR:HB	2:H:117:VAL:HG22	1.92	0.52
1:F:119:LEU:HD23	2:H:94:LEU:HD23	1.92	0.52
1:B:177:TYR:CE2	1:B:181:TYR:CD2	2.98	0.52
2:H:105:LEU:HD23	2:H:105:LEU:C	2.35	0.52
2:H:263:THR:O	2:H:264:ASP:HB3	2.09	0.52
1:A:159:GLY:HA2	1:B:67:PRO:HB3	1.92	0.51
1:B:30:ILE:O	1:B:34:VAL:HG23	2.10	0.51
1:B:32:LEU:CB	1:B:33:PRO:HD3	2.36	0.51
2:C:182:LEU:HD13	2:C:186:ILE:CD1	2.40	0.51
2:D:262:PHE:O	2:D:265:CYS:HB3	2.10	0.51
2:G:271:LEU:HD23	2:G:337:VAL:HG13	1.90	0.51
2:C:182:LEU:HD11	2:C:186:ILE:HD11	1.92	0.51
2:G:107:LEU:O	2:G:110:ASP:O	2.27	0.51
2:G:124:LEU:O	2:G:128:VAL:HG22	2.10	0.51
2:G:287:ALA:O	2:H:288:ARG:NH1	2.43	0.51
2:H:162:LEU:HB2	2:H:192:LEU:HD11	1.91	0.51
2:H:198:THR:OG1	2:H:203:VAL:HG11	2.10	0.51
2:C:9:LYS:HD3	2:C:56:ARG:O	2.10	0.51
2:C:128:VAL:HB	2:C:148:ARG:HB3	1.93	0.51
2:D:11:PHE:CE2	4:D:2000:ADP:C4	2.98	0.51
2:D:34:TYR:OH	2:D:212:ALA:HB2	2.10	0.51
2:D:100:PHE:HD2	2:D:134:HIS:HE1	1.58	0.51
1:E:27:GLY:C	1:E:104:ILE:CD1	2.83	0.51
2:C:53:LEU:HD12	2:C:69:LEU:HB3	1.91	0.51
2:C:234:THR:HG22	2:C:237:ALA:H	1.75	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:116:GLU:HA	2:H:119:ARG:NH1	2.25	0.51
2:C:316:MET:CE	2:C:326:ALA:HB3	2.40	0.51
2:D:105:LEU:HD23	2:D:105:LEU:C	2.36	0.51
2:D:249:ILE:HG12	2:D:313:LEU:HD11	1.92	0.51
1:E:104:ILE:O	1:E:108:VAL:HG23	2.11	0.51
2:C:1:MET:CE	2:C:161:VAL:HG23	2.38	0.51
2:C:220:ILE:HD11	2:C:236:LEU:HD23	1.92	0.51
2:C:299:ALA:HB1	2:C:312:MET:HG2	1.92	0.51
2:D:34:TYR:OH	2:D:222:GLN:HG3	2.10	0.51
2:D:293:ASN:HB2	2:D:317:HIS:HB2	1.92	0.51
1:E:52:LEU:O	1:E:56:VAL:N	2.37	0.51
2:G:49:ARG:HG3	2:G:55:GLU:HB2	1.92	0.51
2:G:302:ASP:OD1	2:G:303:TYR:N	2.41	0.51
1:B:172:LEU:HD21	1:B:192:VAL:HG13	1.92	0.51
2:C:52:ASN:HB3	2:C:54:LEU:HD12	1.93	0.51
1:E:143:LEU:HD23	1:E:214:VAL:CG2	2.38	0.51
1:F:30:ILE:O	1:F:34:VAL:HG23	2.11	0.51
1:F:37:LEU:HD11	1:F:52:LEU:HD21	1.92	0.51
1:A:94:VAL:H	1:A:95:PRO:HD2	1.76	0.51
1:B:180:GLY:CA	1:B:189:MET:HE3	2.40	0.51
2:C:234:THR:HG23	2:C:235:PRO:HD2	1.92	0.51
1:E:189:MET:HB3	1:F:74:TRP:HD1	1.76	0.51
1:B:177:TYR:CE2	1:B:181:TYR:HD2	2.29	0.51
1:E:96:LEU:CD2	1:E:160:TYR:HB3	2.40	0.51
2:G:93:LEU:HD11	2:G:149:VAL:CG1	2.41	0.51
2:G:221:GLU:CD	2:G:233:LYS:H	2.18	0.51
2:H:91:PHE:CD2	2:H:146:LYS:CG	2.82	0.51
2:H:91:PHE:HD2	2:H:146:LYS:CG	2.21	0.51
2:H:283:LEU:HD11	2:H:312:MET:SD	2.50	0.51
1:B:104:ILE:HD11	1:B:153:THR:HG21	1.93	0.50
2:G:91:PHE:CD2	2:G:91:PHE:N	2.60	0.50
2:G:24:VAL:HG11	2:G:47:LEU:HD23	1.93	0.50
1:A:179:TYR:HB3	1:A:188:VAL:HG11	1.92	0.50
2:C:182:LEU:HD13	2:C:182:LEU:C	2.36	0.50
2:D:182:LEU:HD13	2:D:186:ILE:CD1	2.40	0.50
1:A:72:LEU:C	1:A:72:LEU:HD23	2.37	0.50
1:B:57:SER:OG	1:B:106:ARG:NH2	2.33	0.50
2:G:1:MET:CB	2:G:65:ASP:H	2.23	0.50
2:C:44:LYS:HE2	4:C:2000:ADP:O3B	2.11	0.50
2:C:52:ASN:HB3	2:C:54:LEU:CD1	2.41	0.50
2:D:83:GLN:O	2:D:159:PRO:HB2	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:134:VAL:HA	1:E:138:LEU:HB2	1.93	0.50
1:F:28:PHE:HZ	1:F:150:ALA:HA	1.77	0.50
1:A:139:LEU:HB2	1:A:140:PRO:HD3	1.94	0.50
2:G:204:VAL:HA	2:G:208:CYS:SG	2.52	0.50
2:H:49:ARG:HB2	2:H:54:LEU:HB2	1.94	0.50
2:H:110:ASP:O	2:H:111:ASN:CB	2.56	0.50
2:G:105:LEU:N	2:G:106:PRO:HD2	2.26	0.50
2:G:296:ILE:HG13	2:G:296:ILE:O	2.12	0.50
1:A:127:GLY:HA2	2:C:81:ARG:HD3	1.93	0.50
2:C:234:THR:HG22	2:C:236:LEU:N	2.27	0.50
1:E:6:MET:C	1:E:8:LEU:H	2.18	0.50
1:F:200:VAL:O	1:F:203:ILE:HG13	2.11	0.50
2:H:198:THR:HG21	2:H:203:VAL:HG13	1.94	0.49
2:H:268:MET:HE3	2:H:313:LEU:HD13	1.93	0.49
2:H:303:TYR:CZ	2:H:308:LYS:HD2	2.46	0.49
1:A:2:SER:HB2	1:A:4:PRO:HD2	1.94	0.49
1:B:101:ALA:HB3	1:B:102:PRO:HD3	1.94	0.49
2:C:164:CYS:HB2	2:C:196:LEU:HD13	1.94	0.49
1:E:58:ALA:O	1:E:62:ILE:HG12	2.13	0.49
2:G:30:ALA:HA	2:G:193:THR:CG2	2.42	0.49
2:H:273:PHE:HE1	2:H:282:LEU:HD12	1.76	0.49
2:H:319:THR:HG23	2:H:322:ASP:H	1.77	0.49
2:H:333:HIS:O	2:H:334:HIS:ND1	2.45	0.49
2:G:50:CYS:SG	2:G:57:PRO:HG3	2.52	0.49
2:H:120:ARG:NH2	2:H:156:ALA:O	2.45	0.49
1:B:8:LEU:O	1:B:11:ARG:HB3	2.12	0.49
1:B:17:LEU:HD21	1:B:199:LEU:HD11	1.94	0.49
2:C:73:SER:OG	2:C:76:GLU:HB2	2.12	0.49
2:C:286:THR:O	2:C:292:VAL:HG22	2.13	0.49
2:D:100:PHE:CD2	2:D:134:HIS:CE1	2.94	0.49
1:E:118:GLY:HA2	1:E:121:GLU:HB2	1.93	0.49
1:F:30:ILE:HD11	1:F:101:ALA:HB2	1.93	0.49
2:G:155:LEU:HD11	2:G:186:ILE:HD13	1.93	0.49
2:G:236:LEU:O	2:G:239:LYS:HB2	2.12	0.49
2:H:273:PHE:CE1	2:H:282:LEU:HD12	2.47	0.49
2:C:128:VAL:HG11	2:C:152:ALA:HB2	1.95	0.49
2:D:225:VAL:HG13	2:D:229:PHE:CE2	2.47	0.49
1:E:25:PHE:C	1:E:25:PHE:CD2	2.90	0.49
2:D:327:ILE:HG23	2:D:337:VAL:HG11	1.94	0.49
1:E:207:GLY:O	1:E:211:VAL:HG23	2.13	0.49
2:C:13:GLN:CD	4:C:2000:ADP:H2	2.20	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:G:301:MET:HG3	2:H:297:ILE:HG23	1.95	0.49
2:C:205:LYS:HB3	2:C:229:PHE:CZ	2.48	0.49
1:F:133:ILE:HG22	1:F:138:LEU:HG	1.94	0.49
2:G:36:VAL:HB	2:G:197:ILE:HG22	1.94	0.49
1:A:41:THR:HG21	1:A:52:LEU:HD12	1.95	0.48
1:B:211:VAL:CA	1:B:214:VAL:HG12	2.39	0.48
2:G:226:SER:HB2	2:G:307:VAL:HG11	1.95	0.48
2:H:29:PRO:HG2	2:H:32:GLN:OE1	2.13	0.48
2:H:89:GLN:HE21	2:H:169:SER:HB3	1.77	0.48
2:G:105:LEU:O	2:G:105:LEU:HD23	2.13	0.48
1:B:45:GLN:CG	1:B:46:ILE:H	2.26	0.48
1:E:63:PHE:HB3	1:E:102:PRO:HG2	1.94	0.48
2:G:299:ALA:HB2	2:G:312:MET:HG3	1.95	0.48
2:H:105:LEU:O	2:H:108:GLU:HB2	2.13	0.48
2:H:114:LYS:CG	2:H:115:ASP:N	2.70	0.48
1:A:45:GLN:HG3	1:A:46:ILE:H	1.78	0.48
2:C:8:THR:HG23	2:C:21:LEU:O	2.13	0.48
2:C:145:GLN:O	2:C:149:VAL:HG23	2.13	0.48
1:B:42:ARG:HE	1:B:113:LEU:HD21	1.77	0.48
1:B:66:ILE:HG22	1:B:71:LEU:HB2	1.96	0.48
1:F:68:PHE:CZ	1:F:96:LEU:HD23	2.48	0.48
2:G:93:LEU:HD11	2:G:149:VAL:HG13	1.94	0.48
2:G:316:MET:HE3	2:G:323:THR:HG23	1.95	0.48
1:A:211:VAL:O	1:A:214:VAL:HG22	2.14	0.48
2:C:252:ASP:O	2:C:256:ARG:HG3	2.14	0.48
2:D:93:LEU:HB3	2:D:102:ASN:HD21	1.77	0.48
2:D:234:THR:HG23	2:D:235:PRO:HD2	1.94	0.48
2:D:268:MET:HE3	2:D:297:ILE:CD1	2.43	0.48
2:G:93:LEU:CD1	2:G:149:VAL:CG1	2.92	0.48
1:A:115:ILE:HG23	1:A:119:LEU:HD12	1.96	0.48
1:B:89:LEU:HD12	1:B:90:GLN:N	2.29	0.48
1:E:69:ILE:HG23	1:E:70:ILE:HD12	1.96	0.48
1:F:46:ILE:O	1:F:47:ILE:HG22	2.13	0.48
2:G:83:GLN:HA	2:G:159:PRO:HA	1.96	0.48
2:D:296:ILE:CG2	2:D:312:MET:HE1	2.41	0.48
2:D:329:TRP:O	2:D:333:HIS:CD2	2.66	0.48
1:E:100:ALA:O	1:E:103:PHE:HB2	2.13	0.48
2:D:130:LEU:HD11	2:D:148:ARG:HB2	1.96	0.47
2:G:89:GLN:HB2	2:G:166:GLN:HB2	1.96	0.47
2:H:0:HIS:CB	2:H:29:PRO:HA	2.44	0.47
2:C:53:LEU:HD11	2:C:69:LEU:HB3	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:225:VAL:O	2:D:229:PHE:HD2	1.97	0.47
1:E:30:ILE:HD11	1:E:101:ALA:HB2	1.95	0.47
1:E:35:GLY:HA2	1:E:38:LEU:HD12	1.96	0.47
2:G:278:VAL:HG12	2:G:301:MET:SD	2.54	0.47
1:A:30:ILE:O	1:A:33:PRO:HD2	2.14	0.47
1:A:103:PHE:HE2	1:A:160:TYR:OH	1.95	0.47
2:G:301:MET:O	2:G:302:ASP:HB2	2.15	0.47
2:H:72:LEU:HG	2:H:76:GLU:HB2	1.95	0.47
2:C:163:LEU:HD23	2:C:195:LEU:HB3	1.95	0.47
1:E:162:ALA:C	1:F:69:ILE:HD11	2.39	0.47
2:H:89:GLN:HB2	2:H:166:GLN:HB2	1.97	0.47
1:E:3:GLU:N	1:E:4:PRO:CD	2.78	0.47
2:G:30:ALA:HA	2:G:193:THR:HG21	1.96	0.47
2:G:168:THR:HB	2:G:176:THR:HG23	1.97	0.47
2:G:248:ASP:O	2:G:249:ILE:CG2	2.62	0.47
1:A:19:MET:HG3	1:A:89:LEU:HD22	1.97	0.47
1:A:155:ILE:HG12	1:A:203:ILE:HG21	1.96	0.47
1:E:148:ASN:O	1:E:151:THR:HB	2.14	0.47
1:F:118:GLY:HA3	2:H:92:ASN:ND2	2.25	0.47
2:G:11:PHE:HB2	2:G:18:ILE:HG22	1.95	0.47
2:G:250:PRO:HG2	2:H:301:MET:HE2	1.96	0.47
1:B:172:LEU:HD23	1:B:172:LEU:O	2.14	0.47
2:C:172:ASP:O	2:C:176:THR:HG23	2.14	0.47
1:B:179:TYR:O	1:B:184:TYR:HA	2.15	0.47
2:C:4:LEU:HD11	2:C:51:VAL:HG22	1.96	0.47
2:C:164:CYS:HB2	2:C:196:LEU:HD12	1.96	0.47
2:D:63:LEU:O	2:D:63:LEU:HD12	2.15	0.47
1:F:12:GLY:O	1:F:172:LEU:HD12	2.15	0.47
1:B:32:LEU:HD12	1:B:32:LEU:HA	1.68	0.47
2:D:124:LEU:HD22	2:D:152:ALA:O	2.15	0.47
2:G:37:ILE:HG12	2:G:240:PHE:CE1	2.50	0.47
1:A:3:GLU:N	1:A:4:PRO:CD	2.78	0.46
1:A:129:THR:OG1	1:A:132:GLN:HG3	2.15	0.46
2:C:3:LYS:HG2	2:C:27:HIS:CD2	2.50	0.46
2:C:171:LEU:HD13	2:C:179:ILE:HG13	1.97	0.46
1:F:129:THR:N	1:F:132:GLN:HE21	2.04	0.46
2:H:45:SER:HB2	4:H:2000:ADP:PA	2.55	0.46
1:E:92:ALA:O	1:E:95:PRO:HD2	2.16	0.46
1:A:155:ILE:HD11	1:A:203:ILE:HG23	1.97	0.46
1:B:41:THR:O	1:B:41:THR:CG2	2.62	0.46
1:E:141:GLU:OE2	2:G:95:SER:HB2	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:313:LEU:HD23	2:H:313:LEU:HA	1.76	0.46
1:B:8:LEU:O	1:B:11:ARG:CA	2.63	0.46
1:B:38:LEU:HD21	1:B:56:VAL:HG12	1.97	0.46
2:C:130:LEU:CD1	2:C:148:ARG:HB2	2.44	0.46
2:D:11:PHE:CZ	4:D:2000:ADP:C5	3.03	0.46
2:G:296:ILE:HG22	2:G:314:THR:CG2	2.45	0.46
2:H:273:PHE:CZ	3:H:1000:MSE:HG3	2.51	0.46
1:B:107:MET:HE1	1:B:152:ILE:HD12	1.96	0.46
2:D:119:ARG:O	2:D:122:THR:OG1	2.33	0.46
2:D:176:THR:HG22	2:D:180:LEU:CD1	2.46	0.46
1:E:68:PHE:CZ	1:E:96:LEU:HD23	2.51	0.46
1:F:50:ALA:O	1:F:54:ARG:HB2	2.14	0.46
1:B:66:ILE:CG2	1:B:71:LEU:HB2	2.46	0.46
2:C:293:ASN:OD1	2:D:288:ARG:NH1	2.44	0.46
2:G:331:GLN:HA	2:G:335:VAL:O	2.15	0.46
2:H:312:MET:CE	2:H:314:THR:HB	2.46	0.46
1:B:27:GLY:O	1:B:31:GLY:N	2.48	0.46
2:C:154:ALA:O	2:C:159:PRO:HG3	2.16	0.46
1:B:2:SER:CA	1:B:4:PRO:HD2	2.44	0.46
1:B:19:MET:HE2	1:B:96:LEU:CD1	2.42	0.46
1:E:5:MET:O	1:E:8:LEU:HB3	2.14	0.46
1:E:42:ARG:HB3	1:E:43:PRO:CD	2.38	0.46
2:D:230:SER:HB3	2:D:309:PHE:HE2	1.77	0.46
2:D:329:TRP:CE2	2:D:333:HIS:NE2	2.84	0.46
1:E:37:LEU:HA	1:E:40:VAL:HG22	1.97	0.46
1:F:119:LEU:HD23	2:H:94:LEU:CD2	2.46	0.46
2:H:286:THR:O	2:H:292:VAL:HG22	2.16	0.46
2:C:3:LYS:CG	2:C:27:HIS:HD2	2.29	0.46
2:D:198:THR:HG21	2:D:203:VAL:HG23	1.96	0.46
1:F:16:THR:CB	1:F:172:LEU:HD13	2.44	0.46
2:H:45:SER:HB2	4:H:2000:ADP:O2A	2.15	0.46
2:H:270:ARG:HA	2:H:313:LEU:HD23	1.98	0.46
1:A:60:VAL:O	1:A:64:ARG:HB3	2.15	0.45
1:A:154:LEU:HB3	1:A:203:ILE:CD1	2.42	0.45
2:C:2:ILE:CG2	2:C:64:VAL:HG13	2.42	0.45
2:C:11:PHE:CD2	4:C:2000:ADP:N3	2.84	0.45
2:D:49:ARG:HD3	2:D:86:MET:HE1	1.97	0.45
1:F:3:GLU:N	1:F:4:PRO:CD	2.79	0.45
2:H:312:MET:HE2	2:H:314:THR:HB	1.98	0.45
2:C:160:LYS:O	2:C:192:LEU:HD12	2.16	0.45
1:F:178:GLN:O	1:F:182:ILE:HG13	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:268:MET:O	2:H:339:VAL:HA	2.16	0.45
2:D:234:THR:HG22	2:D:236:LEU:H	1.81	0.45
2:G:47:LEU:O	2:G:50:CYS:HB3	2.17	0.45
1:A:20:THR:HG22	1:A:157:LEU:C	2.42	0.45
1:A:92:ALA:C	1:A:94:VAL:H	2.24	0.45
1:A:155:ILE:CG1	1:A:203:ILE:HG21	2.47	0.45
2:C:156:ALA:C	2:C:158:ASN:H	2.23	0.45
2:D:9:LYS:HA	2:D:58:THR:CG2	2.31	0.45
1:E:181:TYR:CD1	1:F:167:VAL:HG13	2.52	0.45
1:F:41:THR:HG21	1:F:52:LEU:HD23	1.98	0.45
2:H:273:PHE:HB3	2:H:278:VAL:HG22	1.98	0.45
1:A:62:ILE:O	1:A:66:ILE:HG13	2.17	0.45
1:F:101:ALA:N	1:F:102:PRO:CD	2.79	0.45
2:H:52:ASN:OD1	2:H:54:LEU:HG	2.17	0.45
2:H:288:ARG:O	2:H:289:ARG:C	2.60	0.45
1:B:76:ILE:HD12	1:B:86:SER:OG	2.17	0.45
1:F:128:ALA:HA	2:H:109:LEU:CD2	2.46	0.45
2:H:93:LEU:HD22	2:H:102:ASN:CG	2.42	0.45
1:A:77:PRO:O	1:A:81:VAL:HG23	2.17	0.45
1:A:175:ILE:O	1:A:175:ILE:HG22	2.17	0.45
1:B:9:LEU:HA	1:B:175:ILE:HD13	1.98	0.45
2:C:172:ASP:HB2	2:C:173:PRO:HD2	1.98	0.45
1:E:13:VAL:CG2	1:E:195:LEU:HD21	2.43	0.45
1:E:80:ARG:HD3	1:E:86:SER:HB2	1.98	0.45
2:H:180:LEU:CD1	2:H:203:VAL:HG23	2.46	0.45
1:E:100:ALA:HB2	1:E:157:LEU:HD21	1.98	0.45
1:F:28:PHE:CZ	1:F:150:ALA:HA	2.52	0.45
2:G:293:ASN:HD21	2:H:285:GLU:N	2.14	0.45
1:A:139:LEU:N	1:A:140:PRO:CD	2.80	0.45
2:C:270:ARG:HB3	2:C:340:LEU:HD11	1.99	0.45
2:H:4:LEU:HD13	2:H:47:LEU:HD11	1.99	0.45
2:H:271:LEU:CD1	2:H:283:LEU:HD21	2.46	0.45
1:B:41:THR:CG2	1:B:53:TYR:HB2	2.47	0.45
2:D:10:VAL:HG11	2:D:17:THR:HG23	1.99	0.45
1:E:2:SER:HB2	1:E:4:PRO:HD2	1.99	0.45
1:E:144:PRO:O	1:E:148:ASN:HB2	2.17	0.45
1:F:62:ILE:O	1:F:65:SER:HB3	2.17	0.45
2:H:165:ASP:O	2:H:166:GLN:C	2.59	0.45
2:D:11:PHE:CE2	4:D:2000:ADP:C5	3.04	0.44
2:G:292:VAL:O	2:G:292:VAL:HG13	2.17	0.44
2:H:68:GLU:O	2:H:72:LEU:HD13	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:312:MET:C	2:C:313:LEU:HD12	2.42	0.44
1:F:127:GLY:O	2:H:78:THR:HG23	2.16	0.44
2:H:232:PRO:HB2	2:H:238:GLN:HG3	1.98	0.44
2:H:300:GLN:NE2	2:H:300:GLN:HA	2.32	0.44
2:H:314:THR:HG23	2:H:314:THR:O	2.17	0.44
1:B:13:VAL:HG22	1:B:172:LEU:CD1	2.47	0.44
2:C:52:ASN:HD22	2:C:54:LEU:HD11	1.83	0.44
2:H:0:HIS:HB3	2:H:29:PRO:HA	1.99	0.44
1:B:4:PRO:HG2	1:B:5:MET:H	1.83	0.44
2:D:82:ARG:HH21	2:D:110:ASP:HB2	1.82	0.44
1:F:13:VAL:N	1:F:172:LEU:HD11	2.32	0.44
1:B:147:VAL:HG11	1:B:210:ILE:HG23	2.00	0.44
2:D:10:VAL:HG13	2:D:18:ILE:O	2.18	0.44
2:D:93:LEU:HD22	2:D:102:ASN:HD21	1.82	0.44
2:D:114:LYS:HA	2:D:117:VAL:HG22	1.99	0.44
2:D:310:GLY:HA3	3:D:1000:MSE:HE2	1.98	0.44
2:H:72:LEU:HG	2:H:76:GLU:CB	2.48	0.44
2:C:220:ILE:HD13	2:C:236:LEU:HD23	1.99	0.44
2:C:231:HIS:N	2:C:232:PRO:CD	2.79	0.44
2:D:234:THR:CG2	2:D:235:PRO:HD2	2.47	0.44
2:D:105:LEU:N	2:D:106:PRO:CD	2.80	0.44
2:D:249:ILE:HG12	2:D:313:LEU:CD1	2.47	0.44
1:F:106:ARG:HA	1:F:106:ARG:HD3	1.65	0.44
1:B:4:PRO:CD	1:B:5:MET:H	2.31	0.44
2:C:201:MET:HE2	2:C:240:PHE:CB	2.47	0.44
1:F:121:GLU:OE1	2:H:49:ARG:NH2	2.51	0.44
2:G:78:THR:O	2:G:82:ARG:HG3	2.17	0.44
2:H:260:GLU:O	2:H:262:PHE:HD2	2.01	0.44
1:B:123:SER:HA	1:B:126:MET:HE2	1.99	0.43
1:B:134:VAL:HA	1:B:138:LEU:HB2	1.98	0.43
2:D:230:SER:OG	2:D:231:HIS:ND1	2.21	0.43
2:D:253:TYR:O	2:D:257:LEU:N	2.51	0.43
1:F:116:PRO:C	1:F:118:GLY:N	2.74	0.43
2:G:251:GLU:OE2	2:H:206:ARG:NH2	2.48	0.43
2:H:112:THR:CB	2:H:117:VAL:HG23	2.48	0.43
2:H:333:HIS:C	2:H:334:HIS:CG	2.96	0.43
1:B:4:PRO:HG2	1:B:5:MET:N	2.33	0.43
1:B:129:THR:O	1:B:133:ILE:N	2.44	0.43
1:B:154:LEU:HD13	1:B:154:LEU:HA	1.87	0.43
2:C:56:ARG:NH1	2:C:70:THR:O	2.50	0.43
1:E:137:VAL:HG23	1:E:138:LEU:N	2.33	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:G:275:GLY:O	2:G:278:VAL:HG23	2.18	0.43
1:A:19:MET:HG3	1:A:89:LEU:CD2	2.48	0.43
1:B:72:LEU:CG	1:B:167:VAL:HG11	2.48	0.43
2:C:83:GLN:HA	2:C:159:PRO:HA	2.00	0.43
2:C:116:GLU:O	2:C:120:ARG:HG2	2.18	0.43
2:D:22:ASN:H	2:D:217:GLY:C	2.23	0.43
1:E:69:ILE:HD12	1:F:163:MET:HA	2.01	0.43
1:E:69:ILE:CG1	1:E:163:MET:HE2	2.46	0.43
1:F:94:VAL:HB	1:F:95:PRO:HD3	2.01	0.43
2:H:266:VAL:CG1	2:H:316:MET:C	2.92	0.43
2:H:303:TYR:CZ	2:H:308:LYS:HB2	2.51	0.43
1:B:134:VAL:HA	1:B:138:LEU:HD12	1.99	0.43
1:B:167:VAL:O	1:B:167:VAL:CG2	2.67	0.43
1:B:188:VAL:CA	1:B:191:THR:HG22	2.46	0.43
2:D:312:MET:HE3	2:D:312:MET:HB2	1.83	0.43
2:G:269:LEU:HD23	2:G:339:VAL:HG13	2.01	0.43
2:H:195:LEU:C	2:H:195:LEU:HD13	2.43	0.43
1:A:85:THR:HG22	1:A:87:ILE:H	1.84	0.43
1:B:2:SER:HB3	1:B:4:PRO:CD	2.47	0.43
1:E:118:GLY:C	1:E:120:ILE:N	2.72	0.43
2:G:281:PRO:CB	2:H:293:ASN:HB3	2.48	0.43
2:G:334:HIS:N	2:G:334:HIS:CD2	2.84	0.43
1:A:59:ILE:HG13	1:A:60:VAL:N	2.33	0.43
1:A:178:GLN:HA	1:A:182:ILE:HD12	2.01	0.43
1:B:42:ARG:HG2	1:B:43:PRO:N	2.34	0.43
1:B:52:LEU:O	1:B:56:VAL:HG23	2.18	0.43
2:G:145:GLN:HA	2:G:148:ARG:HB2	2.00	0.43
2:G:251:GLU:HG2	2:H:177:ARG:HH21	1.83	0.43
2:H:288:ARG:HG3	2:H:289:ARG:N	2.34	0.43
2:C:1:MET:HG2	2:C:193:THR:HG21	2.00	0.43
2:C:10:VAL:C	2:C:11:PHE:HD1	2.27	0.43
1:F:47:ILE:O	1:F:47:ILE:CG1	2.63	0.43
2:G:91:PHE:O	2:G:92:ASN:CB	2.62	0.43
1:B:8:LEU:O	1:B:10:VAL:N	2.52	0.43
1:B:66:ILE:HA	1:B:67:PRO:HD3	1.86	0.43
2:D:247:LEU:O	2:D:247:LEU:HG	2.19	0.43
1:E:162:ALA:HB1	1:F:70:ILE:HD13	2.01	0.43
1:F:13:VAL:HG12	1:F:172:LEU:HD21	2.00	0.43
1:B:8:LEU:C	1:B:10:VAL:N	2.76	0.43
1:B:41:THR:HA	1:B:47:ILE:HB	2.01	0.43
2:C:21:LEU:HD23	2:C:217:GLY:HA2	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:15:GLU:CD	1:F:89:LEU:HD21	2.43	0.43
2:H:48:ILE:HD11	2:H:163:LEU:HB3	2.00	0.43
2:D:56:ARG:HG3	2:D:57:PRO:HD2	2.00	0.43
1:E:134:VAL:HG23	1:E:135:ARG:N	2.34	0.43
1:F:61:ASN:OD1	1:F:106:ARG:NH2	2.45	0.43
2:G:168:THR:HA	2:G:171:LEU:HD12	2.00	0.43
2:H:112:THR:C	2:H:117:VAL:HG23	2.44	0.43
2:C:53:LEU:HD11	2:C:69:LEU:CB	2.48	0.42
2:D:61:SER:HB2	2:D:68:GLU:OE2	2.19	0.42
1:F:177:TYR:O	1:F:181:TYR:HB3	2.19	0.42
2:G:24:VAL:HG11	2:G:47:LEU:CD2	2.49	0.42
2:G:124:LEU:HD12	2:G:124:LEU:N	2.34	0.42
2:H:21:LEU:HD11	2:H:47:LEU:HA	2.01	0.42
2:H:91:PHE:CZ	2:H:171:LEU:HD21	2.51	0.42
1:B:4:PRO:CG	1:B:5:MET:N	2.81	0.42
2:C:182:LEU:CD1	2:C:186:ILE:CD1	2.96	0.42
2:G:52:ASN:HB3	2:G:54:LEU:HG	2.00	0.42
1:A:80:ARG:CG	1:A:86:SER:HB2	2.49	0.42
1:B:133:ILE:O	1:B:137:VAL:CG1	2.66	0.42
1:B:188:VAL:HA	1:B:191:THR:CG2	2.46	0.42
2:C:11:PHE:HD2	4:C:2000:ADP:N3	2.17	0.42
2:C:130:LEU:C	2:C:132:ASP:N	2.76	0.42
2:C:339:VAL:HG12	2:C:341:GLY:H	1.83	0.42
2:D:34:TYR:CD1	2:D:34:TYR:C	2.97	0.42
1:E:17:LEU:HD21	1:E:199:LEU:CD2	2.49	0.42
1:E:162:ALA:HB1	1:F:70:ILE:CD1	2.49	0.42
1:F:38:LEU:HD13	1:F:56:VAL:HG12	2.01	0.42
2:G:98:THR:HG23	2:G:101:GLY:N	2.34	0.42
2:H:272:GLU:CD	2:H:336:LYS:HD3	2.43	0.42
1:B:4:PRO:CG	1:B:5:MET:H	2.32	0.42
1:B:45:GLN:CG	1:B:46:ILE:N	2.82	0.42
2:C:10:VAL:O	2:C:11:PHE:HD1	2.02	0.42
2:D:201:MET:HE2	2:D:241:ILE:HD13	2.01	0.42
1:F:80:ARG:CB	1:F:86:SER:HB3	2.43	0.42
2:G:330:LEU:O	2:G:335:VAL:HB	2.19	0.42
2:H:86:MET:HG3	2:H:163:LEU:HB2	2.00	0.42
1:B:2:SER:C	1:B:4:PRO:CD	2.77	0.42
2:D:302:ASP:O	2:D:309:PHE:HB2	2.19	0.42
1:E:132:GLN:O	1:E:136:LYS:HB2	2.20	0.42
1:E:143:LEU:HA	1:E:143:LEU:HD12	1.79	0.42
2:G:9:LYS:CE	2:G:56:ARG:O	2.59	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:G:236:LEU:HA	2:G:239:LYS:HD2	2.00	0.42
1:A:178:GLN:HB2	1:A:179:TYR:CD1	2.54	0.42
1:B:40:VAL:CG1	1:B:47:ILE:HD12	2.47	0.42
1:B:75:MET:O	1:B:79:THR:N	2.40	0.42
1:B:157:LEU:HA	1:B:160:TYR:HB2	2.01	0.42
2:C:3:LYS:CG	2:C:27:HIS:CD2	3.02	0.42
2:C:227:GLU:O	2:C:230:SER:O	2.38	0.42
2:D:7:ILE:HA	2:D:60:GLY:HA3	2.01	0.42
2:G:229:PHE:HA	2:G:241:ILE:CD1	2.49	0.42
1:A:26:PHE:HB2	1:A:97:THR:HG21	2.02	0.42
1:A:155:ILE:HG13	1:A:203:ILE:CG2	2.49	0.42
2:H:123:GLU:O	2:H:127:LEU:HG	2.20	0.42
1:A:20:THR:HG21	1:A:158:VAL:HA	2.01	0.42
2:H:251:GLU:HA	2:H:254:GLN:HB3	2.01	0.42
1:A:104:ILE:O	1:A:108:VAL:HG23	2.20	0.42
1:B:20:THR:HG22	1:B:154:LEU:CD1	2.42	0.42
2:C:72:LEU:HG	2:C:76:GLU:HB3	2.02	0.42
2:D:1:MET:O	2:D:64:VAL:HA	2.19	0.42
1:E:116:PRO:O	1:E:119:LEU:HB2	2.20	0.42
2:H:93:LEU:HD12	2:H:146:LYS:HB2	2.01	0.42
2:H:114:LYS:CG	2:H:115:ASP:H	2.12	0.42
2:H:144:GLY:O	2:H:148:ARG:HG3	2.20	0.42
1:B:115:ILE:HA	1:B:116:PRO:HD3	1.86	0.42
1:B:130:PRO:O	1:B:134:VAL:HG23	2.20	0.42
1:B:42:ARG:CG	1:B:43:PRO:CD	2.97	0.41
2:C:89:GLN:HB2	2:C:166:GLN:HB2	2.02	0.41
2:C:283:LEU:HA	2:C:286:THR:HG22	2.02	0.41
1:F:2:SER:O	1:F:6:MET:HG2	2.20	0.41
1:F:13:VAL:HG23	1:F:14:TRP:N	2.35	0.41
2:G:103:VAL:O	2:G:106:PRO:HD2	2.20	0.41
2:H:319:THR:HG23	2:H:322:ASP:CB	2.44	0.41
1:F:116:PRO:C	1:F:118:GLY:H	2.28	0.41
2:G:33:ILE:HG12	2:G:194:ILE:HB	2.01	0.41
2:G:214:ILE:HA	2:G:218:GLU:O	2.20	0.41
1:A:157:LEU:HA	1:A:157:LEU:HD23	1.81	0.41
1:B:147:VAL:HB	1:B:211:VAL:HG22	2.00	0.41
1:B:210:ILE:O	1:B:214:VAL:HG12	2.21	0.41
2:C:44:LYS:CE	4:C:2000:ADP:O3B	2.67	0.41
2:D:24:VAL:HG13	2:D:219:LEU:HD12	2.01	0.41
2:D:195:LEU:HD13	2:D:195:LEU:C	2.46	0.41
1:E:195:LEU:HD23	1:E:195:LEU:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:150:ALA:O	1:F:153:THR:HB	2.21	0.41
2:H:90:HIS:C	2:H:91:PHE:CD1	2.98	0.41
2:H:107:LEU:HA	2:H:110:ASP:OD1	2.20	0.41
2:H:241:ILE:O	2:H:244:THR:HG22	2.20	0.41
1:A:68:PHE:CZ	1:A:96:LEU:HD12	2.46	0.41
1:A:75:MET:HE2	1:A:75:MET:HB3	1.96	0.41
2:C:11:PHE:CD2	4:C:2000:ADP:C4	3.08	0.41
2:C:182:LEU:HD13	2:C:182:LEU:O	2.20	0.41
2:G:99:VAL:HG12	2:G:141:LEU:HD11	1.96	0.41
2:H:28:VAL:HA	2:H:29:PRO:HD2	1.92	0.41
2:H:298:SER:O	2:H:312:MET:HB2	2.21	0.41
1:A:71:LEU:C	1:A:71:LEU:HD13	2.46	0.41
2:G:99:VAL:HG21	2:G:133:LYS:O	2.21	0.41
2:H:257:LEU:HD12	2:H:258:GLN:H	1.84	0.41
1:A:104:ILE:HD11	1:A:153:THR:CG2	2.51	0.41
1:F:54:ARG:HD2	1:F:54:ARG:HA	1.85	0.41
2:G:53:LEU:HD23	2:G:53:LEU:HA	1.84	0.41
2:H:300:GLN:HE21	2:H:300:GLN:HA	1.86	0.41
1:A:184:TYR:OH	1:A:186:ALA:HA	2.21	0.41
2:D:104:ALA:O	2:D:108:GLU:HG3	2.20	0.41
2:G:4:LEU:O	2:G:4:LEU:HD12	2.20	0.41
2:H:180:LEU:HD11	2:H:203:VAL:HG23	2.02	0.41
2:H:200:GLU:O	2:H:203:VAL:HG12	2.21	0.41
1:A:70:ILE:HG22	1:B:193:LEU:HD12	2.03	0.41
1:E:125:ALA:CB	2:G:52:ASN:HD21	2.32	0.41
1:F:30:ILE:CD1	1:F:101:ALA:HB2	2.50	0.41
1:F:80:ARG:O	1:F:84:GLY:HA2	2.20	0.41
1:F:118:GLY:O	1:F:119:LEU:C	2.62	0.41
1:F:122:ALA:O	1:F:126:MET:HG3	2.21	0.41
2:G:142:SER:O	2:G:143:GLY:C	2.63	0.41
2:G:241:ILE:HA	2:G:244:THR:HG22	2.03	0.41
2:H:77:LEU:HD21	2:H:81:ARG:NH1	2.35	0.41
1:A:13:VAL:CG2	1:A:172:LEU:HD11	2.45	0.41
1:A:15:GLU:HB2	1:A:171:GLY:HA2	2.03	0.41
1:A:41:THR:HA	1:A:47:ILE:O	2.21	0.41
1:B:13:VAL:HG22	1:B:172:LEU:HD12	2.03	0.41
1:B:39:TYR:HE1	1:B:113:LEU:HA	1.85	0.41
1:B:42:ARG:HG2	1:B:43:PRO:HD2	2.03	0.41
1:B:104:ILE:O	1:B:107:MET:N	2.54	0.41
2:C:180:LEU:O	2:C:207:ILE:HG21	2.20	0.41
2:C:295:ASN:HB3	2:C:315:GLU:HG3	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:296:ILE:HD11	2:D:312:MET:SD	2.61	0.41
2:D:230:SER:CB	2:D:309:PHE:HE2	2.34	0.41
1:E:9:LEU:O	1:E:13:VAL:HG23	2.21	0.41
1:E:15:GLU:OE2	1:E:171:GLY:N	2.54	0.41
1:E:94:VAL:HB	1:E:95:PRO:HD3	2.02	0.41
2:G:56:ARG:HA	2:G:57:PRO:HD3	1.87	0.41
2:G:296:ILE:HG12	3:H:1000:MSE:OXT	2.20	0.41
2:H:93:LEU:HA	2:H:93:LEU:HD23	1.89	0.41
2:H:266:VAL:HG21	2:H:317:HIS:HA	2.03	0.41
1:A:49:ASN:HB2	1:A:52:LEU:HB3	2.03	0.41
1:B:162:ALA:HA	1:B:173:GLY:HA3	2.03	0.41
2:H:49:ARG:HG3	2:H:55:GLU:HB2	2.03	0.41
2:H:244:THR:HG23	2:H:245:LEU:HD13	2.03	0.41
2:C:4:LEU:HD21	2:C:51:VAL:HG22	2.02	0.40
2:C:130:LEU:CD1	2:C:149:VAL:HG22	2.49	0.40
2:G:48:ILE:HD12	2:G:195:LEU:HD12	2.03	0.40
2:G:245:LEU:HD21	2:G:309:PHE:CE2	2.56	0.40
2:H:46:THR:HG23	2:H:55:GLU:HG2	2.03	0.40
2:H:100:PHE:CD1	2:H:100:PHE:C	2.99	0.40
2:H:148:ARG:HA	2:H:151:ILE:HD12	2.02	0.40
2:C:231:HIS:NE2	2:C:274:THR:HG21	2.35	0.40
2:H:280:ALA:O	2:H:282:LEU:N	2.50	0.40
1:A:60:VAL:O	1:A:64:ARG:CB	2.70	0.40
1:A:64:ARG:HG3	1:A:64:ARG:HH11	1.85	0.40
1:B:69:ILE:HD12	1:B:69:ILE:HA	1.93	0.40
1:E:104:ILE:HD13	1:E:104:ILE:HG21	1.80	0.40
2:H:35:GLY:CA	2:H:196:LEU:HD13	2.47	0.40
1:A:184:TYR:CD2	1:A:184:TYR:C	2.99	0.40
2:D:22:ASN:ND2	2:D:218:GLU:OE1	2.55	0.40
2:G:105:LEU:HD23	2:G:105:LEU:C	2.47	0.40
2:H:257:LEU:HG	2:H:258:GLN:N	2.36	0.40
2:H:283:LEU:CD1	2:H:312:MET:SD	3.09	0.40
2:C:187:ASN:O	2:C:191:GLY:N	2.48	0.40
2:C:228:VAL:O	2:C:232:PRO:HG3	2.21	0.40
2:C:283:LEU:C	2:C:286:THR:HG22	2.46	0.40
1:F:39:TYR:OH	1:F:120:ILE:HD11	2.22	0.40
2:G:301:MET:HE2	2:H:296:ILE:O	2.20	0.40
2:H:270:ARG:HE	2:H:311:ILE:HD11	1.87	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	214/217 (99%)	208 (97%)	4 (2%)	2 (1%)	14	42
1	B	214/217 (99%)	206 (96%)	7 (3%)	1 (0%)	24	54
1	E	213/217 (98%)	202 (95%)	11 (5%)	0	100	100
1	F	214/217 (99%)	206 (96%)	8 (4%)	0	100	100
2	C	342/366 (93%)	333 (97%)	9 (3%)	0	100	100
2	D	341/366 (93%)	322 (94%)	19 (6%)	0	100	100
2	G	341/366 (93%)	320 (94%)	20 (6%)	1 (0%)	36	65
2	H	338/366 (92%)	321 (95%)	15 (4%)	2 (1%)	21	50
All	All	2217/2332 (95%)	2118 (96%)	93 (4%)	6 (0%)	36	65

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	H	111	ASN
2	H	113	PRO
1	B	9	LEU
2	G	250	PRO
1	A	93	ILE
1	A	94	VAL

5.3.2 Protein sidechains

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	171/172 (99%)	166 (97%)	5 (3%)	37	60
1	B	171/172 (99%)	167 (98%)	4 (2%)	44	63
1	E	170/172 (99%)	168 (99%)	2 (1%)	63	72
1	F	171/172 (99%)	171 (100%)	0	100	100
2	C	299/319 (94%)	296 (99%)	3 (1%)	68	75
2	D	298/319 (93%)	292 (98%)	6 (2%)	48	65
2	G	298/319 (93%)	286 (96%)	12 (4%)	28	53
2	H	296/319 (93%)	286 (97%)	10 (3%)	32	57
All	All	1874/1964 (95%)	1832 (98%)	42 (2%)	45	63

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	46	ILE
1	A	69	ILE
1	A	89	LEU
1	A	106	ARG
1	A	202	LEU
1	B	30	ILE
1	B	62	ILE
1	B	137	VAL
1	B	179	TYR
2	C	140	ASN
2	C	182	LEU
2	C	269	LEU
2	D	182	LEU
2	D	192	LEU
2	D	195	LEU
2	D	196	LEU
2	D	219	LEU
2	D	283	LEU
1	E	187	THR
1	E	202	LEU
2	G	47	LEU
2	G	48	ILE
2	G	91	PHE
2	G	107	LEU
2	G	145	GLN
2	G	179	ILE
2	G	192	LEU

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Mol	Chain	Res	Type
2	G	196	LEU
2	G	198	THR
2	G	278	VAL
2	G	327	ILE
2	G	339	VAL
2	H	0	HIS
2	H	47	LEU
2	H	182	LEU
2	H	192	LEU
2	H	195	LEU
2	H	196	LEU
2	H	222	GLN
2	H	260	GLU
2	H	291	ASN
2	H	312	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (37) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	174	GLN
1	A	178	GLN
1	B	45	GLN
2	C	13	GLN
2	D	22	ASN
2	D	27	HIS
2	D	102	ASN
2	D	134	HIS
2	D	291	ASN
2	D	331	GLN
2	D	333	HIS
1	E	178	GLN
1	F	45	GLN
1	F	110	ASN
1	F	132	GLN
1	F	204	GLN
2	G	13	GLN
2	G	52	ASN
2	G	83	GLN
2	G	92	ASN
2	G	140	ASN
2	G	145	GLN
2	G	147	GLN

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Mol	Chain	Res	Type
2	G	166	GLN
2	G	199	HIS
2	G	276	GLN
2	G	334	HIS
2	H	67	GLN
2	H	89	GLN
2	H	92	ASN
2	H	147	GLN
2	H	246	HIS
2	H	254	GLN
2	H	258	GLN
2	H	294	ASN
2	H	300	GLN
2	H	324	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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
3	MSE	C	1000	-	5,8,8	1.15	1 (20%)	3,9,9	1.80	1 (33%)
4	ADP	C	2000	-	28,29,29	1.65	4 (14%)	43,45,45	1.88	7 (16%)
3	MSE	G	1000	-	5,8,8	1.07	1 (20%)	3,9,9	2.23	1 (33%)
4	ADP	D	2000	-	28,29,29	1.69	4 (14%)	43,45,45	1.80	10 (23%)
3	MSE	D	1000	-	5,8,8	0.94	0	3,9,9	1.82	1 (33%)
4	ADP	G	2000	-	28,29,29	1.74	4 (14%)	43,45,45	1.77	11 (25%)
4	ADP	H	2000	-	28,29,29	1.90	4 (14%)	43,45,45	1.98	10 (23%)
3	MSE	H	1000	-	5,8,8	1.18	1 (20%)	3,9,9	1.70	1 (33%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	MSE	C	1000	-	-	2/7/8/8	-
4	ADP	C	2000	-	-	8/16/32/32	0/3/3/3
3	MSE	G	1000	-	-	0/7/8/8	-
4	ADP	D	2000	-	-	4/16/32/32	0/3/3/3
3	MSE	D	1000	-	-	3/7/8/8	-
4	ADP	G	2000	-	-	1/16/32/32	0/3/3/3
4	ADP	H	2000	-	-	1/16/32/32	0/3/3/3
3	MSE	H	1000	-	-	6/7/8/8	-

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	H	2000	ADP	C5-C4	5.81	1.49	1.39
4	G	2000	ADP	C5-C4	5.70	1.49	1.39
4	C	2000	ADP	C5-C4	5.60	1.49	1.39
4	H	2000	ADP	PA-O3A	5.04	1.64	1.59
4	D	2000	ADP	C5-C4	5.04	1.48	1.39
4	D	2000	ADP	PA-O3A	4.56	1.64	1.59
4	G	2000	ADP	PA-O3A	3.90	1.63	1.59
4	H	2000	ADP	C5-C6	3.79	1.51	1.41
4	C	2000	ADP	C5-C6	3.66	1.51	1.41
4	G	2000	ADP	C5-C6	3.38	1.50	1.41
4	D	2000	ADP	C5-C6	3.01	1.49	1.41
4	H	2000	ADP	C8-N7	2.89	1.37	1.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	2000	ADP	PA-O3A	2.77	1.62	1.59
4	C	2000	ADP	C8-N7	2.52	1.36	1.31
4	G	2000	ADP	C8-N7	2.51	1.36	1.31
3	H	1000	MSE	OXT-C	-2.48	1.22	1.30
4	D	2000	ADP	C8-N7	2.47	1.36	1.31
3	C	1000	MSE	OXT-C	-2.47	1.22	1.30
3	G	1000	MSE	OXT-C	-2.32	1.23	1.30

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	2000	ADP	C5-C4-N3	-6.43	117.86	126.72
4	H	2000	ADP	C5-C4-N3	-6.13	118.28	126.72
4	G	2000	ADP	C5-C4-N3	-5.42	119.25	126.72
4	D	2000	ADP	C5-C4-N3	-5.05	119.76	126.72
4	C	2000	ADP	C4-C5-N7	-4.83	105.06	110.58
4	H	2000	ADP	N3-C4-N9	4.50	134.82	127.17
4	C	2000	ADP	N3-C4-N9	4.49	134.79	127.17
4	H	2000	ADP	C4-C5-N7	-4.44	105.51	110.58
4	G	2000	ADP	N3-C4-N9	4.22	134.35	127.17
4	D	2000	ADP	N3-C4-N9	4.11	134.16	127.17
4	H	2000	ADP	C2-N3-C4	3.94	121.46	111.83
4	G	2000	ADP	C4-C5-N7	-3.72	106.32	110.58
3	G	1000	MSE	OXT-C-O	-3.68	115.73	124.08
4	D	2000	ADP	N3-C2-N1	-3.58	123.16	128.58
4	D	2000	ADP	C2-N3-C4	3.56	120.53	111.83
4	C	2000	ADP	C2-N3-C4	3.54	120.47	111.83
4	D	2000	ADP	C4-C5-N7	-3.49	106.59	110.58
4	G	2000	ADP	C2-N3-C4	3.40	120.13	111.83
4	H	2000	ADP	N3-C2-N1	-3.33	123.54	128.58
4	C	2000	ADP	C5-N7-C8	3.14	108.38	103.45
3	D	1000	MSE	OXT-C-O	-3.14	116.97	124.08
4	D	2000	ADP	C4-N9-C8	3.13	109.02	105.74
3	C	1000	MSE	OXT-C-O	-3.12	117.00	124.08
4	H	2000	ADP	C5-N7-C8	3.10	108.32	103.45
4	G	2000	ADP	N3-C2-N1	-2.91	124.17	128.58
3	H	1000	MSE	OXT-C-O	-2.85	117.61	124.08
4	D	2000	ADP	C5-N7-C8	2.60	107.53	103.45
4	H	2000	ADP	C6-C5-N7	2.58	137.06	132.09
4	G	2000	ADP	C5-N7-C8	2.53	107.43	103.45
4	D	2000	ADP	C6-C5-N7	2.51	136.92	132.09
4	H	2000	ADP	C4-N9-C8	2.40	108.26	105.74

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	2000	ADP	C6-C5-N7	2.37	136.66	132.09
4	D	2000	ADP	C2-N1-C6	2.35	122.59	118.73
4	G	2000	ADP	C6-C5-N7	2.31	136.54	132.09
4	G	2000	ADP	C4-N9-C8	2.25	108.10	105.74
4	D	2000	ADP	N9-C8-N7	-2.22	110.78	113.94
4	C	2000	ADP	N3-C2-N1	-2.17	125.30	128.58
4	G	2000	ADP	C3'-C2'-C1'	2.16	105.55	101.46
4	G	2000	ADP	O4'-C1'-N9	2.16	112.23	108.09
4	H	2000	ADP	C3'-C2'-C1'	2.08	105.40	101.46
4	H	2000	ADP	N9-C8-N7	-2.06	111.02	113.94
4	G	2000	ADP	C2-N1-C6	2.05	122.10	118.73

There are no chirality outliers.

All (25) torsion outliers are listed below:

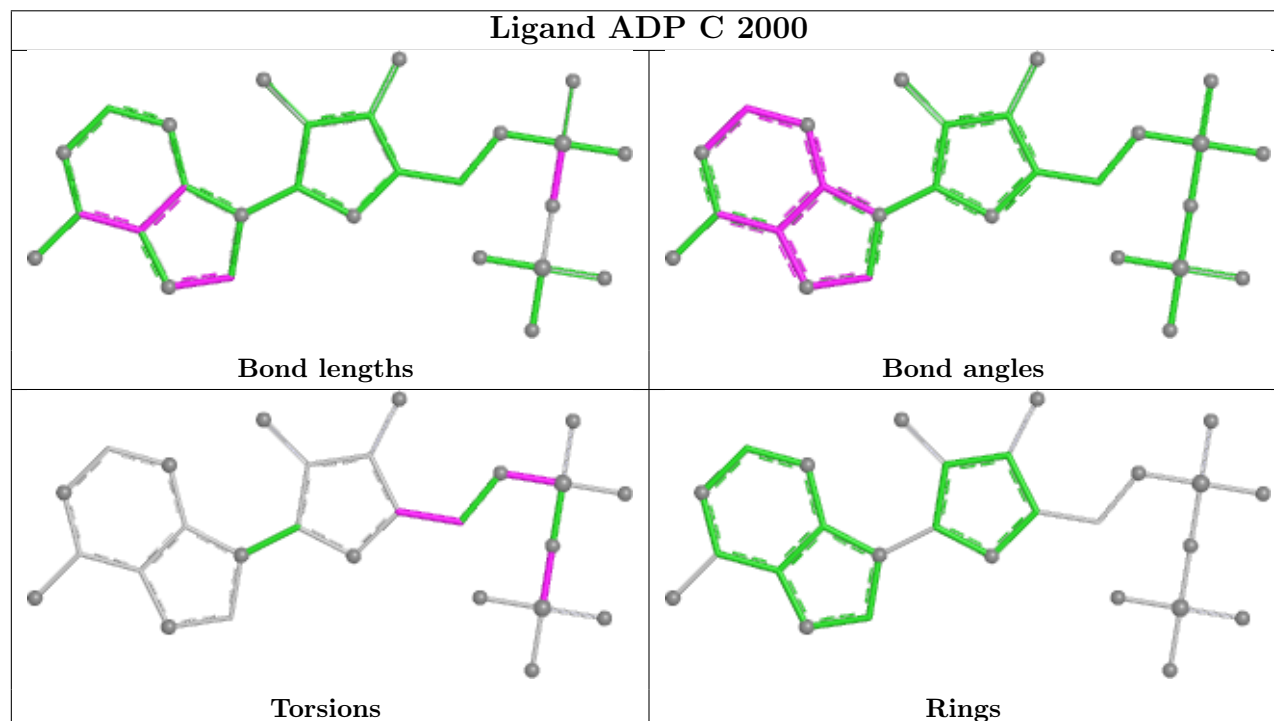
Mol	Chain	Res	Type	Atoms
3	H	1000	MSE	O-C-CA-N
3	H	1000	MSE	N-CA-CB-CG
4	C	2000	ADP	PA-O3A-PB-O2B
4	C	2000	ADP	C5'-O5'-PA-O1A
4	C	2000	ADP	C5'-O5'-PA-O2A
4	C	2000	ADP	C5'-O5'-PA-O3A
4	C	2000	ADP	O4'-C4'-C5'-O5'
3	H	1000	MSE	OXT-C-CA-N
4	C	2000	ADP	C3'-C4'-C5'-O5'
3	C	1000	MSE	OXT-C-CA-N
4	D	2000	ADP	O4'-C4'-C5'-O5'
4	D	2000	ADP	PB-O3A-PA-O1A
3	C	1000	MSE	O-C-CA-N
3	D	1000	MSE	CA-CB-CG-SE
3	H	1000	MSE	C-CA-CB-CG
4	G	2000	ADP	C5'-O5'-PA-O1A
3	D	1000	MSE	O-C-CA-CB
4	D	2000	ADP	PB-O3A-PA-O2A
3	D	1000	MSE	OXT-C-CA-CB
3	H	1000	MSE	OXT-C-CA-CB
4	C	2000	ADP	PA-O3A-PB-O3B
4	D	2000	ADP	C2'-C1'-N9-C8
4	H	2000	ADP	PB-O3A-PA-O1A
4	C	2000	ADP	PA-O3A-PB-O1B
3	H	1000	MSE	O-C-CA-CB

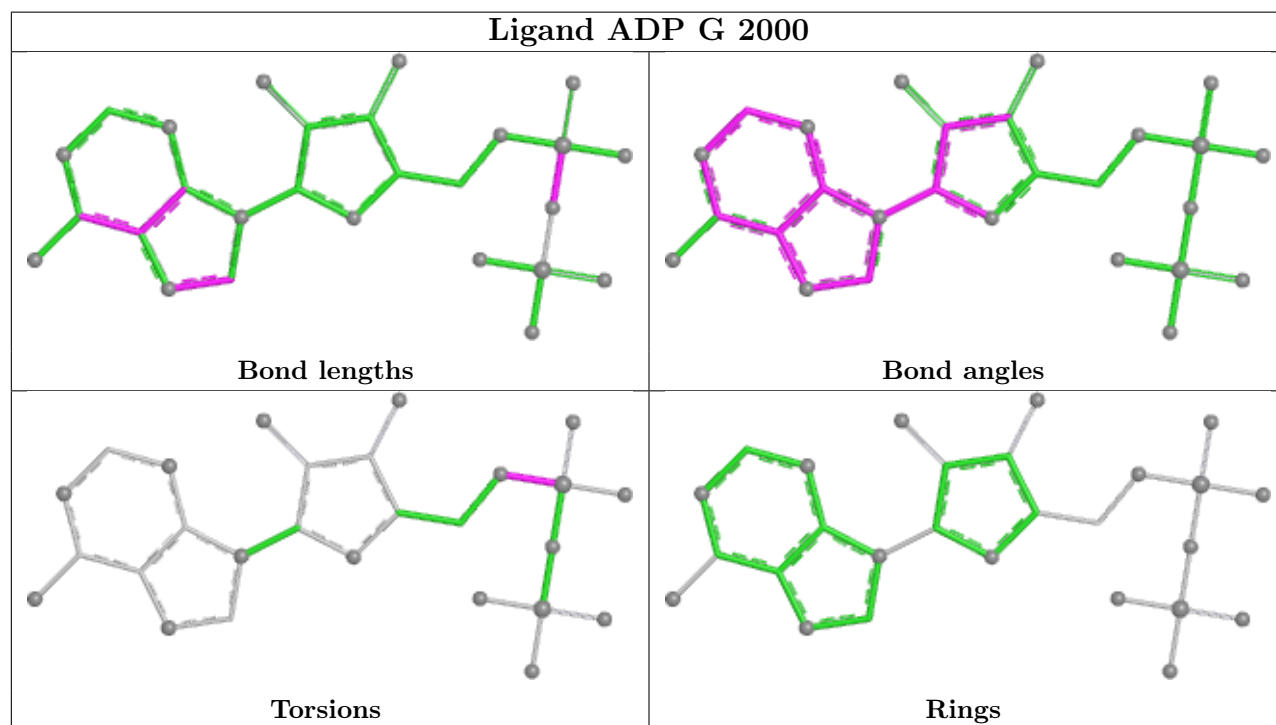
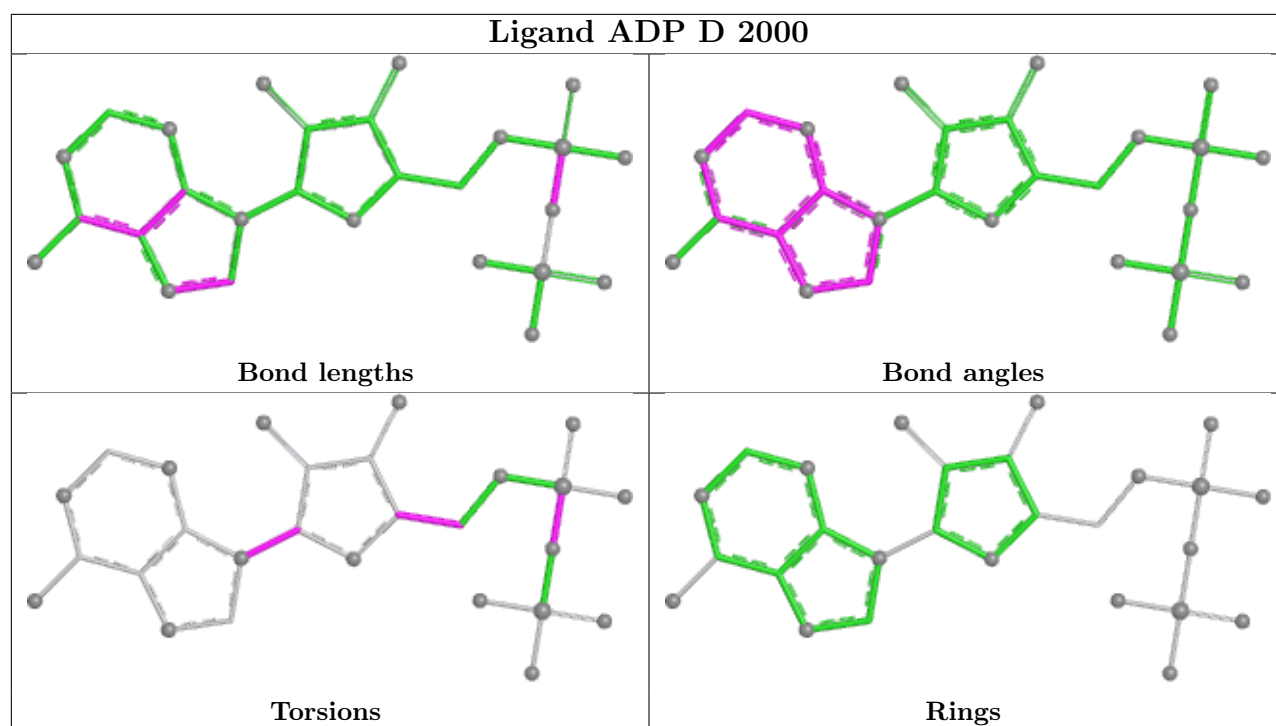
There are no ring outliers.

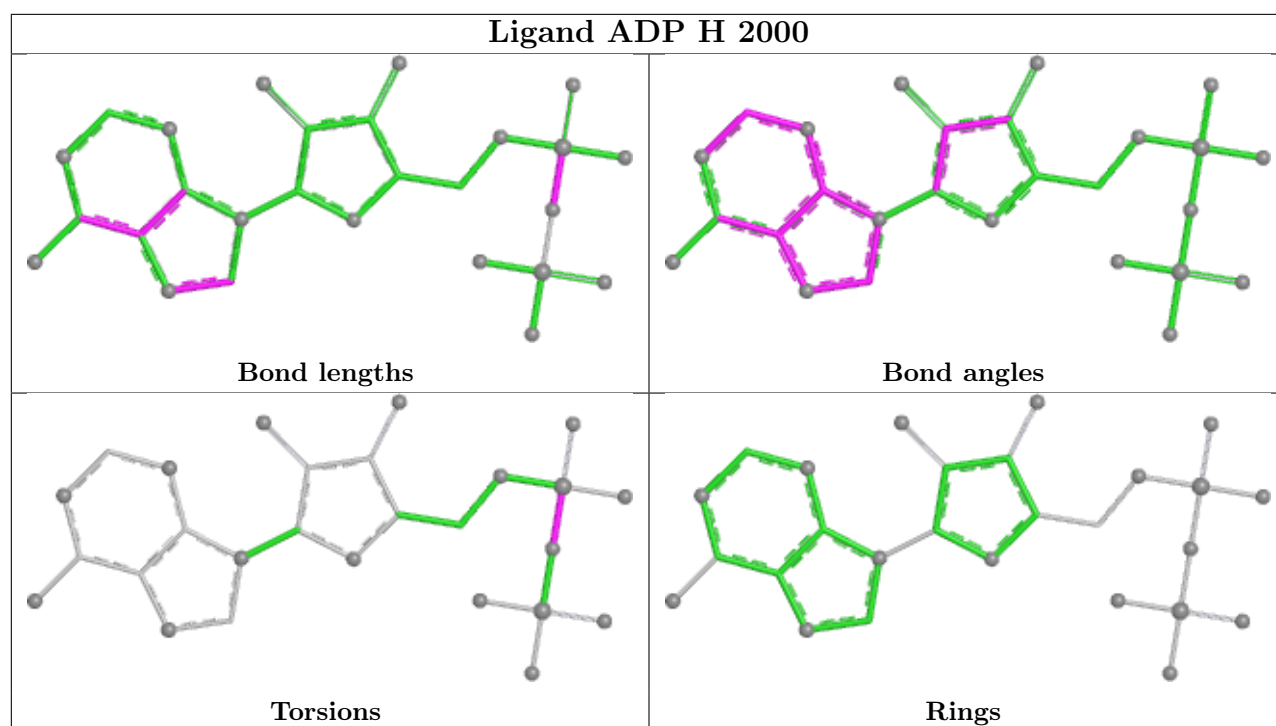
8 monomers are involved in 34 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	1000	MSE	2	0
4	C	2000	ADP	12	0
3	G	1000	MSE	2	0
4	D	2000	ADP	5	0
3	D	1000	MSE	1	0
4	G	2000	ADP	1	0
4	H	2000	ADP	5	0
3	H	1000	MSE	6	0

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







5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	216/217 (99%)	0.05	1 (0%) 87 78	43, 67, 100, 113	0
1	B	216/217 (99%)	0.13	1 (0%) 87 78	48, 70, 109, 150	0
1	E	215/217 (99%)	0.08	1 (0%) 87 78	48, 72, 98, 119	0
1	F	216/217 (99%)	0.17	8 (3%) 45 32	51, 68, 104, 148	0
2	C	344/366 (93%)	0.11	3 (0%) 81 68	38, 73, 121, 146	0
2	D	343/366 (93%)	0.63	22 (6%) 25 20	48, 89, 150, 198	0
2	G	343/366 (93%)	0.20	10 (2%) 53 39	41, 69, 112, 144	0
2	H	340/366 (92%)	0.48	20 (5%) 28 22	39, 83, 152, 168	0
All	All	2233/2332 (95%)	0.26	66 (2%) 52 39	38, 74, 130, 198	0

All (66) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	170	GLY	5.2
2	D	239	LYS	4.3
2	D	81	ARG	4.2
2	G	2	ILE	4.0
2	D	248	ASP	3.9
2	H	331	GLN	3.7
2	G	139	SER	3.7
2	D	249	ILE	3.6
2	D	250	PRO	3.6
2	D	245	LEU	3.4
2	H	249	ILE	3.4
2	H	257	LEU	3.3
2	H	243	SER	3.2
2	D	241	ILE	3.2
2	H	290	PHE	3.0
2	H	261	PRO	3.0

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Mol	Chain	Res	Type	RSRZ
2	H	252	ASP	3.0
2	G	32	GLN	3.0
2	G	258	GLN	3.0
1	F	15	GLU	2.9
2	H	255	GLU	2.9
2	D	229	PHE	2.9
2	D	313	LEU	2.9
2	D	18	ILE	2.8
2	D	247	LEU	2.8
2	H	267	PRO	2.8
2	H	248	ASP	2.8
2	D	284	SER	2.7
2	G	284	SER	2.7
2	H	284	SER	2.7
2	G	30	ALA	2.7
2	C	284	SER	2.7
2	G	71	THR	2.6
2	H	185	ASP	2.6
2	D	132	ASP	2.6
1	E	11	ARG	2.6
1	B	46	ILE	2.5
2	H	332	GLU	2.5
2	D	240	PHE	2.5
2	H	94	LEU	2.5
2	G	259	ALA	2.4
2	H	334	HIS	2.4
2	D	85	GLY	2.4
1	F	5	MET	2.4
2	H	134	HIS	2.4
1	F	89	LEU	2.4
2	D	122	THR	2.3
2	D	25	SER	2.3
2	G	261	PRO	2.3
1	F	19	MET	2.3
1	F	4	PRO	2.2
1	F	7	TRP	2.2
2	D	52	ASN	2.2
2	H	266	VAL	2.1
2	D	252	ASP	2.1
2	D	10	VAL	2.1
2	D	297	ILE	2.1
1	F	90	GLN	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	78	PHE	2.1
2	C	340	LEU	2.1
2	G	14	GLY	2.1
2	H	190	LEU	2.1
2	H	247	LEU	2.1
2	C	123	GLU	2.1
2	H	199	HIS	2.0
2	D	135	ASP	2.0

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

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

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

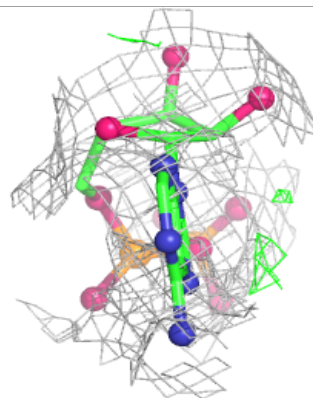
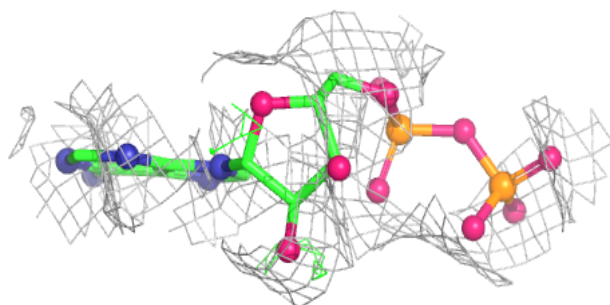
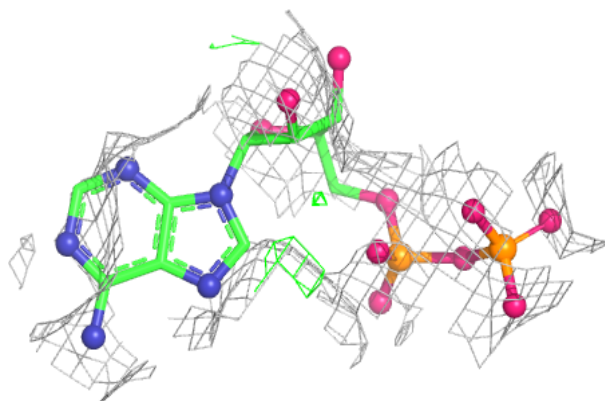
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	ADP	D	2000	27/27	0.66	0.14	103,152,169,172	0
4	ADP	G	2000	27/27	0.84	0.12	84,124,142,145	0
3	MSE	H	1000	9/9	0.91	0.15	89,98,115,130	0
4	ADP	C	2000	27/27	0.91	0.08	82,118,144,149	0
4	ADP	H	2000	27/27	0.94	0.08	69,84,91,94	0
3	MSE	D	1000	9/9	0.95	0.12	73,75,78,123	0
3	MSE	G	1000	9/9	0.96	0.11	61,71,77,106	0
3	MSE	C	1000	9/9	0.96	0.11	77,84,87,113	0

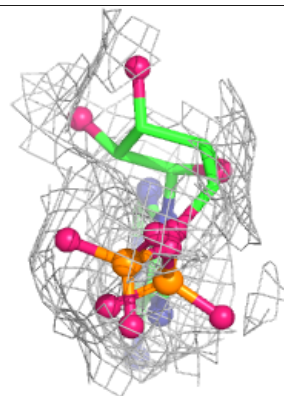
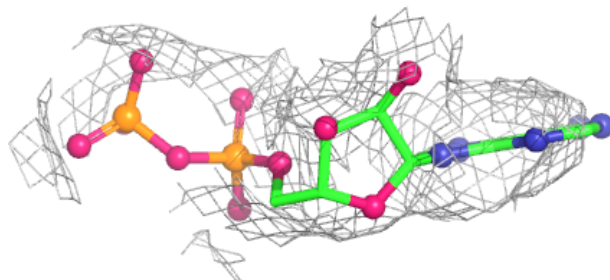
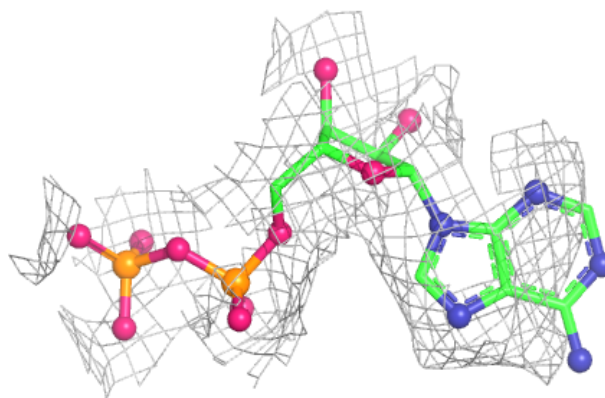
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around ADP D 2000:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

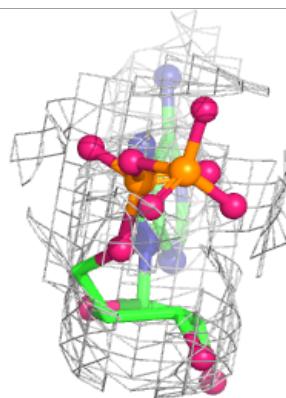
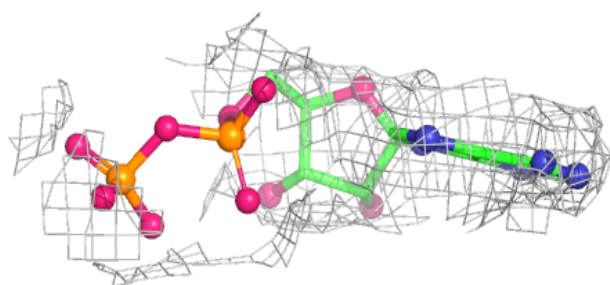
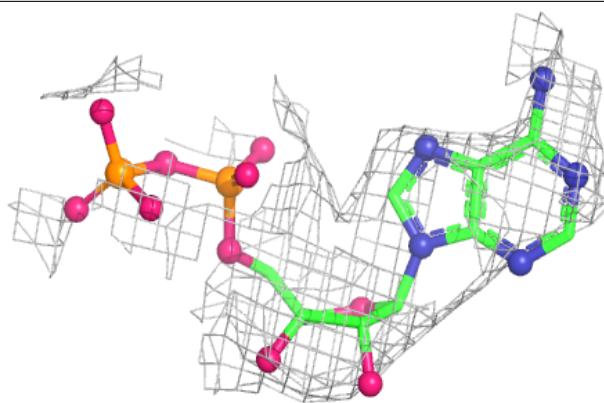
**Electron density around ADP G 2000:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

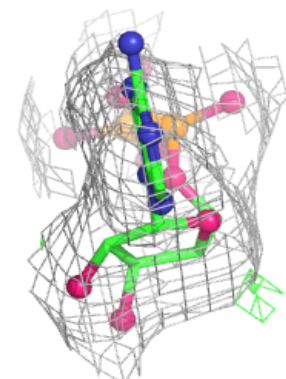
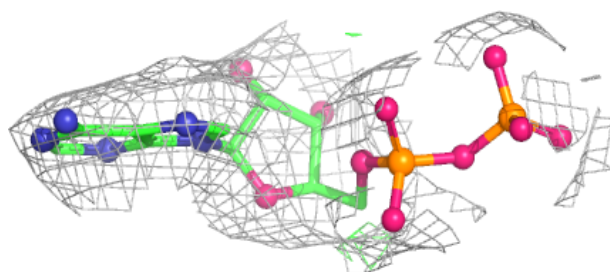
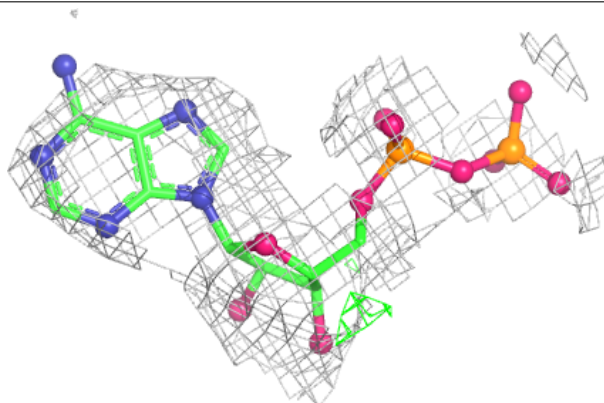


Electron density around ADP C 2000:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around ADP H 2000:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.