



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 10:15 PM UTC

PDB ID : 1Z7N / pdb_00001z7n
Title : ATP Phosphoribosyl transferase (HisZG ATP-PRTase) from *Lactococcus lactis* with bound PRPP substrate
Authors : Champagne, K.S.; Sissler, M.; Larrabee, Y.; Doublié, S.; Francklyn, C.S.
Deposited on : 2005-03-25
Resolution : 3.25 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

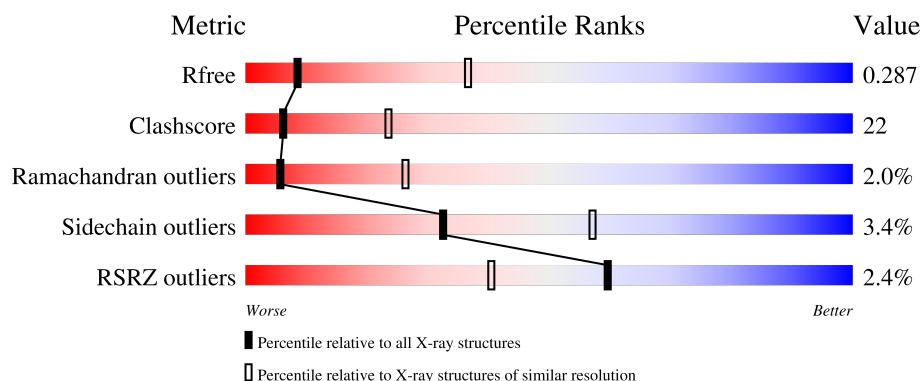
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.25 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	1605 (3.30-3.22)
Clashscore	190562	1660 (3.30-3.22)
Ramachandran outliers	187476	1630 (3.30-3.22)
Sidechain outliers	187428	1629 (3.30-3.22)
RSRZ outliers	180081	1605 (3.30-3.22)

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	344	
1	B	344	
1	C	344	
1	D	344	
2	E	208	

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
2	F	208	4% 54% 41% ..
2	G	208	3% 52% 43% ..
2	H	208	% 52% 40% . .

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
4	PO4	G	1001	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 16750 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 ATP phosphoribosyltransferase regulatory subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	313	Total	C	N	O	S	21	0	0
			2547	1636	415	485	11			
1	B	318	Total	C	N	O	S	12	0	0
			2584	1657	424	492	11			
1	C	304	Total	C	N	O	S	10	0	0
			2472	1591	399	471	11			
1	D	312	Total	C	N	O	S	5	0	0
			2541	1633	416	481	11			

There are 64 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-15	MET	-	cloning artifact	UNP Q02147
A	-14	ARG	-	cloning artifact	UNP Q02147
A	-13	GLY	-	cloning artifact	UNP Q02147
A	-12	SER	-	cloning artifact	UNP Q02147
A	-11	HIS	-	expression tag	UNP Q02147
A	-10	HIS	-	expression tag	UNP Q02147
A	-9	HIS	-	expression tag	UNP Q02147
A	-8	HIS	-	expression tag	UNP Q02147
A	-7	HIS	-	expression tag	UNP Q02147
A	-6	HIS	-	expression tag	UNP Q02147
A	-5	GLY	-	cloning artifact	UNP Q02147
A	-4	SER	-	cloning artifact	UNP Q02147
A	-3	ILE	-	cloning artifact	UNP Q02147
A	-2	GLU	-	cloning artifact	UNP Q02147
A	-1	GLY	-	cloning artifact	UNP Q02147
A	0	ARG	-	cloning artifact	UNP Q02147
B	-15	MET	-	cloning artifact	UNP Q02147
B	-14	ARG	-	cloning artifact	UNP Q02147
B	-13	GLY	-	cloning artifact	UNP Q02147
B	-12	SER	-	cloning artifact	UNP Q02147
B	-11	HIS	-	expression tag	UNP Q02147

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
B	-10	HIS	-	expression tag	UNP Q02147
B	-9	HIS	-	expression tag	UNP Q02147
B	-8	HIS	-	expression tag	UNP Q02147
B	-7	HIS	-	expression tag	UNP Q02147
B	-6	HIS	-	expression tag	UNP Q02147
B	-5	GLY	-	cloning artifact	UNP Q02147
B	-4	SER	-	cloning artifact	UNP Q02147
B	-3	ILE	-	cloning artifact	UNP Q02147
B	-2	GLU	-	cloning artifact	UNP Q02147
B	-1	GLY	-	cloning artifact	UNP Q02147
B	0	ARG	-	cloning artifact	UNP Q02147
C	-15	MET	-	cloning artifact	UNP Q02147
C	-14	ARG	-	cloning artifact	UNP Q02147
C	-13	GLY	-	cloning artifact	UNP Q02147
C	-12	SER	-	cloning artifact	UNP Q02147
C	-11	HIS	-	expression tag	UNP Q02147
C	-10	HIS	-	expression tag	UNP Q02147
C	-9	HIS	-	expression tag	UNP Q02147
C	-8	HIS	-	expression tag	UNP Q02147
C	-7	HIS	-	expression tag	UNP Q02147
C	-6	HIS	-	expression tag	UNP Q02147
C	-5	GLY	-	cloning artifact	UNP Q02147
C	-4	SER	-	cloning artifact	UNP Q02147
C	-3	ILE	-	cloning artifact	UNP Q02147
C	-2	GLU	-	cloning artifact	UNP Q02147
C	-1	GLY	-	cloning artifact	UNP Q02147
C	0	ARG	-	cloning artifact	UNP Q02147
D	-15	MET	-	cloning artifact	UNP Q02147
D	-14	ARG	-	cloning artifact	UNP Q02147
D	-13	GLY	-	cloning artifact	UNP Q02147
D	-12	SER	-	cloning artifact	UNP Q02147
D	-11	HIS	-	expression tag	UNP Q02147
D	-10	HIS	-	expression tag	UNP Q02147
D	-9	HIS	-	expression tag	UNP Q02147
D	-8	HIS	-	expression tag	UNP Q02147
D	-7	HIS	-	expression tag	UNP Q02147
D	-6	HIS	-	expression tag	UNP Q02147
D	-5	GLY	-	cloning artifact	UNP Q02147
D	-4	SER	-	cloning artifact	UNP Q02147
D	-3	ILE	-	cloning artifact	UNP Q02147
D	-2	GLU	-	cloning artifact	UNP Q02147
D	-1	GLY	-	cloning artifact	UNP Q02147

Continued on next page...

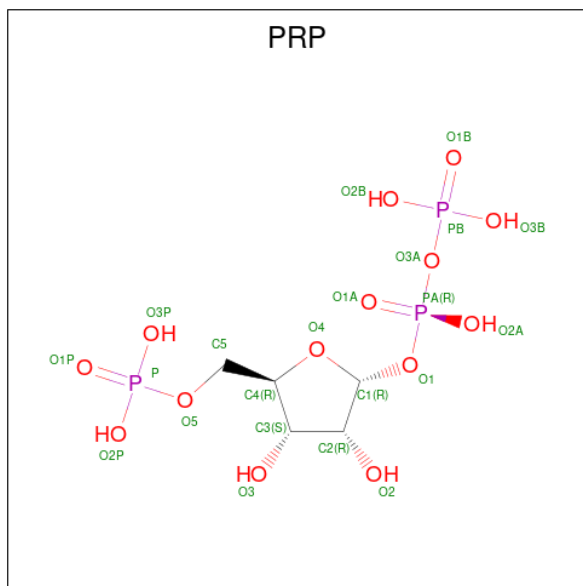
Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
D	0	ARG	-	cloning artifact	UNP Q02147

- Molecule 2 is a protein called ATP phosphoribosyltransferase.

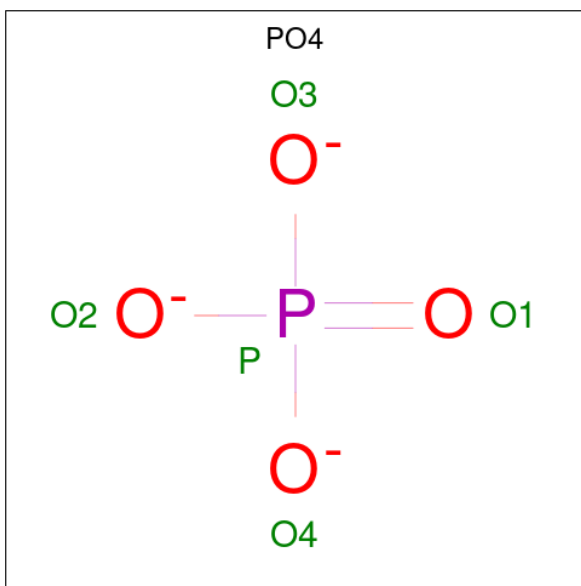
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	203	Total	C	N	O	S	9	0	0
			1630	1048	266	312	4			
2	F	205	Total	C	N	O	S	0	0	0
			1643	1055	269	315	4			
2	G	203	Total	C	N	O	S	4	0	0
			1632	1048	267	313	4			
2	H	200	Total	C	N	O	S	0	0	0
			1603	1030	261	308	4			

- Molecule 3 is 1-O-pyrophosphono-5-O-phosphono-alpha-D-ribofuranose (CCD ID: PRP) (formula: $C_5H_{13}O_{14}P_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	E	1	Total	C	O	P	0	0
			22	5	14	3		
3	F	1	Total	C	O	P	0	0
			22	5	14	3		
3	G	1	Total	C	O	P	0	0
			22	5	14	3		
3	H	1	Total	C	O	P	0	0
			22	5	14	3		

- Molecule 4 is PHOSPHATE ION (CCD ID: PO4) (formula: O_4P).

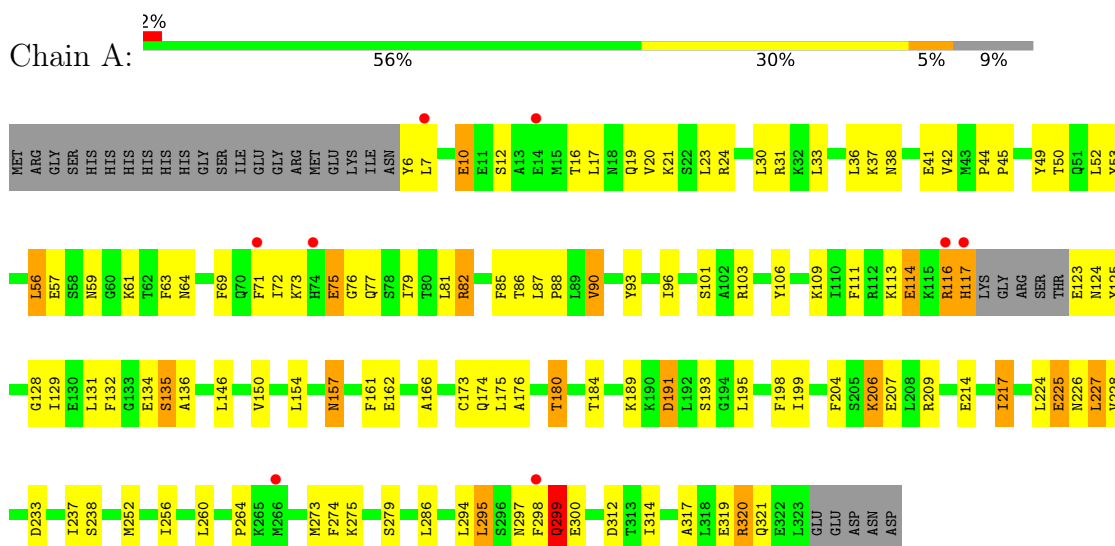


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	F	1	Total	O	P	0	0
			5	4	1		
4	G	1	Total	O	P	0	0
			5	4	1		

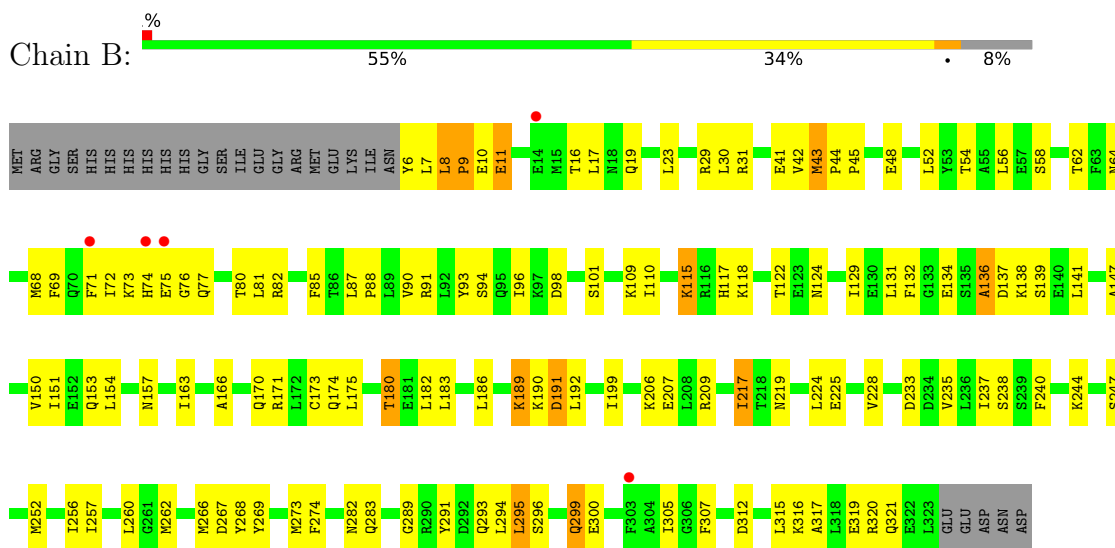
3 Residue-property plots [i](#)

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: ATP phosphoribosyltransferase regulatory subunit



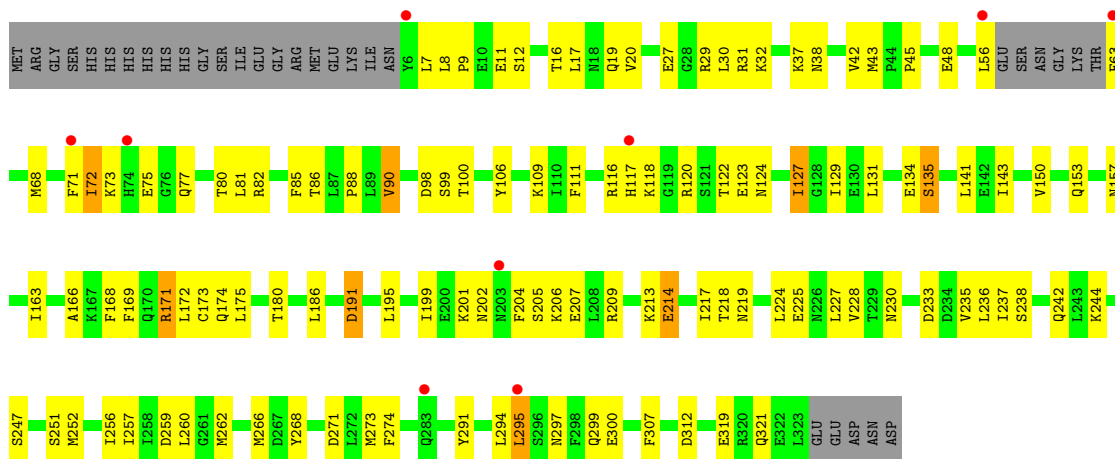
- Molecule 1: ATP phosphoribosyltransferase regulatory subunit



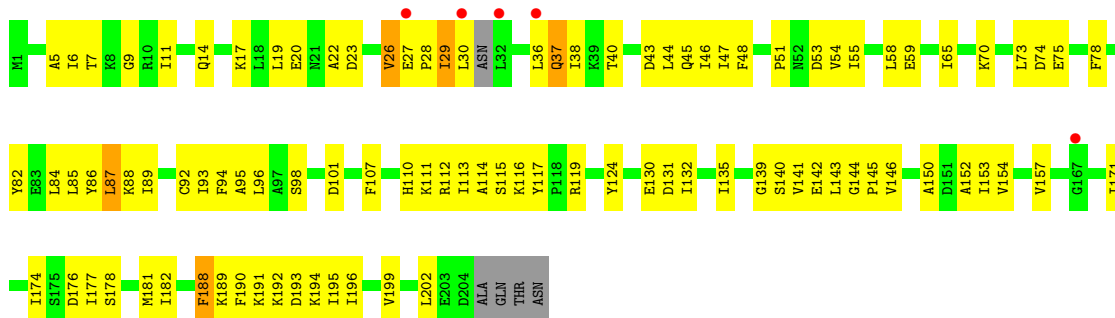
- Molecule 1: ATP phosphoribosyltransferase regulatory subunit



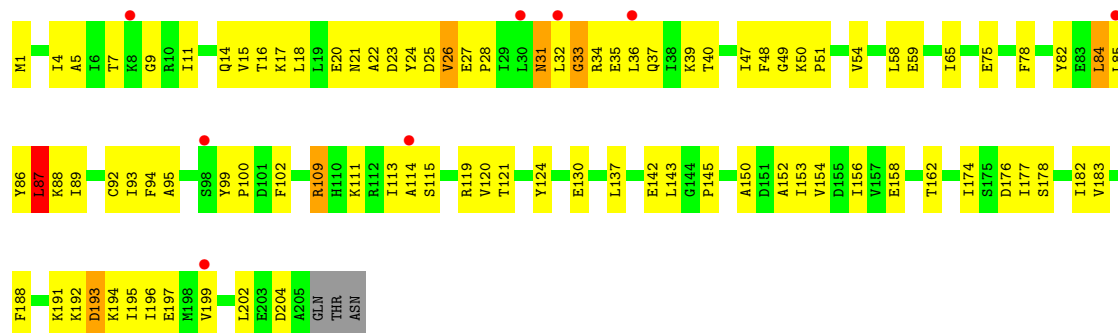
- Molecule 1: ATP phosphoribosyltransferase regulatory subunit



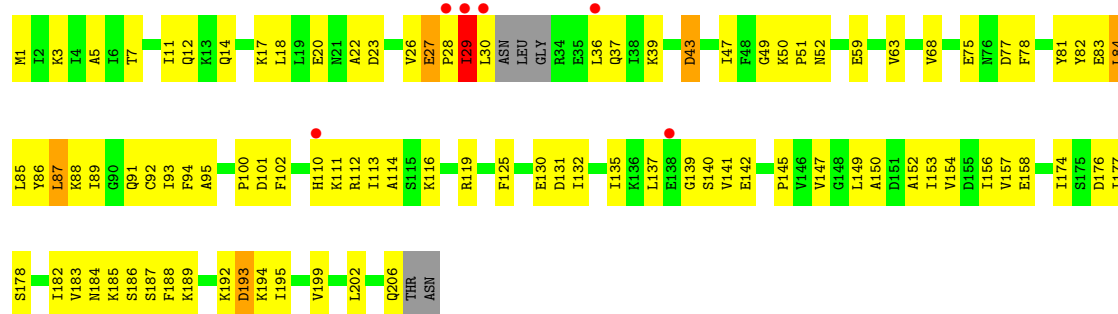
- Molecule 2: ATP phosphoribosyltransferase



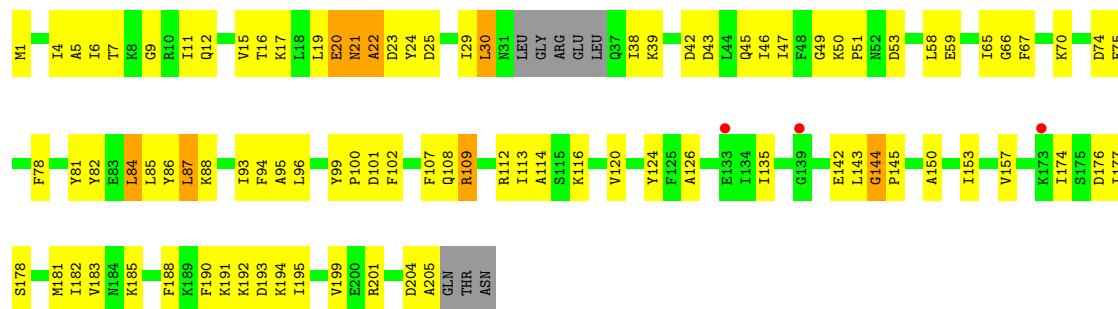
- Molecule 2: ATP phosphoribosyltransferase



• Molecule 2: ATP phosphoribosyltransferase



• Molecule 2: ATP phosphoribosyltransferase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	151.68Å 222.94Å 86.38Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.00 – 3.25 29.00 – 3.26	Depositor EDS
% Data completeness (in resolution range)	85.7 (29.00-3.25) 85.9 (29.00-3.26)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.86 (at 3.24Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.247 , 0.301 0.241 , 0.287	Depositor DCC
R_{free} test set	1981 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å ²)	61.3	Xtriage
Anisotropy	0.073	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 43.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	16750	wwPDB-VP
Average B, all atoms (Å ²)	66.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.03% 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: PRP, PO4

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	0.68	1/2589 (0.0%)	1.01	15/3475 (0.4%)
1	B	0.59	0/2627	0.94	3/3526 (0.1%)
1	C	0.59	0/2511	0.97	3/3370 (0.1%)
1	D	0.51	0/2583	0.93	12/3466 (0.3%)
2	E	0.81	4/1652 (0.2%)	1.01	13/2220 (0.6%)
2	F	0.51	0/1666	0.93	4/2241 (0.2%)
2	G	0.54	0/1654	0.97	3/2223 (0.1%)
2	H	0.58	1/1625 (0.1%)	0.99	9/2185 (0.4%)
All	All	0.61	6/16907 (0.0%)	0.97	62/22706 (0.3%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
2	E	0	2
All	All	0	3

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	29	ILE	C-N	-20.89	1.04	1.33
2	E	26	VAL	CA-C	8.96	1.63	1.52
1	A	114	GLU	C-N	-8.85	1.20	1.33
2	H	21	ASN	N-CA	6.28	1.54	1.46
2	E	26	VAL	C-N	5.29	1.41	1.33
2	E	188	PHE	CA-C	5.10	1.59	1.52

All (62) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	299	GLN	OE1-CD-NE2	-9.80	112.80	122.60
1	A	299	GLN	OE1-CD-NE2	-9.77	112.83	122.60
1	A	64	ASN	N-CA-C	9.52	124.24	108.73
1	D	43	MET	CA-C-N	9.50	126.59	119.66
1	D	43	MET	C-N-CA	9.50	126.59	119.66
1	A	114	GLU	CA-C-N	-9.39	103.27	121.58
1	A	114	GLU	C-N-CA	-9.39	103.27	121.58
2	E	37	GLN	OE1-CD-NE2	-8.53	114.07	122.60
2	H	22	ALA	N-CA-C	-7.69	96.36	108.90
2	E	117	TYR	CA-C-N	7.67	126.94	118.97
2	E	117	TYR	C-N-CA	7.67	126.94	118.97
1	A	116	ARG	CA-C-N	7.56	135.30	121.70
1	A	116	ARG	C-N-CA	7.56	135.30	121.70
2	G	29	ILE	N-CA-C	-7.03	94.72	109.34
2	E	29	ILE	CA-C-N	-7.01	109.09	121.70
2	E	29	ILE	C-N-CA	-7.01	109.09	121.70
2	E	26	VAL	CA-C-N	-6.91	109.91	122.35
2	E	26	VAL	C-N-CA	-6.91	109.91	122.35
1	A	206	LYS	N-CA-C	-6.86	103.74	111.14
1	A	10	GLU	N-CA-C	-6.77	105.03	113.28
2	F	32	LEU	N-CA-C	6.77	118.34	110.97
1	D	118	LYS	N-CA-C	-6.72	104.78	114.12
1	B	299	GLN	CG-CD-NE2	6.70	126.45	116.40
1	D	201	LYS	N-CA-C	6.66	119.39	111.33
1	A	299	GLN	CG-CD-NE2	6.66	126.39	116.40
1	C	43	MET	CA-C-N	6.53	124.42	119.66
1	C	43	MET	C-N-CA	6.53	124.42	119.66
1	A	214	GLU	N-CA-C	6.47	119.30	111.40
2	E	37	GLN	CG-CD-NE2	6.38	125.97	116.40
1	C	217	ILE	N-CA-C	6.35	122.54	109.34
1	A	217	ILE	N-CA-C	6.29	122.43	109.34
2	F	22	ALA	N-CA-C	-6.25	105.29	113.17
1	B	217	ILE	N-CA-C	6.21	122.26	109.34
2	H	30	LEU	N-CA-C	6.20	119.24	110.50
2	G	116	LYS	N-CA-C	-6.08	105.86	113.28
2	E	191	LYS	N-CA-C	-6.03	106.17	112.93
1	D	217	ILE	N-CA-C	6.01	121.84	109.34
2	F	33	GLY	N-CA-C	5.94	127.27	113.18
2	E	188	PHE	CB-CA-C	5.76	120.65	110.85
2	E	189	LYS	CA-C-N	-5.76	112.53	122.79
2	E	189	LYS	C-N-CA	-5.76	112.53	122.79
1	A	56	LEU	N-CA-C	5.60	117.46	111.36
1	D	171	ARG	N-CA-C	-5.48	105.20	111.07

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	90	VAL	N-CA-C	-5.39	105.25	110.42
2	F	54	VAL	N-CA-C	-5.37	105.48	110.53
2	H	144	GLY	CA-C-N	5.35	124.86	119.19
2	H	144	GLY	C-N-CA	5.35	124.86	119.19
1	A	116	ARG	O-C-N	-5.34	114.89	122.15
2	H	20	GLU	CA-C-N	5.34	129.23	120.63
2	H	20	GLU	C-N-CA	5.34	129.23	120.63
1	D	72	ILE	N-CA-C	5.29	116.59	108.71
1	A	157	ASN	N-CA-C	5.28	116.89	111.03
1	D	127	ILE	N-CA-C	-5.26	101.07	108.27
2	H	21	ASN	CA-C-N	5.22	130.35	122.99
2	H	21	ASN	C-N-CA	5.22	130.35	122.99
2	E	54	VAL	N-CA-C	-5.21	104.89	110.36
1	A	90	VAL	N-CA-C	-5.20	105.46	110.72
1	D	214	GLU	N-CA-C	5.17	117.71	111.40
2	G	195	ILE	N-CA-C	-5.09	105.37	110.30
1	D	218	THR	N-CA-C	5.06	116.76	108.76
1	D	242	GLN	N-CA-C	-5.05	105.05	111.11
2	H	116	LYS	N-CA-C	-5.03	107.14	113.28

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	114	GLU	Mainchain
2	E	26	VAL	Mainchain
2	E	29	ILE	Mainchain

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	2547	0	2574	117	0
1	B	2584	0	2616	121	0
1	C	2472	0	2504	131	0
1	D	2541	0	2575	106	0
2	E	1630	0	1679	92	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	F	1643	0	1692	85	0
2	G	1632	0	1679	80	0
2	H	1603	0	1647	85	0
3	E	22	0	8	0	0
3	F	22	0	8	1	0
3	G	22	0	8	1	0
3	H	22	0	8	0	0
4	F	5	0	0	0	0
4	G	5	0	0	2	0
All	All	16750	0	16998	754	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 22.

All (754) 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:A:116:ARG:O	1:A:117:HIS:CD2	1.76	1.37
1:A:116:ARG:C	1:A:117:HIS:CD2	2.08	1.30
2:E:27:GLU:O	2:E:30:LEU:HG	1.24	1.28
1:A:116:ARG:O	1:A:117:HIS:CG	1.91	1.22
2:H:19:LEU:O	2:H:22:ALA:O	1.63	1.15
2:E:36:LEU:HD13	2:F:143:LEU:HG	1.29	1.13
1:D:73:LYS:HE3	1:D:75:GLU:HB3	1.39	1.04
1:C:173:CYS:HB3	1:C:180:THR:HG22	1.40	1.00
1:B:73:LYS:HE3	1:B:75:GLU:HB3	1.44	0.99
2:G:86:TYR:HE1	2:G:178:SER:HG	1.09	0.97
2:G:111:LYS:HE2	2:G:130:GLU:OE2	1.65	0.97
1:A:299:GLN:HG2	1:A:300:GLU:H	1.26	0.96
1:B:115:LYS:HG3	2:G:77:ASP:OD2	1.66	0.95
1:C:73:LYS:HE3	1:C:75:GLU:HB3	1.47	0.95
2:E:27:GLU:O	2:E:30:LEU:CG	2.14	0.95
1:C:299:GLN:HG2	1:C:300:GLU:H	1.32	0.94
2:E:86:TYR:HE1	2:E:178:SER:HG	1.16	0.93
1:A:116:ARG:HB3	1:A:117:HIS:CD2	2.04	0.93
2:H:87:LEU:HD23	2:H:87:LEU:H	1.32	0.93
1:A:73:LYS:HE3	1:A:75:GLU:HB3	1.50	0.92
2:G:59:GLU:HG3	2:G:78:PHE:CE1	2.06	0.91
1:D:173:CYS:HB3	1:D:180:THR:HG22	1.52	0.89
1:A:73:LYS:HE2	1:A:77:GLN:HB2	1.55	0.88
1:C:12:SER:HB2	1:D:45:PRO:HB2	1.55	0.88

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:86:TYR:HE1	2:H:178:SER:HG	1.17	0.87
1:A:116:ARG:HB3	1:A:117:HIS:NE2	1.89	0.87
2:E:36:LEU:CD1	2:F:143:LEU:HG	2.03	0.87
1:A:73:LYS:HG2	1:A:77:GLN:H	1.40	0.86
2:F:93:ILE:HG22	2:F:176:ASP:HA	1.57	0.85
2:E:36:LEU:HD13	2:F:143:LEU:CG	2.07	0.85
2:E:87:LEU:HD23	2:E:87:LEU:H	1.41	0.84
1:A:109:LYS:HD3	1:A:123:GLU:HG2	1.59	0.83
1:D:299:GLN:HG2	1:D:300:GLU:H	1.42	0.82
2:E:59:GLU:HG3	2:E:78:PHE:CE1	2.15	0.82
2:G:193:ASP:OD2	2:G:194:LYS:HG3	1.80	0.82
2:G:27:GLU:H	2:G:28:PRO:CD	1.92	0.81
1:B:199:ILE:HG21	1:B:209:ARG:HG3	1.62	0.81
2:F:86:TYR:HE1	2:F:178:SER:HG	1.26	0.80
2:F:195:ILE:O	2:F:199:VAL:HG23	1.82	0.80
1:A:90:VAL:HG13	1:A:295:LEU:HD22	1.64	0.80
1:C:73:LYS:HE2	1:C:77:GLN:HB2	1.62	0.80
2:H:193:ASP:OD2	2:H:194:LYS:HG3	1.82	0.80
2:E:111:LYS:HE2	2:E:130:GLU:OE2	1.82	0.80
2:F:59:GLU:HG3	2:F:78:PHE:CE1	2.17	0.80
2:F:111:LYS:HE2	2:F:130:GLU:OE2	1.82	0.80
1:C:199:ILE:HD13	1:C:209:ARG:HG3	1.64	0.79
1:D:202:ASN:HB2	1:D:204:PHE:HE1	1.47	0.78
1:C:45:PRO:HG3	1:D:8:LEU:HD11	1.64	0.78
1:A:299:GLN:HG2	1:A:300:GLU:N	1.99	0.78
1:C:81:LEU:HD11	1:D:81:LEU:HD21	1.64	0.78
1:C:193:SER:OG	2:H:86:TYR:HB3	1.85	0.77
1:C:199:ILE:HG21	1:C:209:ARG:HG3	1.67	0.77
2:E:113:ILE:HG23	2:E:153:ILE:HA	1.66	0.77
2:H:193:ASP:CG	2:H:194:LYS:H	1.93	0.76
1:A:116:ARG:C	1:A:117:HIS:HD2	1.90	0.76
1:B:16:THR:HG23	1:B:19:GLN:H	1.51	0.75
1:C:131:LEU:HD21	1:C:134:GLU:HB3	1.68	0.75
1:A:173:CYS:HB3	1:A:180:THR:HG22	1.69	0.74
1:A:73:LYS:HE2	1:A:77:GLN:CB	2.18	0.74
2:F:193:ASP:OD2	2:F:194:LYS:HG3	1.87	0.74
2:G:86:TYR:HE1	2:G:178:SER:OG	1.70	0.74
1:B:48:GLU:HG2	1:B:52:LEU:HD22	1.70	0.74
1:D:56:LEU:HD21	1:D:63:PHE:CE1	2.23	0.73
2:E:195:ILE:O	2:E:199:VAL:HG23	1.89	0.73
1:A:199:ILE:HG21	1:A:209:ARG:HG3	1.69	0.73

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:18:LEU:HD23	2:F:202:LEU:HD22	1.70	0.73
1:B:64:ASN:O	1:B:68:MET:HE3	1.88	0.73
2:H:86:TYR:CE2	2:H:88:LYS:HG2	2.24	0.72
1:B:73:LYS:HE2	1:B:77:GLN:HB2	1.71	0.72
2:E:86:TYR:HE1	2:E:178:SER:OG	1.71	0.72
2:F:16:THR:CG2	2:F:28:PRO:HG3	2.20	0.72
2:F:16:THR:HG22	2:F:28:PRO:HG3	1.72	0.72
1:A:299:GLN:CG	1:A:300:GLU:H	2.02	0.71
2:G:87:LEU:HD23	2:G:87:LEU:H	1.55	0.71
1:A:116:ARG:CB	1:A:117:HIS:CD2	2.73	0.71
1:D:7:LEU:HD23	1:D:7:LEU:H	1.54	0.70
2:F:109:ARG:HG2	2:F:109:ARG:HH11	1.57	0.70
2:F:87:LEU:HD23	2:F:87:LEU:H	1.57	0.70
1:C:49:TYR:CE2	1:D:9:PRO:HB3	2.26	0.70
1:B:189:LYS:HE2	2:G:81:TYR:CE1	2.28	0.69
2:E:95:ALA:HA	2:E:174:ILE:HG12	1.74	0.69
2:E:193:ASP:OD2	2:E:194:LYS:HG3	1.92	0.69
2:G:189:LYS:NZ	4:G:1001:PO4:O2	2.22	0.69
1:B:30:LEU:HD21	1:B:150:VAL:HG13	1.75	0.68
1:A:73:LYS:HE2	1:A:77:GLN:CG	2.24	0.68
1:A:116:ARG:O	1:A:117:HIS:NE2	2.25	0.68
1:A:131:LEU:HD21	1:A:134:GLU:HB3	1.76	0.68
1:B:96:ILE:HG22	1:B:98:ASP:H	1.59	0.68
2:F:114:ALA:HB1	2:F:137:LEU:HD12	1.74	0.68
2:H:195:ILE:O	2:H:199:VAL:HG23	1.94	0.68
2:H:59:GLU:HG3	2:H:78:PHE:CE1	2.30	0.67
1:C:7:LEU:HG	2:G:131:ASP:HB2	1.77	0.67
1:D:8:LEU:HG	1:D:9:PRO:HD2	1.76	0.67
1:C:90:VAL:HG13	1:C:295:LEU:HD22	1.76	0.67
2:F:58:LEU:C	2:F:58:LEU:HD23	2.20	0.67
1:B:206:LYS:HD3	1:B:206:LYS:C	2.20	0.67
2:E:36:LEU:HB3	2:F:143:LEU:HD21	1.76	0.67
1:D:16:THR:HG23	1:D:19:GLN:H	1.60	0.66
1:D:73:LYS:HE2	1:D:77:GLN:HB2	1.77	0.66
1:C:173:CYS:CB	1:C:180:THR:HG22	2.22	0.66
2:E:36:LEU:CB	2:F:143:LEU:HD21	2.24	0.66
1:A:131:LEU:C	1:A:131:LEU:HD23	2.20	0.66
2:H:38:ILE:HD11	2:H:46:ILE:HB	1.77	0.66
1:B:293:GLN:O	1:B:296:SER:HB2	1.96	0.66
1:B:147:ALA:O	1:B:151:ILE:HG13	1.95	0.66
1:D:131:LEU:HD21	1:D:134:GLU:HB3	1.78	0.66

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:73:LYS:HE2	1:C:77:GLN:CB	2.25	0.66
2:E:30:LEU:HD22	2:E:38:ILE:HG12	1.78	0.66
1:B:30:LEU:CD2	1:B:150:VAL:HG13	2.26	0.66
1:C:45:PRO:HB2	1:D:12:SER:HB2	1.78	0.66
1:A:224:LEU:HD22	1:A:260:LEU:HD12	1.76	0.66
1:A:20:VAL:HG22	1:B:43:MET:HG3	1.78	0.65
2:F:193:ASP:CG	2:F:194:LYS:H	2.03	0.65
2:F:25:ASP:O	2:F:27:GLU:HG2	1.96	0.65
2:G:113:ILE:HG21	2:G:154:VAL:HG23	1.79	0.65
2:H:96:LEU:HD12	2:H:153:ILE:O	1.96	0.65
1:A:116:ARG:C	1:A:117:HIS:CG	2.49	0.65
1:D:206:LYS:HD3	1:D:206:LYS:C	2.21	0.65
1:A:320:ARG:O	1:A:320:ARG:HG2	1.95	0.65
1:C:73:LYS:HE2	1:C:77:GLN:CG	2.26	0.65
1:C:71:PHE:HE2	1:D:71:PHE:HE2	1.44	0.64
1:C:81:LEU:HB3	1:C:111:PHE:CD1	2.32	0.64
1:A:129:ILE:CD1	1:A:146:LEU:HD13	2.27	0.64
2:G:18:LEU:HD23	2:G:202:LEU:HD22	1.79	0.64
1:B:117:HIS:O	1:B:118:LYS:CG	2.46	0.64
2:G:93:ILE:HG13	2:G:157:VAL:O	1.98	0.64
2:H:30:LEU:HD11	2:H:38:ILE:HG23	1.78	0.64
2:H:38:ILE:HD12	2:H:38:ILE:C	2.23	0.64
2:F:84:LEU:HD12	2:F:183:VAL:HG21	1.78	0.64
2:E:28:PRO:C	2:E:30:LEU:N	2.54	0.64
1:C:71:PHE:HD2	1:D:71:PHE:HD2	1.45	0.64
1:D:202:ASN:HB2	1:D:204:PHE:CE1	2.31	0.63
2:F:86:TYR:HE1	2:F:178:SER:OG	1.81	0.63
2:G:189:LYS:HZ1	4:G:1001:PO4:P	2.21	0.63
2:E:38:ILE:HD12	2:E:38:ILE:C	2.24	0.63
2:G:95:ALA:HA	2:G:174:ILE:HG12	1.79	0.63
1:B:163:ILE:HG22	1:B:274:PHE:HB3	1.79	0.63
2:E:38:ILE:HD11	2:E:46:ILE:HB	1.79	0.63
1:D:291:TYR:CD1	1:D:294:LEU:HD12	2.34	0.63
2:E:111:LYS:HE2	2:E:130:GLU:CD	2.25	0.62
1:D:109:LYS:HE2	1:D:123:GLU:HB3	1.80	0.62
1:C:45:PRO:O	1:C:82:ARG:HG3	2.00	0.62
1:A:71:PHE:CE1	1:A:79:ILE:HD11	2.34	0.62
1:B:72:ILE:HG13	1:D:71:PHE:HA	1.82	0.62
1:C:206:LYS:HD3	1:C:206:LYS:C	2.24	0.61
1:C:42:VAL:HG11	1:C:89:LEU:HD22	1.81	0.61
1:D:199:ILE:HD13	1:D:209:ARG:HG3	1.80	0.61

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:29:ARG:HD2	1:B:153:GLN:OE1	2.01	0.61
2:F:37:GLN:HE21	2:F:47:ILE:HD11	1.65	0.61
2:H:94:PHE:CD1	2:H:177:ILE:HD11	2.36	0.61
1:D:202:ASN:O	1:D:204:PHE:HD1	1.84	0.61
2:H:65:ILE:HD13	2:H:181:MET:HE3	1.82	0.61
2:H:58:LEU:HD13	2:H:182:ILE:HG13	1.82	0.60
2:H:84:LEU:HD12	2:H:183:VAL:HG21	1.82	0.60
1:A:16:THR:HG23	1:A:19:GLN:H	1.66	0.60
2:F:109:ARG:HG2	2:F:109:ARG:NH1	2.16	0.60
2:G:27:GLU:N	2:G:28:PRO:CD	2.61	0.60
1:C:224:LEU:HD22	1:C:260:LEU:HD12	1.82	0.60
2:H:193:ASP:CG	2:H:194:LYS:N	2.60	0.60
1:C:129:ILE:HD11	1:C:307:PHE:CE2	2.36	0.60
1:A:71:PHE:CE2	1:B:71:PHE:CE2	2.90	0.60
1:B:173:CYS:HB3	1:B:180:THR:HG22	1.84	0.60
2:G:142:GLU:O	2:G:145:PRO:HD2	2.02	0.60
1:C:81:LEU:CD1	1:D:81:LEU:HD21	2.31	0.60
2:F:145:PRO:HA	2:F:150:ALA:O	2.01	0.60
2:G:113:ILE:HG23	2:G:153:ILE:HA	1.84	0.60
1:B:11:GLU:C	1:B:122:THR:HG21	2.26	0.60
2:G:86:TYR:HE1	2:G:178:SER:CB	2.14	0.60
2:H:1:MET:HE1	2:H:39:LYS:HE3	1.84	0.60
1:A:154:LEU:HD13	1:A:314:ILE:HD13	1.84	0.60
2:F:188:PHE:CE2	2:F:192:LYS:HE2	2.37	0.60
2:G:85:LEU:HG	2:G:86:TYR:N	2.16	0.60
1:A:71:PHE:CD1	1:A:79:ILE:HD11	2.37	0.59
1:B:131:LEU:HD21	1:B:134:GLU:HB3	1.84	0.59
2:F:191:LYS:O	2:F:195:ILE:HG12	2.01	0.59
2:H:85:LEU:HG	2:H:86:TYR:N	2.17	0.59
2:G:84:LEU:HD12	2:G:183:VAL:HG21	1.83	0.59
2:H:188:PHE:O	2:H:192:LYS:HB2	2.01	0.59
1:A:73:LYS:CE	1:A:77:GLN:HB2	2.28	0.59
1:B:6:TYR:HB2	1:B:7:LEU:HD12	1.84	0.59
2:G:185:LYS:O	2:G:189:LYS:HG2	2.02	0.59
1:B:73:LYS:HE2	1:B:77:GLN:CB	2.32	0.59
2:E:143:LEU:HG	2:F:36:LEU:HD13	1.84	0.59
1:B:150:VAL:O	1:B:154:LEU:HG	2.03	0.59
2:H:188:PHE:CE2	2:H:192:LYS:HE2	2.37	0.59
1:C:71:PHE:CE2	1:D:71:PHE:HE2	2.21	0.59
2:F:87:LEU:O	2:F:89:ILE:HG12	2.02	0.59
1:C:204:PHE:O	1:C:209:ARG:NH2	2.36	0.58

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:124:ASN:HA	1:D:312:ASP:OD2	2.02	0.58
1:D:90:VAL:HG13	1:D:295:LEU:HD22	1.83	0.58
2:F:75:GLU:OE1	2:F:120:VAL:HG23	2.02	0.58
2:E:94:PHE:CD1	2:E:177:ILE:HD11	2.38	0.58
2:G:125:PHE:CD1	2:G:132:ILE:HD13	2.38	0.58
2:H:93:ILE:O	2:H:93:ILE:HD12	2.03	0.58
2:E:7:THR:OG1	2:E:11:ILE:HB	2.03	0.58
1:A:49:TYR:CD2	1:B:9:PRO:HG3	2.38	0.58
2:H:93:ILE:HG22	2:H:176:ASP:HA	1.86	0.58
2:G:113:ILE:HD12	2:G:113:ILE:N	2.19	0.58
2:H:17:LYS:HE3	2:H:21:ASN:HD21	1.68	0.58
1:C:94:SER:C	1:C:96:ILE:H	2.12	0.57
2:E:193:ASP:CG	2:E:194:LYS:H	2.11	0.57
2:G:29:ILE:HB	2:G:30:LEU:HD12	1.86	0.57
2:F:7:THR:HA	2:F:49:GLY:O	2.03	0.57
2:G:36:LEU:HD23	2:H:143:LEU:HG	1.85	0.57
1:A:319:GLU:C	1:A:321:GLN:H	2.10	0.57
1:C:49:TYR:CZ	1:D:9:PRO:HB3	2.39	0.57
2:E:143:LEU:O	2:E:144:GLY:C	2.46	0.57
2:H:112:ARG:C	2:H:113:ILE:HD12	2.29	0.57
1:C:275:LYS:HD3	1:C:277:TYR:CE1	2.39	0.57
1:D:199:ILE:HG21	1:D:209:ARG:HG3	1.86	0.57
2:E:188:PHE:O	2:E:192:LYS:HB2	2.05	0.57
1:C:37:LYS:O	1:C:38:ASN:HB2	2.05	0.57
1:C:219:ASN:HB3	1:C:258:ILE:O	2.05	0.57
1:B:237:ILE:HG13	1:B:238:SER:N	2.19	0.57
1:B:73:LYS:HE2	1:B:77:GLN:CG	2.35	0.57
1:B:129:ILE:HD11	1:B:307:PHE:CE2	2.40	0.57
1:C:12:SER:HB2	1:D:45:PRO:CB	2.31	0.57
1:C:71:PHE:HD2	1:D:71:PHE:CD2	2.22	0.57
1:C:283:GLN:HB3	1:C:284:PRO:HD2	1.86	0.57
2:H:86:TYR:HE1	2:H:178:SER:OG	1.85	0.57
1:B:175:LEU:HD11	1:B:233:ASP:OD1	2.05	0.57
2:E:14:GLN:HE21	2:E:89:ILE:HD12	1.70	0.57
2:H:7:THR:OG1	2:H:11:ILE:HB	2.04	0.57
2:H:38:ILE:HD12	2:H:38:ILE:O	2.05	0.57
1:C:71:PHE:HE2	1:D:71:PHE:CE2	2.23	0.57
1:D:56:LEU:HD11	1:D:63:PHE:CG	2.39	0.56
1:C:124:ASN:HA	1:C:312:ASP:OD2	2.06	0.56
1:C:71:PHE:CE2	1:D:71:PHE:CE2	2.93	0.56
1:C:71:PHE:CD2	1:D:71:PHE:HD2	2.22	0.56

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:116:ARG:HG3	1:D:117:HIS:CD2	2.40	0.56
2:F:95:ALA:HA	2:F:174:ILE:HG12	1.88	0.56
2:F:113:ILE:HG23	2:F:153:ILE:HA	1.88	0.56
2:G:87:LEU:HA	2:G:206:GLN:NE2	2.20	0.56
1:D:27:GLU:HG2	1:D:127:ILE:HD11	1.88	0.56
2:H:86:TYR:CZ	2:H:88:LYS:HG2	2.40	0.56
1:B:58:SER:HB3	1:B:266:MET:HA	1.85	0.56
2:F:7:THR:OG1	2:F:11:ILE:HB	2.06	0.56
1:A:175:LEU:HD11	1:A:233:ASP:OD1	2.06	0.56
1:C:71:PHE:CD2	1:D:71:PHE:CD2	2.94	0.56
2:H:145:PRO:HA	2:H:150:ALA:O	2.06	0.56
1:A:217:ILE:HB	2:E:196:ILE:HD11	1.88	0.56
1:C:192:LEU:HD12	2:H:84:LEU:HA	1.88	0.56
1:D:29:ARG:HD2	1:D:153:GLN:OE1	2.07	0.56
2:H:86:TYR:HE1	2:H:178:SER:CB	2.18	0.56
1:A:166:ALA:N	1:A:273:MET:HE2	2.20	0.55
1:A:20:VAL:CG2	1:B:43:MET:HG3	2.36	0.55
2:H:58:LEU:CD1	2:H:182:ILE:HG13	2.36	0.55
1:A:319:GLU:O	1:A:321:GLN:N	2.40	0.55
1:B:224:LEU:O	1:B:228:VAL:HG23	2.07	0.55
2:F:15:VAL:O	2:F:16:THR:C	2.50	0.55
2:G:1:MET:HE1	2:G:39:LYS:HE3	1.87	0.55
1:C:71:PHE:CE1	1:C:79:ILE:HD11	2.41	0.55
1:C:166:ALA:N	1:C:273:MET:HE2	2.21	0.55
2:F:31:ASN:OD1	2:F:35:GLU:HG2	2.07	0.55
1:B:117:HIS:O	1:B:118:LYS:HG3	2.06	0.55
1:D:224:LEU:HD22	1:D:260:LEU:HD12	1.88	0.55
1:A:233:ASP:O	1:A:237:ILE:HG23	2.07	0.55
1:B:16:THR:HG22	1:B:19:GLN:HG3	1.89	0.55
1:C:299:GLN:HG2	1:C:300:GLU:N	2.13	0.55
2:F:86:TYR:HE1	2:F:178:SER:CB	2.19	0.55
1:A:206:LYS:HD3	1:A:206:LYS:C	2.32	0.54
1:B:85:PHE:O	1:B:88:PRO:HG2	2.07	0.54
1:C:16:THR:HG23	1:C:19:GLN:H	1.71	0.54
1:D:244:LYS:O	1:D:247:SER:HB3	2.07	0.54
2:F:58:LEU:CD1	2:F:182:ILE:HG13	2.37	0.54
2:E:86:TYR:HE1	2:E:178:SER:CB	2.20	0.54
1:C:69:PHE:O	1:C:80:THR:HG23	2.07	0.54
1:B:23:LEU:HB2	1:B:315:LEU:HD11	1.89	0.54
1:C:131:LEU:C	1:C:131:LEU:HD23	2.32	0.54
2:H:124:TYR:CG	2:H:174:ILE:HD12	2.43	0.54

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:45:PRO:O	1:A:82:ARG:HG3	2.07	0.54
1:A:86:THR:O	1:A:90:VAL:HG23	2.07	0.54
2:H:86:TYR:O	2:H:86:TYR:CG	2.58	0.54
1:A:12:SER:HB2	1:B:45:PRO:HB2	1.88	0.54
1:D:37:LYS:O	1:D:38:ASN:HB2	2.08	0.54
2:G:7:THR:OG1	2:G:11:ILE:HB	2.08	0.54
2:H:113:ILE:HD12	2:H:113:ILE:N	2.22	0.54
1:A:129:ILE:HD13	1:A:146:LEU:HD13	1.88	0.54
1:B:16:THR:CG2	1:B:19:GLN:HG3	2.38	0.54
1:B:207:GLU:H	1:B:207:GLU:CD	2.15	0.54
1:D:169:PHE:O	1:D:172:LEU:HB3	2.08	0.54
1:D:168:PHE:CE2	1:D:260:LEU:HB3	2.43	0.53
1:D:299:GLN:HG2	1:D:300:GLU:N	2.17	0.53
1:D:131:LEU:C	1:D:131:LEU:HD23	2.33	0.53
2:F:124:TYR:CG	2:F:174:ILE:HD12	2.44	0.53
2:G:29:ILE:C	2:G:30:LEU:HD12	2.34	0.53
1:B:16:THR:O	1:B:17:LEU:C	2.51	0.53
2:E:116:LYS:HG2	2:E:116:LYS:O	2.08	0.53
2:E:58:LEU:HD13	2:E:182:ILE:HG13	1.91	0.53
1:B:16:THR:HG22	1:B:19:GLN:CG	2.39	0.53
2:G:147:VAL:HG23	2:G:149:LEU:H	1.73	0.53
2:G:7:THR:HG22	2:G:51:PRO:HA	1.90	0.53
2:G:37:GLN:HB3	2:G:47:ILE:HG12	1.89	0.53
1:C:275:LYS:HA	1:C:286:LEU:O	2.09	0.53
2:E:143:LEU:HD21	2:F:36:LEU:HB2	1.91	0.53
2:G:188:PHE:O	2:G:192:LYS:HB2	2.09	0.53
1:A:71:PHE:HE2	1:B:71:PHE:CE2	2.25	0.53
1:C:79:ILE:HD13	1:D:11:GLU:HG3	1.90	0.53
1:D:68:MET:HB2	1:D:80:THR:HG21	1.92	0.53
1:B:299:GLN:HG2	1:B:300:GLU:H	1.73	0.52
2:E:65:ILE:HG21	2:E:181:MET:HE3	1.91	0.52
1:A:116:ARG:O	1:A:117:HIS:CE1	2.62	0.52
2:E:114:ALA:HA	2:E:135:ILE:O	2.10	0.52
2:F:75:GLU:CD	2:F:119:ARG:HB2	2.34	0.52
1:D:116:ARG:HG3	1:D:117:HIS:HD2	1.74	0.52
1:B:90:VAL:HG13	1:B:295:LEU:HD22	1.92	0.52
1:C:7:LEU:C	1:C:8:LEU:HD12	2.35	0.52
1:C:81:LEU:HD21	1:D:81:LEU:HD11	1.91	0.52
1:A:124:ASN:HA	1:A:312:ASP:OD2	2.09	0.52
1:B:71:PHE:HA	1:D:72:ILE:HG13	1.92	0.52
1:C:157:ASN:OD1	1:C:321:GLN:NE2	2.43	0.52

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:233:ASP:O	1:C:237:ILE:HG23	2.10	0.52
1:D:85:PHE:O	1:D:88:PRO:HG2	2.10	0.52
1:B:224:LEU:HD22	1:B:260:LEU:HD12	1.91	0.52
1:C:16:THR:O	1:C:19:GLN:N	2.41	0.52
2:H:5:ALA:HA	2:H:47:ILE:O	2.10	0.52
2:H:17:LYS:O	2:H:20:GLU:HB3	2.10	0.52
1:B:157:ASN:OD1	1:B:321:GLN:NE2	2.43	0.52
1:D:81:LEU:O	1:D:82:ARG:C	2.53	0.52
1:B:206:LYS:HA	1:B:209:ARG:NH1	2.24	0.51
1:C:169:PHE:O	1:C:172:LEU:HB3	2.10	0.51
2:F:50:LYS:HB3	2:F:51:PRO:HD2	1.92	0.51
2:G:93:ILE:HG22	2:G:176:ASP:HA	1.92	0.51
1:C:16:THR:O	1:C:17:LEU:C	2.54	0.51
1:D:56:LEU:HD11	1:D:63:PHE:CD1	2.45	0.51
2:E:87:LEU:HD12	2:E:89:ILE:HD12	1.93	0.51
1:C:43:MET:SD	1:D:20:VAL:HG13	2.51	0.51
2:H:100:PRO:C	2:H:102:PHE:H	2.19	0.51
1:C:192:LEU:HD12	2:H:84:LEU:CA	2.41	0.51
1:C:39:TYR:CD2	1:C:103:ARG:HB3	2.44	0.51
2:F:27:GLU:N	2:F:28:PRO:CD	2.74	0.51
2:F:86:TYR:CG	2:F:86:TYR:O	2.63	0.51
2:G:50:LYS:HB3	2:G:51:PRO:HD2	1.93	0.51
2:G:112:ARG:C	2:G:113:ILE:HD12	2.36	0.51
2:G:193:ASP:CG	2:G:194:LYS:H	2.18	0.51
2:H:86:TYR:HE1	2:H:178:SER:HB3	1.76	0.51
2:H:108:GLN:O	2:H:109:ARG:HB3	2.10	0.51
1:B:54:THR:C	1:B:56:LEU:H	2.19	0.51
1:B:69:PHE:O	1:B:80:THR:HG23	2.10	0.51
2:F:37:GLN:NE2	2:F:47:ILE:HD11	2.25	0.51
2:E:113:ILE:HD13	2:E:132:ILE:HD13	1.93	0.50
2:H:38:ILE:C	2:H:38:ILE:CD1	2.84	0.50
2:E:22:ALA:O	2:E:23:ASP:HB2	2.11	0.50
1:A:24:ARG:HE	1:B:41:GLU:CD	2.20	0.50
1:D:81:LEU:HB3	1:D:111:PHE:HD1	1.76	0.50
2:H:66:GLY:O	2:H:182:ILE:HG12	2.11	0.50
1:A:157:ASN:OD1	1:A:321:GLN:NE2	2.44	0.50
1:D:206:LYS:HD3	1:D:206:LYS:O	2.12	0.50
1:A:16:THR:O	1:A:17:LEU:C	2.55	0.50
1:A:81:LEU:O	1:A:82:ARG:C	2.54	0.50
1:B:48:GLU:CG	1:B:52:LEU:HD22	2.41	0.50
1:C:73:LYS:HG2	1:C:77:GLN:H	1.76	0.50

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:93:TYR:CD2	1:A:295:LEU:HD11	2.47	0.50
1:B:109:LYS:HA	1:B:124:ASN:O	2.11	0.50
1:B:189:LYS:C	1:B:191:ASP:N	2.67	0.50
1:B:262:MET:SD	1:B:273:MET:HG3	2.52	0.50
1:C:49:TYR:O	1:C:50:THR:C	2.55	0.50
1:A:224:LEU:O	1:A:225:GLU:C	2.55	0.50
1:B:170:GLN:O	1:B:173:CYS:HB2	2.12	0.50
1:C:52:LEU:C	1:C:54:THR:H	2.19	0.50
2:F:114:ALA:HB1	2:F:137:LEU:CD1	2.41	0.50
1:A:81:LEU:HB3	1:A:111:PHE:CD1	2.47	0.49
1:A:146:LEU:O	1:A:150:VAL:HG23	2.12	0.49
1:D:206:LYS:HA	1:D:209:ARG:NH1	2.27	0.49
1:C:31:ARG:HG3	1:C:31:ARG:HH11	1.77	0.49
1:C:93:TYR:HD2	1:C:295:LEU:HD11	1.77	0.49
2:E:36:LEU:HB2	2:F:143:LEU:HD21	1.94	0.49
1:A:30:LEU:HD21	1:A:150:VAL:HG13	1.93	0.49
1:C:311:MET:O	1:C:312:ASP:C	2.55	0.49
1:D:157:ASN:OD1	1:D:321:GLN:NE2	2.45	0.49
2:E:14:GLN:NE2	2:E:89:ILE:HD12	2.27	0.49
2:F:16:THR:HG21	2:F:28:PRO:HG3	1.95	0.49
1:B:7:LEU:HD21	2:H:126:ALA:HA	1.95	0.49
1:B:81:LEU:O	1:B:82:ARG:C	2.55	0.49
1:C:27:GLU:HG2	1:C:127:ILE:HD11	1.94	0.49
1:A:319:GLU:C	1:A:321:GLN:N	2.67	0.49
1:D:31:ARG:HH11	1:D:31:ARG:HG3	1.78	0.49
2:E:112:ARG:O	2:E:150:ALA:HB1	2.13	0.49
1:C:85:PHE:O	1:C:88:PRO:HG2	2.13	0.49
2:F:4:ILE:HA	2:F:65:ILE:O	2.12	0.49
2:G:7:THR:HA	2:G:49:GLY:O	2.12	0.49
1:D:166:ALA:N	1:D:273:MET:HE2	2.28	0.49
1:C:81:LEU:HB3	1:C:111:PHE:HD1	1.73	0.49
2:G:75:GLU:OE2	2:G:119:ARG:HB2	2.13	0.49
2:H:85:LEU:CG	2:H:86:TYR:N	2.76	0.49
1:D:30:LEU:HD21	1:D:150:VAL:HG13	1.95	0.48
1:D:171:ARG:HG2	1:D:235:VAL:HB	1.94	0.48
2:E:92:CYS:O	2:E:93:ILE:HG23	2.13	0.48
2:F:93:ILE:HA	2:F:177:ILE:HG12	1.94	0.48
1:B:68:MET:HB2	1:B:80:THR:HG21	1.94	0.48
1:B:131:LEU:HD23	1:B:131:LEU:C	2.37	0.48
1:A:193:SER:OG	2:E:86:TYR:HB3	2.13	0.48
2:F:25:ASP:HB3	2:F:40:THR:HB	1.95	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:G:27:GLU:H	2:G:28:PRO:HD2	1.76	0.48
1:B:206:LYS:HD3	1:B:206:LYS:O	2.12	0.48
2:G:27:GLU:H	2:G:28:PRO:HD3	1.75	0.48
1:B:72:ILE:HG22	1:B:76:GLY:HA2	1.95	0.48
2:F:27:GLU:HA	2:F:27:GLU:OE2	2.12	0.48
2:G:112:ARG:HG3	2:G:150:ALA:HA	1.96	0.48
2:H:114:ALA:HA	2:H:135:ILE:O	2.14	0.48
1:A:85:PHE:O	1:A:88:PRO:HG2	2.14	0.48
1:A:116:ARG:CA	1:A:117:HIS:CD2	2.93	0.48
1:A:217:ILE:HB	2:E:196:ILE:CD1	2.44	0.48
1:A:317:ALA:O	1:A:321:GLN:HB2	2.14	0.48
1:B:171:ARG:HG2	1:B:235:VAL:HB	1.96	0.48
1:D:86:THR:O	1:D:90:VAL:HG23	2.13	0.48
2:G:30:LEU:HD12	2:G:30:LEU:N	2.28	0.48
1:A:72:ILE:HG22	1:A:73:LYS:N	2.29	0.48
1:D:214:GLU:HG2	1:D:227:LEU:HD11	1.96	0.48
2:E:112:ARG:HG3	2:E:150:ALA:HA	1.95	0.48
1:A:134:GLU:HG3	1:A:135:SER:N	2.29	0.48
1:A:226:ASN:O	1:A:227:LEU:C	2.57	0.48
1:B:319:GLU:C	1:B:321:GLN:N	2.71	0.48
2:H:58:LEU:C	2:H:58:LEU:HD23	2.39	0.48
1:C:81:LEU:O	1:C:82:ARG:C	2.56	0.48
1:C:148:LEU:HD11	1:C:256:ILE:HD11	1.96	0.48
2:G:92:CYS:O	2:G:93:ILE:HG23	2.14	0.48
1:A:23:LEU:HD11	1:A:125:TYR:CE2	2.49	0.48
1:D:186:LEU:CD2	1:D:191:ASP:HB3	2.44	0.48
2:E:93:ILE:HG13	2:E:157:VAL:O	2.14	0.48
2:G:5:ALA:HA	2:G:47:ILE:O	2.14	0.47
1:A:10:GLU:O	1:A:10:GLU:HG3	2.13	0.47
1:A:73:LYS:HG3	1:A:76:GLY:H	1.79	0.47
1:C:165:SER:C	1:C:273:MET:HE2	2.39	0.47
1:C:244:LYS:O	1:C:247:SER:HB3	2.13	0.47
1:B:117:HIS:O	1:B:118:LYS:HG2	2.14	0.47
1:D:209:ARG:HG2	1:D:213:LYS:HE2	1.96	0.47
2:F:93:ILE:O	2:F:156:ILE:HA	2.13	0.47
2:G:7:THR:CG2	2:G:68:VAL:HG12	2.44	0.47
2:G:100:PRO:C	2:G:102:PHE:H	2.23	0.47
2:H:70:LYS:HE3	2:H:74:ASP:OD1	2.15	0.47
2:H:7:THR:HA	2:H:49:GLY:O	2.15	0.47
1:A:56:LEU:HD22	1:A:297:ASN:CG	2.39	0.47
1:A:131:LEU:HD21	1:A:134:GLU:CB	2.42	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:86:THR:O	1:C:90:VAL:HG23	2.14	0.47
1:C:93:TYR:CD2	1:C:295:LEU:HD11	2.49	0.47
1:A:59:ASN:C	1:A:61:LYS:H	2.22	0.47
1:B:58:SER:C	1:B:267:ASP:HB2	2.39	0.47
1:C:173:CYS:O	1:C:176:ALA:N	2.43	0.47
2:F:113:ILE:N	2:F:113:ILE:HD12	2.30	0.47
1:A:103:ARG:HG2	1:A:103:ARG:HH11	1.78	0.47
1:A:217:ILE:HD12	1:A:217:ILE:C	2.39	0.47
1:C:68:MET:HB2	1:C:80:THR:CG2	2.44	0.47
1:C:108:GLY:O	1:C:126:GLN:N	2.45	0.47
1:C:134:GLU:HG3	1:C:135:SER:N	2.29	0.47
2:E:87:LEU:HD12	2:E:89:ILE:CD1	2.45	0.47
2:G:14:GLN:NE2	2:G:89:ILE:HB	2.29	0.47
1:A:279:SER:HA	2:E:190:PHE:CE1	2.49	0.47
1:C:71:PHE:HE1	1:C:79:ILE:HD11	1.78	0.47
1:D:262:MET:SD	1:D:273:MET:HG3	2.55	0.47
2:G:18:LEU:CD2	2:G:202:LEU:HD22	2.44	0.47
2:H:99:TYR:O	2:H:102:PHE:HB2	2.14	0.47
1:B:73:LYS:HG2	1:B:77:GLN:H	1.79	0.47
1:C:262:MET:SD	1:C:273:MET:HG3	2.55	0.47
2:E:17:LYS:O	2:E:20:GLU:HB3	2.15	0.47
2:E:55:ILE:HD11	2:E:73:LEU:HD23	1.96	0.47
2:F:82:TYR:O	2:F:182:ILE:HA	2.14	0.47
2:H:113:ILE:HG23	2:H:153:ILE:HA	1.96	0.47
1:A:275:LYS:HA	1:A:286:LEU:O	2.15	0.47
1:C:299:GLN:CG	1:C:300:GLU:H	2.10	0.47
2:E:70:LYS:HE3	2:E:74:ASP:OD1	2.15	0.47
2:E:82:TYR:O	2:E:182:ILE:HA	2.15	0.47
2:F:4:ILE:HG12	2:F:65:ILE:HB	1.96	0.47
2:F:84:LEU:CD1	2:F:183:VAL:HG21	2.45	0.47
2:H:142:GLU:O	2:H:145:PRO:HD2	2.15	0.47
1:B:71:PHE:HA	1:D:72:ILE:CG1	2.46	0.46
1:B:72:ILE:CG1	1:D:71:PHE:HA	2.45	0.46
1:D:68:MET:HB2	1:D:80:THR:CG2	2.45	0.46
2:E:113:ILE:HG21	2:E:154:VAL:HG23	1.97	0.46
2:F:94:PHE:CD1	2:F:177:ILE:HD11	2.50	0.46
2:H:85:LEU:HG	2:H:86:TYR:H	1.80	0.46
1:B:124:ASN:HA	1:B:312:ASP:OD2	2.15	0.46
1:D:7:LEU:H	1:D:7:LEU:CD2	2.27	0.46
1:C:8:LEU:HB3	1:C:12:SER:OG	2.15	0.46
1:D:266:MET:HE3	1:D:268:TYR:OH	2.16	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:115:SER:HA	2:E:141:VAL:HG11	1.95	0.46
2:E:143:LEU:HD21	2:F:36:LEU:CB	2.44	0.46
2:G:81:TYR:CD1	2:G:81:TYR:C	2.94	0.46
1:C:48:GLU:HG2	1:C:52:LEU:HD22	1.95	0.46
2:E:111:LYS:CE	2:E:130:GLU:OE2	2.59	0.46
2:F:7:THR:HG22	2:F:51:PRO:HA	1.98	0.46
2:F:121:THR:HG23	2:F:154:VAL:HG21	1.98	0.46
1:B:182:LEU:O	1:B:183:LEU:C	2.57	0.46
1:C:45:PRO:CB	1:D:12:SER:HB2	2.44	0.46
2:E:95:ALA:CA	2:E:174:ILE:HG12	2.45	0.46
1:A:226:ASN:O	1:A:228:VAL:N	2.48	0.46
2:G:85:LEU:CG	2:G:86:TYR:N	2.79	0.46
2:E:47:ILE:HG22	2:E:48:PHE:N	2.30	0.46
1:A:49:TYR:O	1:A:50:THR:C	2.58	0.46
1:B:58:SER:HA	1:B:267:ASP:HB3	1.98	0.46
1:C:16:THR:HG22	1:C:19:GLN:HG3	1.97	0.46
2:F:85:LEU:HG	2:F:86:TYR:N	2.30	0.46
1:B:74:HIS:C	1:B:76:GLY:N	2.71	0.46
1:C:199:ILE:HG21	1:C:209:ARG:CG	2.43	0.46
2:G:184:ASN:O	2:G:187:SER:HB2	2.15	0.46
1:A:20:VAL:O	1:A:21:LYS:C	2.58	0.46
1:A:73:LYS:HG2	1:A:77:GLN:N	2.20	0.46
1:C:279:SER:HA	2:H:190:PHE:CE1	2.51	0.46
1:D:45:PRO:HG2	1:D:48:GLU:OE1	2.16	0.46
1:A:16:THR:HG22	1:A:19:GLN:OE1	2.16	0.45
1:A:73:LYS:HE2	1:A:77:GLN:HG2	1.99	0.45
1:B:58:SER:HA	1:B:267:ASP:CB	2.46	0.45
2:F:100:PRO:C	2:F:102:PHE:H	2.24	0.45
2:G:94:PHE:CE2	2:G:156:ILE:HG23	2.51	0.45
2:H:7:THR:O	2:H:12:GLN:HB2	2.16	0.45
1:C:37:LYS:NZ	1:C:149:GLN:HE22	2.14	0.45
2:G:95:ALA:CA	2:G:174:ILE:HG12	2.46	0.45
2:F:24:TYR:C	2:F:26:VAL:H	2.25	0.45
1:C:206:LYS:HD3	1:C:206:LYS:O	2.17	0.45
2:E:5:ALA:HA	2:E:47:ILE:O	2.16	0.45
2:E:53:ASP:OD1	2:F:137:LEU:HA	2.17	0.45
2:H:75:GLU:OE1	2:H:120:VAL:HG23	2.16	0.45
1:B:94:SER:C	1:B:96:ILE:H	2.25	0.45
1:C:94:SER:O	1:C:96:ILE:N	2.49	0.45
1:C:26:ILE:O	1:C:27:GLU:C	2.60	0.45
1:C:131:LEU:HD23	1:C:131:LEU:O	2.17	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:56:LEU:HD21	1:D:63:PHE:CZ	2.51	0.45
2:G:17:LYS:O	2:G:20:GLU:N	2.49	0.45
1:A:176:ALA:HA	1:A:204:PHE:CZ	2.51	0.45
1:A:189:LYS:C	1:A:191:ASP:N	2.74	0.45
1:B:283:GLN:OE1	1:B:283:GLN:HA	2.17	0.45
2:E:75:GLU:CD	2:E:119:ARG:HB2	2.42	0.45
2:H:23:ASP:O	2:H:201:ARG:NH2	2.50	0.45
1:D:48:GLU:OE2	1:D:48:GLU:HA	2.17	0.45
1:D:175:LEU:HD11	1:D:233:ASP:OD1	2.17	0.45
2:E:37:GLN:HE21	2:E:45:GLN:NE2	2.15	0.45
2:E:93:ILE:HG22	2:E:176:ASP:HA	1.98	0.45
2:H:4:ILE:O	2:H:46:ILE:HA	2.17	0.45
1:C:64:ASN:O	1:C:68:MET:HE3	2.17	0.44
1:C:266:MET:HE3	1:C:268:TYR:OH	2.16	0.44
1:C:291:TYR:CD1	1:C:294:LEU:HD12	2.52	0.44
2:F:92:CYS:O	2:F:93:ILE:HG23	2.18	0.44
2:G:125:PHE:CD1	2:G:132:ILE:CD1	2.99	0.44
1:A:69:PHE:CZ	1:A:113:LYS:HG2	2.52	0.44
1:B:317:ALA:O	1:B:321:GLN:HB2	2.17	0.44
2:E:92:CYS:O	2:E:93:ILE:CG2	2.65	0.44
1:A:162:GLU:O	1:A:274:PHE:HB2	2.17	0.44
2:H:191:LYS:O	2:H:195:ILE:HG12	2.18	0.44
1:B:256:ILE:HG22	1:B:257:ILE:N	2.33	0.44
1:D:207:GLU:HG2	1:D:230:ASN:O	2.18	0.44
2:E:58:LEU:C	2:E:58:LEU:HD23	2.43	0.44
1:A:49:TYR:CZ	1:B:9:PRO:HA	2.52	0.44
1:A:264:PRO:HD3	1:A:273:MET:HE1	2.00	0.44
1:C:49:TYR:O	1:C:51:GLN:N	2.51	0.44
2:E:85:LEU:HG	2:E:86:TYR:N	2.33	0.44
1:A:71:PHE:CD2	1:B:71:PHE:CD2	3.06	0.44
2:F:47:ILE:HG22	2:F:48:PHE:N	2.33	0.44
2:H:15:VAL:O	2:H:16:THR:C	2.60	0.44
1:A:57:GLU:HB3	1:A:63:PHE:CE1	2.53	0.44
1:A:131:LEU:C	1:A:131:LEU:CD2	2.90	0.44
1:A:300:GLU:OE1	1:A:300:GLU:HA	2.17	0.44
2:E:142:GLU:O	2:E:145:PRO:HD2	2.17	0.44
2:H:7:THR:HG22	2:H:51:PRO:HA	1.99	0.44
1:A:71:PHE:CE2	1:B:71:PHE:HE2	2.33	0.44
1:D:247:SER:O	1:D:251:SER:HB3	2.18	0.44
1:D:319:GLU:C	1:D:321:GLN:N	2.75	0.44
2:E:124:TYR:CG	2:E:174:ILE:HD12	2.53	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:86:TYR:HE1	2:F:178:SER:HB3	1.81	0.44
1:B:117:HIS:C	1:B:118:LYS:HG3	2.42	0.43
1:B:320:ARG:O	1:B:320:ARG:HG2	2.18	0.43
2:E:14:GLN:O	2:E:17:LYS:HB3	2.18	0.43
1:A:237:ILE:HG13	1:A:238:SER:N	2.32	0.43
1:C:168:PHE:CE2	1:C:260:LEU:HB3	2.54	0.43
1:C:198:PHE:CE2	1:C:204:PHE:HE1	2.36	0.43
1:C:228:VAL:HG12	1:C:237:ILE:HG22	2.00	0.43
1:D:32:LYS:HE3	1:D:32:LYS:HB2	1.89	0.43
1:D:106:TYR:O	1:D:127:ILE:HA	2.17	0.43
2:E:86:TYR:O	2:E:86:TYR:CG	2.69	0.43
2:E:112:ARG:H	2:E:112:ARG:HG2	1.69	0.43
2:F:1:MET:HE1	2:F:39:LYS:HE3	1.99	0.43
2:F:14:GLN:NE2	2:F:89:ILE:HB	2.33	0.43
2:H:29:ILE:HG12	2:H:29:ILE:O	2.18	0.43
1:A:52:LEU:HD23	1:A:53:TYR:CD2	2.53	0.43
1:B:93:TYR:CD2	1:B:295:LEU:HD11	2.53	0.43
1:B:93:TYR:HD2	1:B:295:LEU:HD11	1.82	0.43
1:C:106:TYR:OH	1:C:308:CYS:HB2	2.18	0.43
1:D:237:ILE:HG13	1:D:238:SER:N	2.32	0.43
2:G:7:THR:O	2:G:12:GLN:HB2	2.18	0.43
1:C:94:SER:C	1:C:96:ILE:N	2.77	0.43
2:E:30:LEU:HD22	2:E:38:ILE:HG21	2.01	0.43
1:B:48:GLU:OE2	1:B:91:ARG:NH1	2.40	0.43
1:D:16:THR:O	1:D:17:LEU:C	2.62	0.43
1:D:16:THR:HG22	1:D:19:GLN:CG	2.49	0.43
1:D:224:LEU:O	1:D:228:VAL:HG23	2.18	0.43
2:F:58:LEU:HD13	2:F:182:ILE:CD1	2.47	0.43
2:G:1:MET:HB2	2:G:43:ASP:O	2.18	0.43
2:G:87:LEU:O	2:G:89:ILE:HG13	2.17	0.43
2:G:158:GLU:HB3	3:G:4009:PRP:O2P	2.18	0.43
2:H:81:TYR:CD1	2:H:81:TYR:C	2.96	0.43
1:B:16:THR:HG22	1:B:19:GLN:OE1	2.19	0.43
2:G:36:LEU:HD12	2:G:36:LEU:N	2.33	0.43
2:G:86:TYR:HE1	2:G:178:SER:HB3	1.82	0.43
2:H:6:ILE:HG22	2:H:7:THR:N	2.33	0.43
2:H:190:PHE:C	2:H:192:LYS:H	2.26	0.43
1:B:189:LYS:O	1:B:190:LYS:C	2.60	0.43
1:C:68:MET:HB2	1:C:80:THR:HG21	2.01	0.43
1:C:205:SER:O	1:C:206:LYS:C	2.61	0.43
1:A:37:LYS:O	1:A:38:ASN:HB2	2.18	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:117:HIS:CD2	1:A:117:HIS:N	2.77	0.43
1:A:207:GLU:CD	1:A:207:GLU:H	2.27	0.43
1:B:101:SER:HA	1:B:132:PHE:O	2.19	0.43
1:B:192:LEU:HD12	2:G:199:VAL:HG11	2.00	0.43
1:B:316:LYS:HB2	1:B:316:LYS:HE3	1.77	0.43
2:E:36:LEU:HD13	2:F:143:LEU:CD1	2.48	0.43
2:E:92:CYS:C	2:E:93:ILE:HG23	2.44	0.43
2:G:114:ALA:HB1	2:G:137:LEU:HD12	2.00	0.43
1:A:206:LYS:HA	1:A:209:ARG:NH1	2.33	0.43
1:C:189:LYS:O	1:C:190:LYS:C	2.62	0.43
2:H:6:ILE:HG21	2:H:12:GLN:HA	2.01	0.43
1:B:217:ILE:C	1:B:217:ILE:HD12	2.43	0.43
1:D:163:ILE:O	1:D:163:ILE:HG13	2.19	0.43
2:F:18:LEU:CD2	2:F:202:LEU:HD22	2.44	0.43
2:F:196:ILE:O	2:F:197:GLU:C	2.61	0.43
1:D:8:LEU:HB3	2:E:131:ASP:OD2	2.18	0.42
1:D:214:GLU:HG2	1:D:227:LEU:CD1	2.49	0.42
2:E:143:LEU:O	2:E:146:VAL:N	2.52	0.42
1:C:7:LEU:N	1:C:7:LEU:HD12	2.33	0.42
1:C:166:ALA:O	1:C:167:LYS:C	2.62	0.42
2:G:59:GLU:HG3	2:G:78:PHE:CZ	2.51	0.42
1:C:275:LYS:HD3	1:C:277:TYR:OH	2.19	0.42
1:D:256:ILE:HG22	1:D:257:ILE:N	2.34	0.42
2:E:86:TYR:CE2	2:E:88:LYS:HG2	2.54	0.42
2:G:114:ALA:HA	2:G:135:ILE:O	2.19	0.42
2:H:24:TYR:HA	2:H:42:ASP:OD2	2.20	0.42
1:A:71:PHE:HE1	1:A:79:ILE:HD11	1.83	0.42
1:B:56:LEU:HD11	1:B:87:LEU:HD11	2.02	0.42
1:B:73:LYS:HG2	1:B:77:GLN:N	2.34	0.42
1:C:144:LEU:HD23	1:C:144:LEU:HA	1.64	0.42
2:E:6:ILE:HG22	2:E:7:THR:N	2.34	0.42
2:E:113:ILE:HG23	2:E:153:ILE:CA	2.45	0.42
1:B:16:THR:O	1:B:19:GLN:N	2.53	0.42
1:B:44:PRO:HG2	1:B:85:PHE:CD2	2.54	0.42
1:C:27:GLU:OE2	1:C:107:PHE:HB3	2.19	0.42
1:C:207:GLU:CD	1:C:207:GLU:H	2.27	0.42
2:E:98:SER:HB2	2:E:152:ALA:CB	2.49	0.42
2:H:143:LEU:O	2:H:144:GLY:C	2.62	0.42
1:A:195:LEU:O	1:A:198:PHE:HB3	2.20	0.42
1:B:319:GLU:C	1:B:321:GLN:H	2.26	0.42
1:C:275:LYS:HD3	1:C:277:TYR:HE1	1.83	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:16:THR:HG22	1:D:19:GLN:HG3	2.00	0.42
2:F:58:LEU:HD23	2:F:58:LEU:O	2.18	0.42
1:C:134:GLU:HB3	1:C:139:SER:OG	2.19	0.42
1:D:129:ILE:HD11	1:D:307:PHE:CE2	2.55	0.42
1:D:186:LEU:HD13	1:D:195:LEU:HA	2.02	0.42
2:E:96:LEU:O	2:E:171:ILE:HG12	2.20	0.42
2:E:141:VAL:O	2:E:141:VAL:HG12	2.20	0.42
1:A:93:TYR:HD2	1:A:295:LEU:HD11	1.82	0.42
1:A:294:LEU:HA	1:A:297:ASN:ND2	2.35	0.42
1:B:266:MET:HE3	1:B:268:TYR:OH	2.18	0.42
1:D:205:SER:O	1:D:206:LYS:C	2.60	0.42
2:G:29:ILE:HG22	2:G:30:LEU:H	1.84	0.42
2:H:85:LEU:CG	2:H:86:TYR:H	2.31	0.42
1:A:6:TYR:HB3	1:A:7:LEU:H	1.59	0.42
1:A:24:ARG:HG3	1:B:41:GLU:OE2	2.19	0.42
1:B:269:TYR:CE1	1:B:291:TYR:HD2	2.37	0.42
1:D:98:ASP:O	1:D:100:THR:N	2.48	0.42
2:G:94:PHE:CD1	2:G:177:ILE:HD11	2.55	0.42
2:E:139:GLY:O	2:E:140:SER:C	2.63	0.41
2:F:5:ALA:HA	2:F:47:ILE:O	2.20	0.41
2:G:51:PRO:HG2	2:G:52:ASN:H	1.84	0.41
1:A:71:PHE:CE1	1:A:79:ILE:CD1	3.02	0.41
1:B:30:LEU:HD22	1:B:150:VAL:HG13	2.02	0.41
1:B:54:THR:C	1:B:56:LEU:N	2.78	0.41
1:C:52:LEU:HD23	1:C:53:TYR:CD2	2.55	0.41
1:C:217:ILE:C	1:C:217:ILE:HD12	2.45	0.41
2:E:193:ASP:CG	2:E:194:LYS:N	2.75	0.41
2:F:20:GLU:O	2:F:23:ASP:N	2.41	0.41
2:F:99:TYR:O	2:F:102:PHE:HB2	2.20	0.41
2:F:113:ILE:HA	2:F:152:ALA:O	2.19	0.41
2:H:82:TYR:CE2	2:H:185:LYS:HE2	2.55	0.41
2:H:95:ALA:HA	2:H:174:ILE:HG12	2.02	0.41
1:A:33:LEU:O	1:A:36:LEU:HB3	2.19	0.41
1:A:298:PHE:HB2	1:A:299:GLN:H	1.48	0.41
1:B:110:ILE:HG13	1:B:124:ASN:HB2	2.02	0.41
1:C:146:LEU:O	1:C:150:VAL:HG23	2.19	0.41
1:C:173:CYS:O	1:C:174:GLN:C	2.63	0.41
2:E:177:ILE:HG22	2:E:178:SER:N	2.35	0.41
2:G:22:ALA:O	2:G:23:ASP:HB2	2.20	0.41
2:G:86:TYR:CE1	2:G:178:SER:HB3	2.55	0.41
1:C:189:LYS:C	1:C:191:ASP:N	2.73	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:7:THR:HG22	2:E:51:PRO:HA	2.01	0.41
2:G:17:LYS:O	2:G:20:GLU:HB3	2.20	0.41
2:G:113:ILE:HA	2:G:152:ALA:O	2.20	0.41
2:H:87:LEU:HD23	2:H:87:LEU:N	2.16	0.41
1:B:8:LEU:H	1:B:8:LEU:HG	1.73	0.41
1:B:117:HIS:C	1:B:118:LYS:CG	2.93	0.41
1:B:191:ASP:HA	2:G:83:GLU:O	2.20	0.41
1:D:120:ARG:HG2	1:D:120:ARG:HH11	1.83	0.41
1:D:274:PHE:N	1:D:274:PHE:CD2	2.88	0.41
2:E:87:LEU:H	2:E:87:LEU:CD2	2.15	0.41
2:H:85:LEU:HD12	2:H:86:TYR:H	1.86	0.41
1:B:68:MET:HB2	1:B:80:THR:CG2	2.51	0.41
1:C:82:ARG:NH2	1:C:109:LYS:O	2.44	0.41
1:C:199:ILE:HD13	1:C:209:ARG:CG	2.42	0.41
1:C:231:THR:O	1:C:232:LYS:C	2.62	0.41
1:C:298:PHE:CD1	1:C:298:PHE:N	2.88	0.41
2:E:19:LEU:HD23	2:E:202:LEU:HD21	2.01	0.41
2:E:38:ILE:C	2:E:38:ILE:CD1	2.93	0.41
2:F:58:LEU:HD13	2:F:182:ILE:HG13	2.02	0.41
2:F:158:GLU:HB3	3:F:4008:PRP:O3P	2.21	0.41
2:H:30:LEU:HD11	2:H:38:ILE:CG2	2.47	0.41
2:H:112:ARG:HG3	2:H:150:ALA:HA	2.02	0.41
1:B:136:ALA:O	1:B:139:SER:HB2	2.21	0.41
1:C:214:GLU:C	1:C:216:PHE:N	2.78	0.41
1:C:214:GLU:O	1:C:215:ILE:C	2.63	0.41
1:D:134:GLU:HG3	1:D:135:SER:N	2.36	0.41
1:D:319:GLU:C	1:D:321:GLN:H	2.27	0.41
1:A:101:SER:HA	1:A:132:PHE:O	2.20	0.41
1:A:116:ARG:O	1:A:117:HIS:ND1	2.45	0.41
1:A:161:PHE:HD2	1:A:256:ILE:HD12	1.86	0.41
1:B:244:LYS:O	1:B:247:SER:HB3	2.21	0.41
1:B:268:TYR:CE1	1:B:294:LEU:HD11	2.56	0.41
1:C:6:TYR:HB3	1:C:7:LEU:H	1.69	0.41
1:C:90:VAL:HG22	1:C:132:PHE:CZ	2.56	0.41
1:C:147:ALA:O	1:C:151:ILE:HG13	2.20	0.41
2:E:40:THR:OG1	2:E:44:LEU:HB2	2.21	0.41
2:F:188:PHE:O	2:F:192:LYS:HB2	2.20	0.41
2:H:38:ILE:O	2:H:45:GLN:HA	2.20	0.41
2:H:67:PHE:CE2	2:H:181:MET:HG3	2.55	0.41
2:H:86:TYR:CD1	2:H:86:TYR:C	2.96	0.41
2:H:204:ASP:O	2:H:205:ALA:HB2	2.21	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:44:PRO:HB2	1:A:82:ARG:HD3	2.02	0.41
1:B:87:LEU:HD23	1:B:87:LEU:HA	1.81	0.41
1:C:73:LYS:CE	1:C:77:GLN:HB2	2.40	0.41
1:C:131:LEU:HD21	1:C:134:GLU:CB	2.46	0.41
1:D:30:LEU:CD2	1:D:150:VAL:HG13	2.51	0.41
1:D:175:LEU:HD12	1:D:236:LEU:HD11	2.03	0.41
2:G:139:GLY:O	2:G:141:VAL:N	2.54	0.41
2:H:50:LYS:O	2:H:53:ASP:HB2	2.20	0.41
1:B:31:ARG:HH11	1:B:31:ARG:HG3	1.85	0.40
1:B:73:LYS:HE3	1:B:75:GLU:CB	2.33	0.40
1:C:224:LEU:O	1:C:225:GLU:C	2.63	0.40
2:E:58:LEU:CD1	2:E:182:ILE:HG13	2.50	0.40
2:H:93:ILE:HG13	2:H:157:VAL:O	2.20	0.40
1:A:31:ARG:HH21	1:A:41:GLU:CD	2.29	0.40
1:A:87:LEU:N	1:A:88:PRO:HD2	2.36	0.40
1:B:186:LEU:HA	1:B:186:LEU:HD23	1.88	0.40
1:B:282:ASN:HA	2:G:186:SER:HB2	2.03	0.40
1:D:259:ASP:OD1	1:D:262:MET:N	2.54	0.40
1:D:168:PHE:CZ	1:D:260:LEU:HD22	2.57	0.40
2:F:142:GLU:HG2	2:F:162:THR:HG22	2.03	0.40
2:G:3:LYS:HB2	2:G:63:VAL:HA	2.04	0.40
1:A:16:THR:O	1:A:19:GLN:N	2.55	0.40
1:D:143:ILE:HD13	1:D:143:ILE:HA	1.97	0.40
2:E:86:TYR:CZ	2:E:88:LYS:HG2	2.57	0.40
2:F:17:LYS:O	2:F:20:GLU:HB3	2.22	0.40
2:F:17:LYS:HE3	2:F:21:ASN:HD21	1.87	0.40
1:A:106:TYR:CZ	1:A:128:GLY:HA3	2.57	0.40
1:B:137:ASP:O	1:B:138:LYS:C	2.64	0.40
1:B:166:ALA:N	1:B:273:MET:HE2	2.37	0.40
1:B:240:PHE:CD2	1:B:240:PHE:C	2.99	0.40
1:B:289:GLY:O	1:B:305:ILE:HG13	2.22	0.40
2:G:82:TYR:O	2:G:182:ILE:HA	2.21	0.40
2:H:86:TYR:CE1	2:H:178:SER:HB3	2.54	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	309/344 (90%)	272 (88%)	32 (10%)	5 (2%)	7	30
1	B	316/344 (92%)	275 (87%)	37 (12%)	4 (1%)	9	34
1	C	298/344 (87%)	265 (89%)	29 (10%)	4 (1%)	9	34
1	D	308/344 (90%)	285 (92%)	20 (6%)	3 (1%)	12	40
2	E	199/208 (96%)	165 (83%)	31 (16%)	3 (2%)	8	32
2	F	203/208 (98%)	170 (84%)	25 (12%)	8 (4%)	2	14
2	G	199/208 (96%)	167 (84%)	24 (12%)	8 (4%)	2	14
2	H	196/208 (94%)	162 (83%)	29 (15%)	5 (3%)	4	21
All	All	2028/2208 (92%)	1761 (87%)	227 (11%)	40 (2%)	6	26

All (40) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	F	84	LEU
2	G	84	LEU
2	G	140	SER
1	A	191	ASP
1	B	115	LYS
1	B	136	ALA
1	B	191	ASP
1	C	95	GLN
1	C	136	ALA
1	C	191	ASP
1	D	191	ASP
2	E	84	LEU
2	F	9	GLY
2	F	88	LYS
2	G	88	LYS
2	H	9	GLY
2	H	25	ASP

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
2	H	84	LEU
1	A	227	LEU
1	A	320	ARG
2	F	26	VAL
2	F	33	GLY
2	F	34	ARG
2	F	87	LEU
2	F	193	ASP
2	G	101	ASP
1	A	136	ALA
1	A	299	GLN
2	H	101	ASP
1	C	203	ASN
1	D	99	SER
2	E	101	ASP
2	G	29	ILE
2	G	91	GLN
2	H	109	ARG
1	D	271	ASP
2	G	193	ASP
1	B	9	PRO
2	E	9	GLY
2	G	27	GLU

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	284/311 (91%)	272 (96%)	12 (4%)	26	53
1	B	288/311 (93%)	274 (95%)	14 (5%)	22	50
1	C	276/311 (89%)	265 (96%)	11 (4%)	28	54
1	D	283/311 (91%)	273 (96%)	10 (4%)	32	56
2	E	182/186 (98%)	178 (98%)	4 (2%)	45	65
2	F	183/186 (98%)	178 (97%)	5 (3%)	39	61

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
2	G	182/186 (98%)	178 (98%)	4 (2%)	45 65
2	H	179/186 (96%)	176 (98%)	3 (2%)	53 69
All	All	1857/1988 (93%)	1794 (97%)	63 (3%)	32 57

All (63) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	42	VAL
1	A	75	GLU
1	A	82	ARG
1	A	96	ILE
1	A	117	HIS
1	A	135	SER
1	A	174	GLN
1	A	180	THR
1	A	184	THR
1	A	225	GLU
1	A	252	MET
1	A	295	LEU
1	B	8	LEU
1	B	10	GLU
1	B	11	GLU
1	B	42	VAL
1	B	43	MET
1	B	62	THR
1	B	141	LEU
1	B	174	GLN
1	B	180	THR
1	B	189	LYS
1	B	219	ASN
1	B	225	GLU
1	B	252	MET
1	B	295	LEU
1	C	42	VAL
1	C	52	LEU
1	C	75	GLU
1	C	113	LYS
1	C	174	GLN
1	C	219	ASN
1	C	225	GLU
1	C	238	SER

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	C	252	MET
1	C	295	LEU
1	C	298	PHE
1	D	42	VAL
1	D	122	THR
1	D	135	SER
1	D	141	LEU
1	D	174	GLN
1	D	219	ASN
1	D	225	GLU
1	D	252	MET
1	D	295	LEU
1	D	297	ASN
2	E	43	ASP
2	E	87	LEU
2	E	107	PHE
2	E	110	HIS
2	F	31	ASN
2	F	87	LEU
2	F	109	ARG
2	F	115	SER
2	F	204	ASP
2	G	26	VAL
2	G	43	ASP
2	G	87	LEU
2	G	110	HIS
2	H	43	ASP
2	H	87	LEU
2	H	107	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (45) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	70	GLN
1	A	117	HIS
1	A	124	ASN
1	A	126	GLN
1	A	149	GLN
1	B	64	ASN
1	B	65	GLN
1	B	70	GLN
1	B	124	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	196	ASN
1	C	124	ASN
1	C	149	GLN
1	C	174	GLN
1	C	299	GLN
1	D	40	GLN
1	D	70	GLN
1	D	95	GLN
1	D	124	ASN
1	D	174	GLN
1	D	196	ASN
1	D	297	ASN
1	D	299	GLN
1	D	310	HIS
2	E	14	GLN
2	E	21	ASN
2	E	37	GLN
2	E	45	GLN
2	F	14	GLN
2	F	21	ASN
2	F	31	ASN
2	F	37	GLN
2	F	45	GLN
2	F	60	HIS
2	F	106	ASN
2	F	161	ASN
2	G	14	GLN
2	G	21	ASN
2	G	37	GLN
2	G	45	GLN
2	G	161	ASN
2	G	206	GLN
2	H	14	GLN
2	H	21	ASN
2	H	37	GLN
2	H	60	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

6 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	PRP	E	4007	-	20,22,22	1.27	2 (10%)	32,35,35	1.02	1 (3%)
4	PO4	F	1002	-	4,4,4	1.59	1 (25%)	6,6,6	0.45	0
3	PRP	F	4008	-	20,22,22	1.04	1 (5%)	32,35,35	0.94	2 (6%)
4	PO4	G	1001	-	4,4,4	1.59	0	6,6,6	0.45	0
3	PRP	H	4010	-	20,22,22	1.17	2 (10%)	32,35,35	0.98	1 (3%)
3	PRP	G	4009	-	20,22,22	1.10	2 (10%)	32,35,35	0.97	1 (3%)

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	PRP	H	4010	-	-	2/16/33/33	0/1/1/1
3	PRP	E	4007	-	-	4/16/33/33	0/1/1/1
3	PRP	G	4009	-	-	5/16/33/33	0/1/1/1
3	PRP	F	4008	-	-	3/16/33/33	0/1/1/1

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	4007	PRP	O4-C4	3.23	1.52	1.45
3	E	4007	PRP	C1-C2	2.71	1.56	1.52
3	H	4010	PRP	O4-C4	2.64	1.50	1.45
3	H	4010	PRP	PA-O3A	2.58	1.62	1.59
3	F	4008	PRP	O4-C4	2.56	1.50	1.45
3	G	4009	PRP	O4-C4	2.50	1.50	1.45
3	G	4009	PRP	PA-O3A	2.15	1.61	1.59
4	F	1002	PO4	P-O2	-2.00	1.48	1.54

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	4008	PRP	O2A-PA-O3A	2.14	113.05	107.27
3	G	4009	PRP	O2A-PA-O3A	2.11	112.99	107.27
3	F	4008	PRP	O3P-P-O5	-2.10	101.18	106.67
3	E	4007	PRP	O2A-PA-O3A	2.05	112.83	107.27
3	H	4010	PRP	O2A-PA-O3A	2.03	112.77	107.27

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	E	4007	PRP	PA-O3A-PB-O2B
3	F	4008	PRP	C5-O5-P-O1P
3	E	4007	PRP	O4-C4-C5-O5
3	E	4007	PRP	C3-C4-C5-O5
3	G	4009	PRP	C3-C4-C5-O5
3	F	4008	PRP	C5-O5-P-O3P
3	G	4009	PRP	C1-O1-PA-O1A
3	G	4009	PRP	C1-O1-PA-O2A
3	H	4010	PRP	C1-O1-PA-O1A
3	E	4007	PRP	PA-O3A-PB-O3B
3	G	4009	PRP	O4-C4-C5-O5
3	F	4008	PRP	C4-C5-O5-P
3	G	4009	PRP	C2-C1-O1-PA
3	H	4010	PRP	C2-C1-O1-PA

There are no ring outliers.

3 monomers are involved in 4 short contacts:

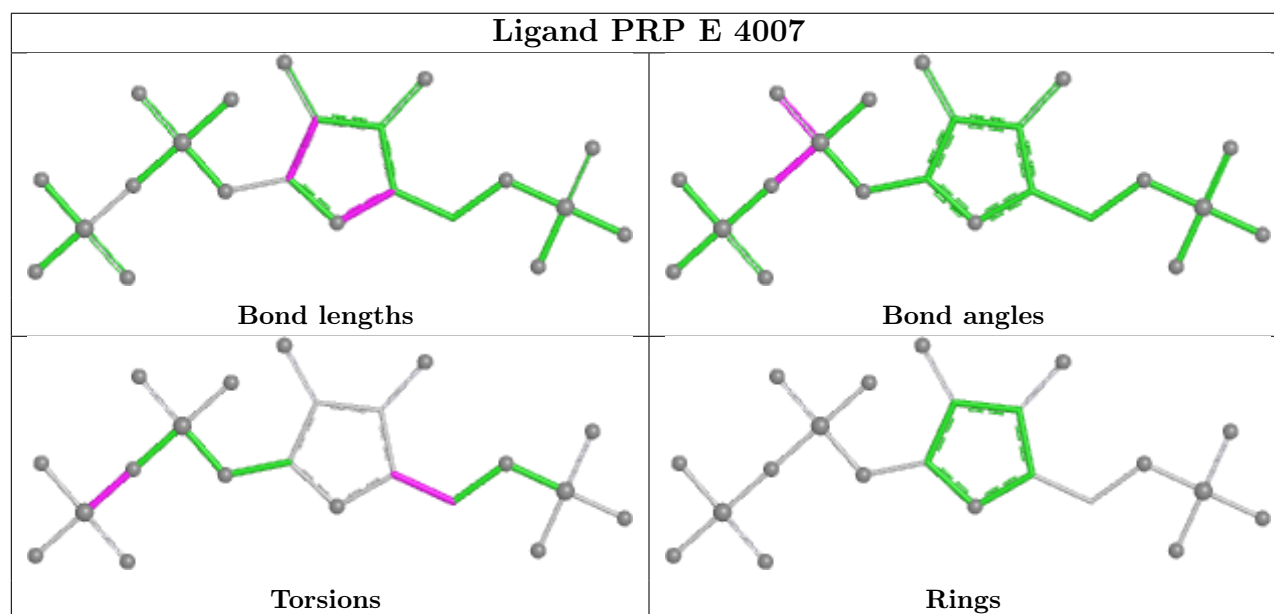
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	F	4008	PRP	1	0

Continued on next page...

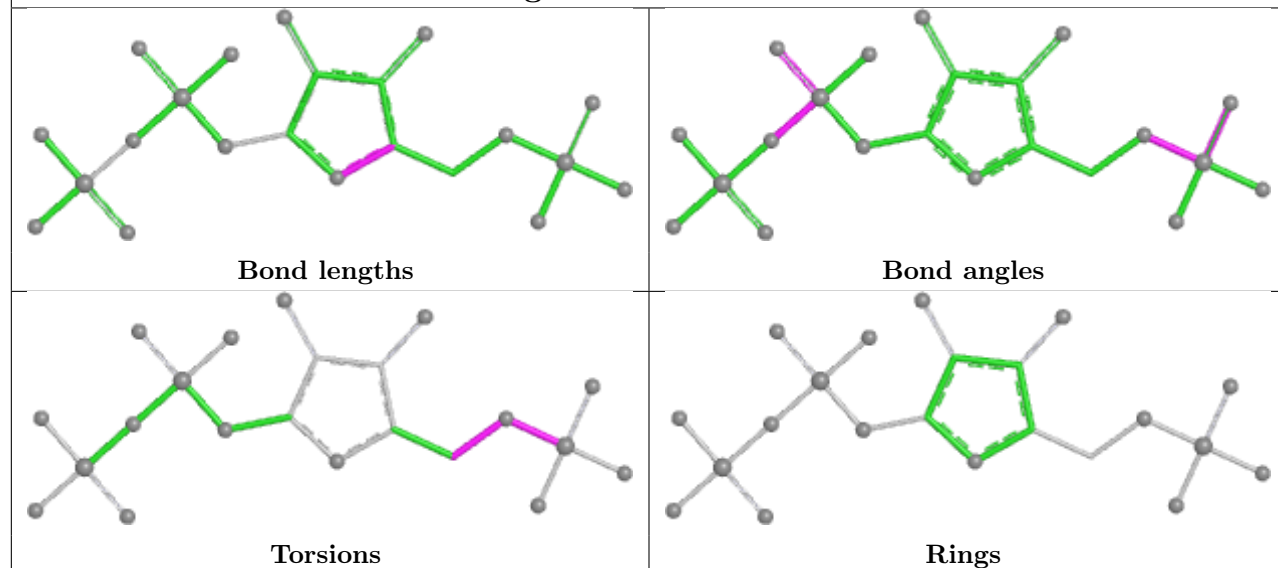
Continued from previous page...

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	G	1001	PO4	2	0
3	G	4009	PRP	1	0

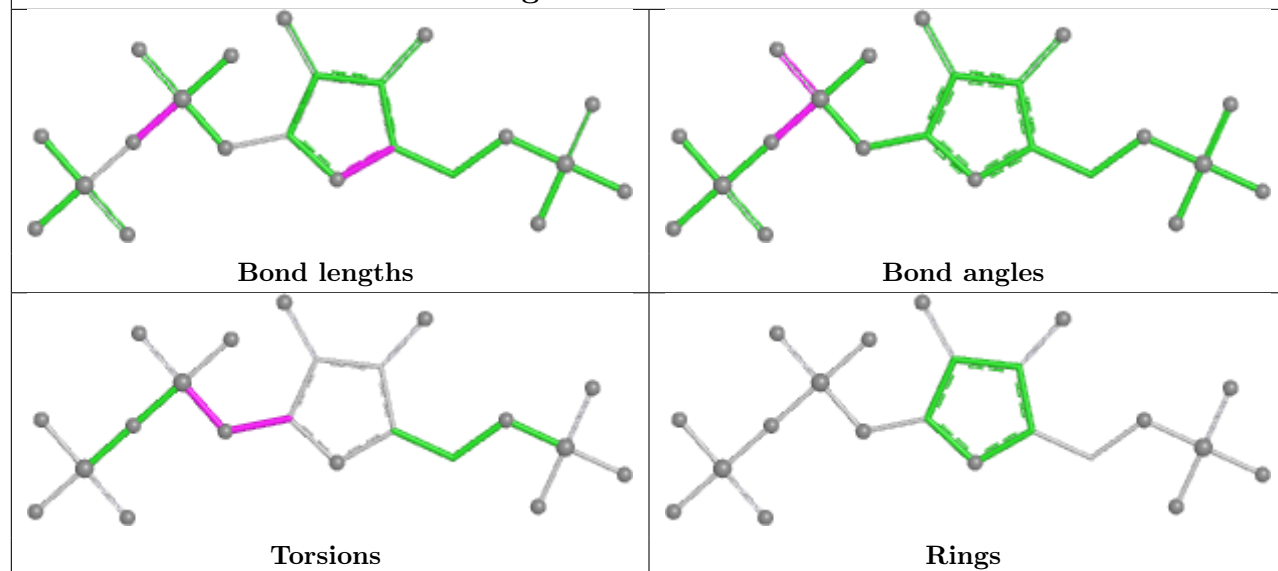
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

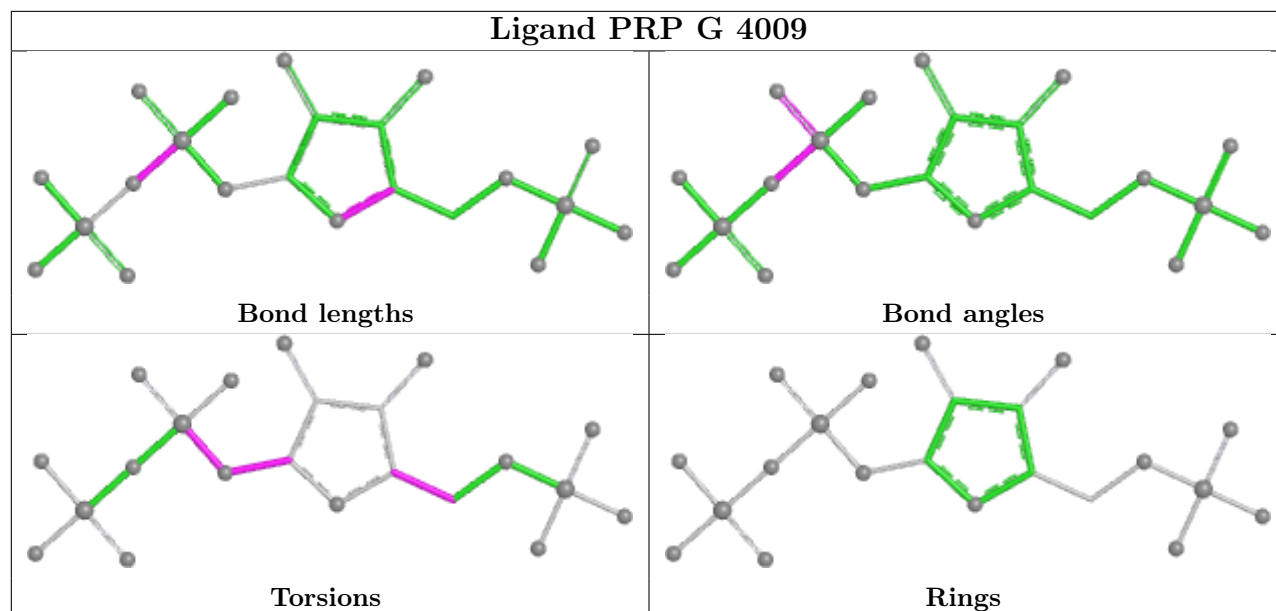


Ligand PRP F 4008



Ligand PRP H 4010





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
2	E	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	E	29:ILE	C	30:LEU	N	1.04

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	311/344 (90%)	-0.07	8 (2%)	57	38	14, 42, 117, 169	2 (0%)
1	B	316/344 (91%)	0.12	5 (1%)	70	52	17, 55, 119, 176	0
1	C	304/344 (88%)	0.02	5 (1%)	70	52	17, 50, 109, 182	3 (0%)
1	D	312/344 (90%)	0.26	9 (2%)	53	36	22, 70, 126, 166	1 (0%)
2	E	202/208 (97%)	0.17	5 (2%)	58	39	25, 61, 120, 166	0
2	F	205/208 (98%)	0.40	8 (3%)	43	29	26, 71, 126, 162	0
2	G	202/208 (97%)	0.26	6 (2%)	52	35	25, 62, 119, 164	0
2	H	200/208 (96%)	0.30	3 (1%)	72	53	27, 65, 129, 168	0
All	All	2052/2208 (92%)	0.16	49 (2%)	59	40	14, 59, 123, 182	6 (0%)

All (49) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	71	PHE	5.4
1	A	117	HIS	4.5
1	A	116	ARG	3.7
1	D	74	HIS	3.6
1	B	303	PHE	3.5
2	E	30	LEU	3.4
2	G	36	LEU	3.3
1	C	7	LEU	3.3
1	B	74	HIS	3.2
1	A	298	PHE	3.2
1	A	74	HIS	3.1
1	C	74	HIS	3.0
2	F	98	SER	3.0
1	C	56	LEU	3.0
2	F	85	LEU	2.8
2	G	28	PRO	2.8

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	C	71	PHE	2.7
2	G	29	ILE	2.7
2	H	139	GLY	2.7
1	D	117	HIS	2.6
1	B	14	GLU	2.6
2	F	36	LEU	2.5
2	G	30	LEU	2.5
1	A	71	PHE	2.5
1	D	6	TYR	2.5
2	E	32	LEU	2.4
2	E	27	GLU	2.4
2	H	173	LYS	2.4
1	A	266	MET	2.4
2	F	114	ALA	2.3
2	G	138	GLU	2.3
1	C	177	ASP	2.3
1	D	295	LEU	2.3
2	F	30	LEU	2.3
1	D	203	ASN	2.3
1	D	283	GLN	2.2
2	F	32	LEU	2.2
2	G	110	HIS	2.2
1	A	7	LEU	2.2
1	B	75	GLU	2.2
2	E	167	GLY	2.2
2	F	199	VAL	2.2
1	D	71	PHE	2.2
1	A	14	GLU	2.1
1	D	56	LEU	2.1
1	D	63	PHE	2.1
2	F	8	LYS	2.0
2	H	133	GLU	2.0
2	E	36	LEU	2.0

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

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

6.3 Carbohydrates [i](#)

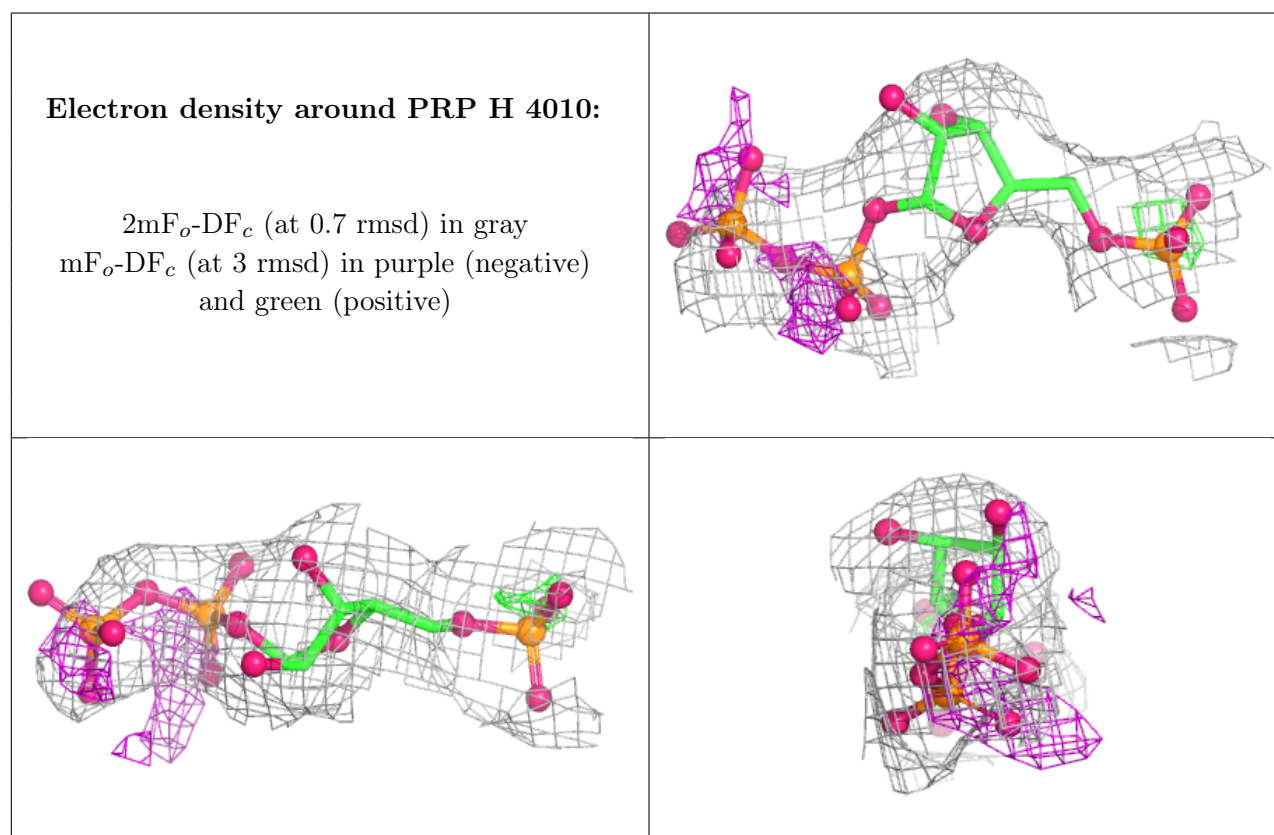
There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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.

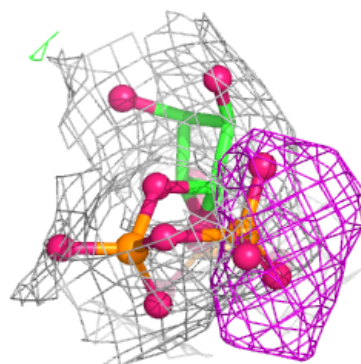
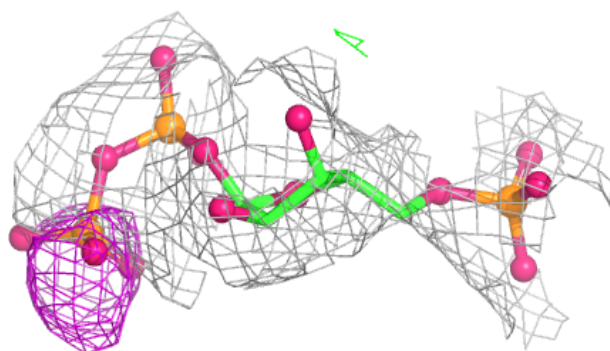
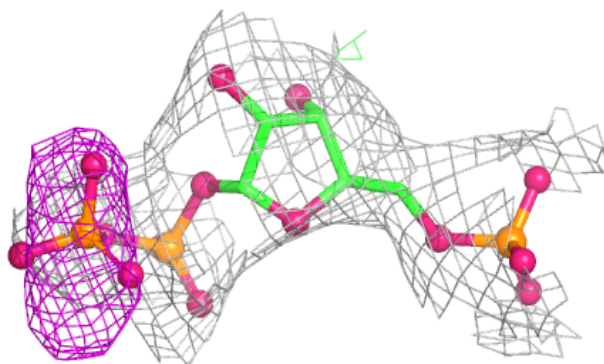
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	PRP	H	4010	22/22	0.56	0.17	138,138,138,138	0
3	PRP	G	4009	22/22	0.57	0.15	118,118,118,118	0
3	PRP	E	4007	22/22	0.62	0.14	113,113,113,113	0
3	PRP	F	4008	22/22	0.63	0.14	117,117,117,117	0
4	PO4	F	1002	5/5	0.87	0.13	111,111,111,111	0
4	PO4	G	1001	5/5	0.88	0.15	127,127,127,127	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.

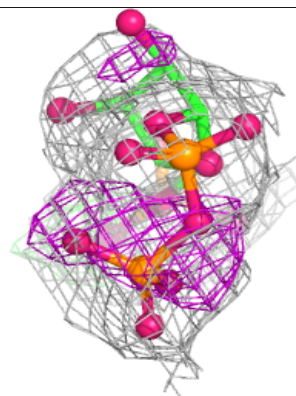
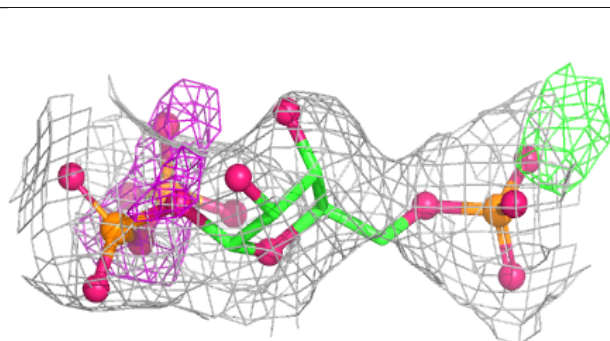
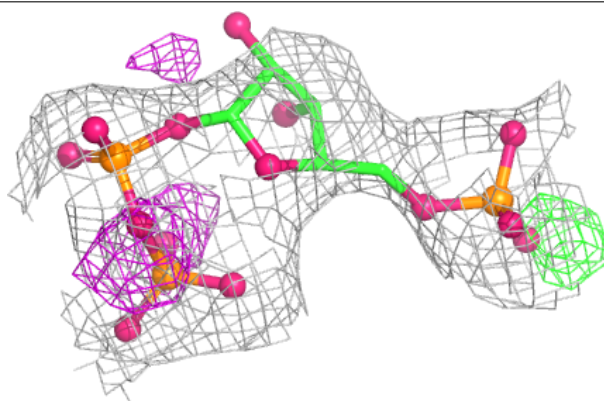


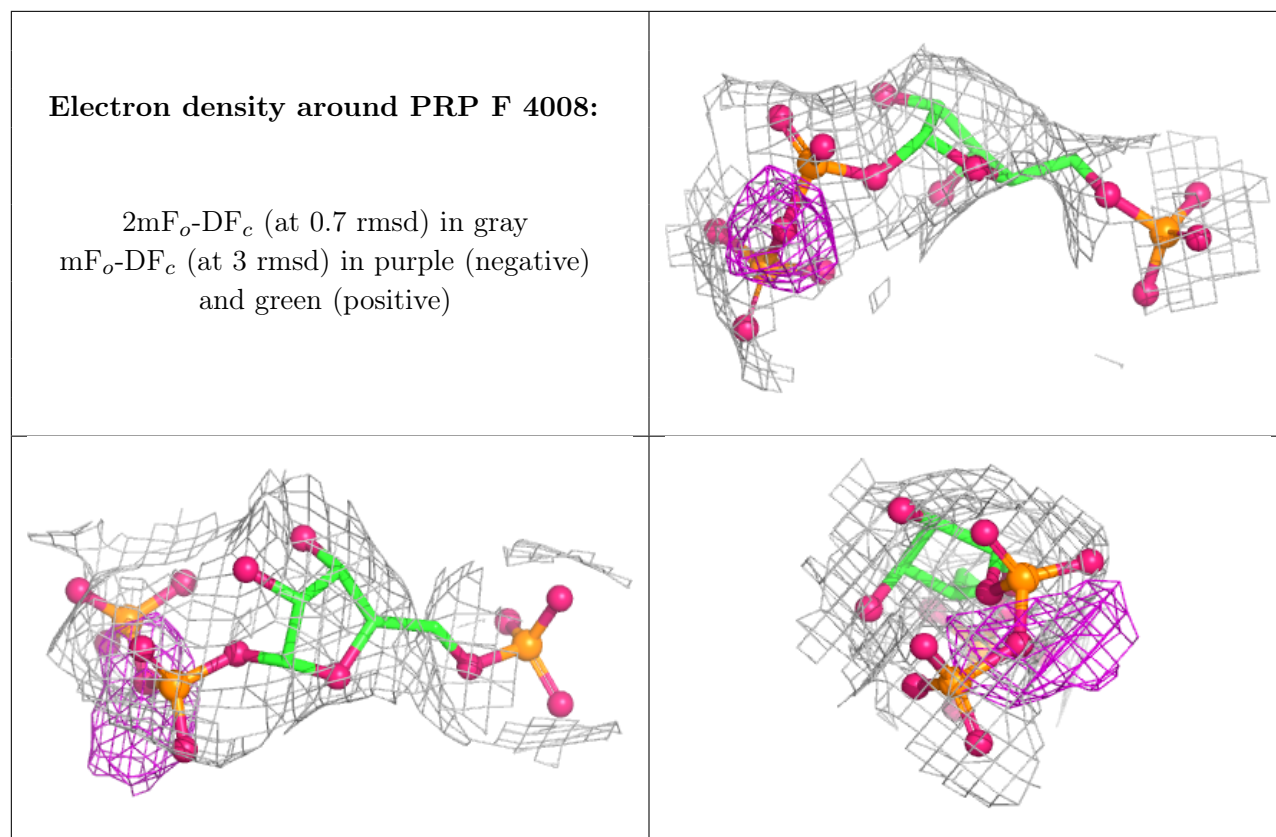
Electron density around PRP G 4009:

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

**Electron density around PRP E 4007:**

$2mF_o-DF_c$ (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.