



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 04:07 AM UTC

PDB ID : 3E1F / pdb_00003e1f
Title : E.Coli (lacZ) beta-galactosidase (H418E) in complex with galactose
Authors : Huber, R.E.; Dugdale, M.L.
Deposited on : 2008-08-04
Resolution : 3.00 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
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

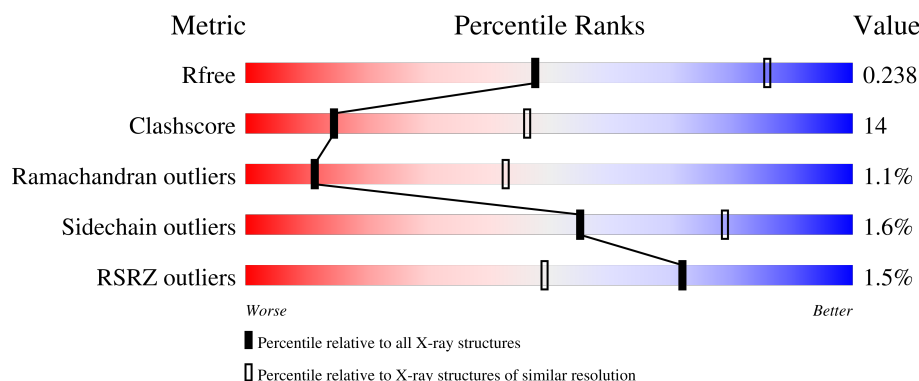
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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	1	1023	2% 66% 30% ..
1	2	1023	3% 67% 30% ..
1	3	1023	% 67% 30% ..
1	4	1023	% 68% 29% ..

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	DMS	1	8502	-	-	X	-
4	DMS	2	8001	-	-	X	-
4	DMS	3	8415	-	-	-	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 33235 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 Beta-galactosidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	1	1011	Total	C	N	O	S	0	0	0
			8124	5137	1438	1511	38			
1	2	1011	Total	C	N	O	S	0	0	0
			8124	5137	1438	1511	38			
1	3	1011	Total	C	N	O	S	0	0	0
			8124	5137	1438	1511	38			
1	4	1011	Total	C	N	O	S	0	0	0
			8124	5137	1438	1511	38			

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
1	1	GLY	-	expression tag	UNP P00722
1	2	SER	-	expression tag	UNP P00722
1	3	HIS	-	expression tag	UNP P00722
1	4	MET	-	expression tag	UNP P00722
1	5	LEU	-	expression tag	UNP P00722
1	6	GLU	-	expression tag	UNP P00722
1	7	ASP	-	expression tag	UNP P00722
1	8	PRO	-	expression tag	UNP P00722
1	418	GLU	HIS	engineered mutation	UNP P00722
2	1	GLY	-	expression tag	UNP P00722
2	2	SER	-	expression tag	UNP P00722
2	3	HIS	-	expression tag	UNP P00722
2	4	MET	-	expression tag	UNP P00722
2	5	LEU	-	expression tag	UNP P00722
2	6	GLU	-	expression tag	UNP P00722
2	7	ASP	-	expression tag	UNP P00722
2	8	PRO	-	expression tag	UNP P00722
2	418	GLU	HIS	engineered mutation	UNP P00722
3	1	GLY	-	expression tag	UNP P00722
3	2	SER	-	expression tag	UNP P00722
3	3	HIS	-	expression tag	UNP P00722

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Chain	Residue	Modelled	Actual	Comment	Reference
3	4	MET	-	expression tag	UNP P00722
3	5	LEU	-	expression tag	UNP P00722
3	6	GLU	-	expression tag	UNP P00722
3	7	ASP	-	expression tag	UNP P00722
3	8	PRO	-	expression tag	UNP P00722
3	418	GLU	HIS	engineered mutation	UNP P00722
4	1	GLY	-	expression tag	UNP P00722
4	2	SER	-	expression tag	UNP P00722
4	3	HIS	-	expression tag	UNP P00722
4	4	MET	-	expression tag	UNP P00722
4	5	LEU	-	expression tag	UNP P00722
4	6	GLU	-	expression tag	UNP P00722
4	7	ASP	-	expression tag	UNP P00722
4	8	PRO	-	expression tag	UNP P00722
4	418	GLU	HIS	engineered mutation	UNP P00722

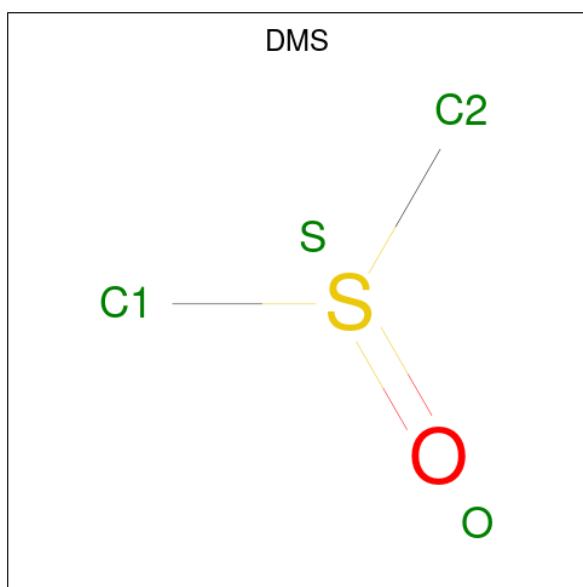
- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	1	3	Total Mg 3 3	0	0
2	2	3	Total Mg 3 3	0	0
2	3	2	Total Mg 2 2	0	0
2	4	2	Total Mg 2 2	0	0

- Molecule 3 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	1	3	Total Na 3 3	0	0
3	2	4	Total Na 4 4	0	0
3	3	4	Total Na 4 4	0	0
3	4	2	Total Na 2 2	0	0

- Molecule 4 is DIMETHYL SULFOXIDE (CCD ID: DMS) (formula: C₂H₆OS).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	1	1	Total	C	O	S	0	0
			4	2	1	1		
4	1	1	Total	C	O	S	0	0
			4	2	1	1		
4	1	1	Total	C	O	S	0	0
			4	2	1	1		
4	1	1	Total	C	O	S	0	0
			4	2	1	1		
4	1	1	Total	C	O	S	0	0
			4	2	1	1		
4	1	1	Total	C	O	S	0	0
			4	2	1	1		
4	1	1	Total	C	O	S	0	0
			4	2	1	1		
4	1	1	Total	C	O	S	0	0
			4	2	1	1		
4	1	1	Total	C	O	S	0	0
			4	2	1	1		
4	1	1	Total	C	O	S	0	0
			4	2	1	1		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	1	1	Total 4	C 2	O 1	S 1	0	0
4	1	1	Total 4	C 2	O 1	S 1	0	0
4	1	1	Total 4	C 2	O 1	S 1	0	0
4	1	1	Total 4	C 2	O 1	S 1	0	0
4	1	1	Total 4	C 2	O 1	S 1	0	0
4	1	1	Total 4	C 2	O 1	S 1	0	0
4	1	1	Total 4	C 2	O 1	S 1	0	0
4	1	1	Total 4	C 2	O 1	S 1	0	0
4	1	1	Total 4	C 2	O 1	S 1	0	0
4	1	1	Total 4	C 2	O 1	S 1	0	0
4	2	1	Total 4	C 2	O 1	S 1	0	0
4	2	1	Total 4	C 2	O 1	S 1	0	0
4	2	1	Total 4	C 2	O 1	S 1	0	0
4	2	1	Total 4	C 2	O 1	S 1	0	0
4	2	1	Total 4	C 2	O 1	S 1	0	0
4	2	1	Total 4	C 2	O 1	S 1	0	0
4	2	1	Total 4	C 2	O 1	S 1	0	0
4	2	1	Total 4	C 2	O 1	S 1	0	0
4	2	1	Total 4	C 2	O 1	S 1	0	0
4	2	1	Total 4	C 2	O 1	S 1	0	0

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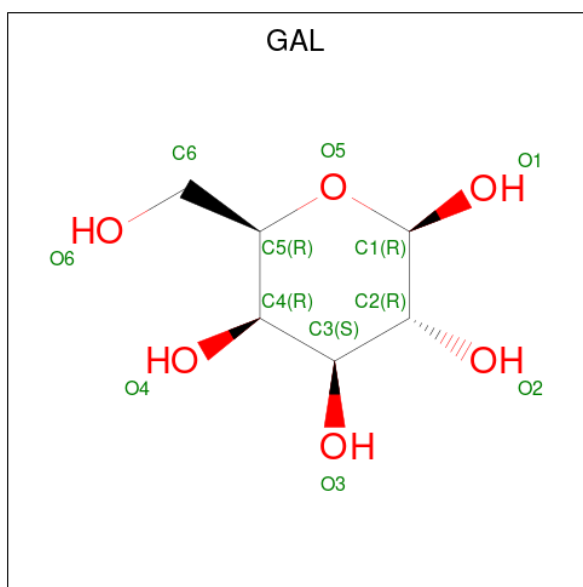
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	2	1	Total 4	C 2	O 1	S 1	0	0
4	2	1	Total 4	C 2	O 1	S 1	0	0
4	2	1	Total 4	C 2	O 1	S 1	0	0
4	2	1	Total 4	C 2	O 1	S 1	0	0
4	2	1	Total 4	C 2	O 1	S 1	0	0
4	2	1	Total 4	C 2	O 1	S 1	0	0
4	2	1	Total 4	C 2	O 1	S 1	0	0
4	2	1	Total 4	C 2	O 1	S 1	0	0
4	2	1	Total 4	C 2	O 1	S 1	0	0
4	3	1	Total 4	C 2	O 1	S 1	0	0
4	3	1	Total 4	C 2	O 1	S 1	0	0
4	3	1	Total 4	C 2	O 1	S 1	0	0
4	3	1	Total 4	C 2	O 1	S 1	0	0
4	3	1	Total 4	C 2	O 1	S 1	0	0
4	3	1	Total 4	C 2	O 1	S 1	0	0
4	3	1	Total 4	C 2	O 1	S 1	0	0
4	3	1	Total 4	C 2	O 1	S 1	0	0
4	3	1	Total 4	C 2	O 1	S 1	0	0
4	3	1	Total 4	C 2	O 1	S 1	0	0
4	3	1	Total 4	C 2	O 1	S 1	0	0

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	3	1	Total 4	C 2	O 1	S 1	0	0
4	3	1	Total 4	C 2	O 1	S 1	0	0
4	3	1	Total 4	C 2	O 1	S 1	0	0
4	3	1	Total 4	C 2	O 1	S 1	0	0
4	3	1	Total 4	C 2	O 1	S 1	0	0
4	3	1	Total 4	C 2	O 1	S 1	0	0
4	3	1	Total 4	C 2	O 1	S 1	0	0
4	3	1	Total 4	C 2	O 1	S 1	0	0
4	3	1	Total 4	C 2	O 1	S 1	0	0
4	3	1	Total 4	C 2	O 1	S 1	0	0
4	3	1	Total 4	C 2	O 1	S 1	0	0
4	4	1	Total 4	C 2	O 1	S 1	0	0
4	4	1	Total 4	C 2	O 1	S 1	0	0
4	4	1	Total 4	C 2	O 1	S 1	0	0
4	4	1	Total 4	C 2	O 1	S 1	0	0
4	4	1	Total 4	C 2	O 1	S 1	0	0

- Molecule 5 is beta-D-galactopyranose (CCD ID: GAL) (formula: C₆H₁₂O₆).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	1	1	Total C O 12 6 6	0	0
5	1	1	Total C O 12 6 6	0	0
5	2	1	Total C O 12 6 6	0	0
5	2	1	Total C O 12 6 6	0	0
5	3	1	Total C O 12 6 6	0	0
5	3	1	Total C O 12 6 6	0	0
5	4	1	Total C O 12 6 6	0	0
5	4	1	Total C O 12 6 6	0	0

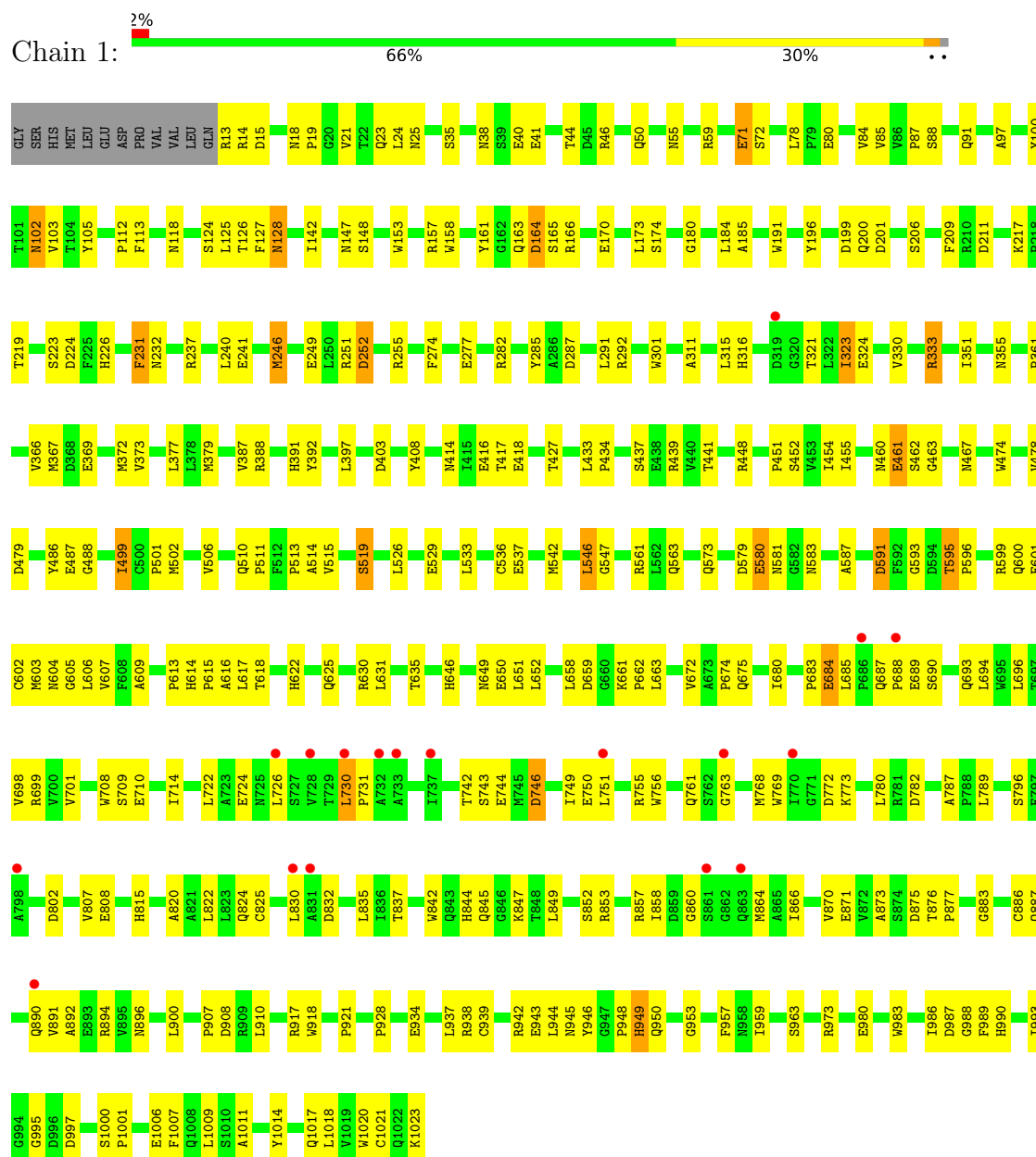
- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	1	56	Total O 56 56	0	0
6	2	59	Total O 59 59	0	0
6	3	109	Total O 109 109	0	0
6	4	104	Total O 104 104	0	0

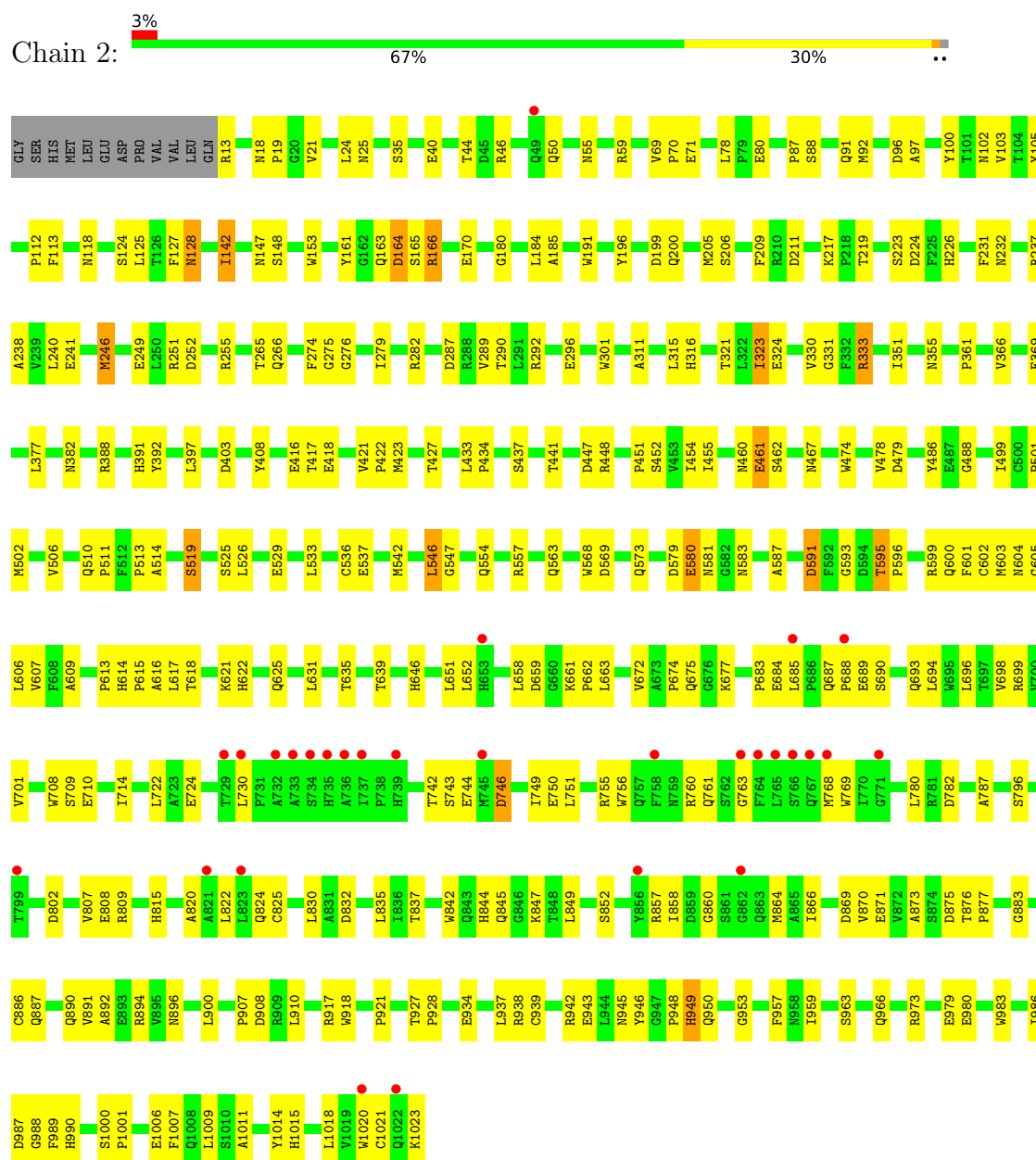
3 Residue-property plots [i](#)

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

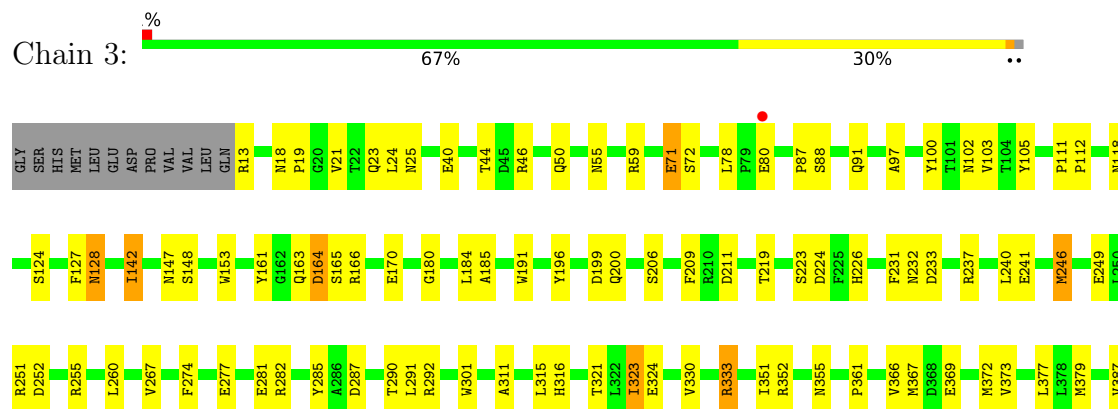
- Molecule 1: Beta-galactosidase

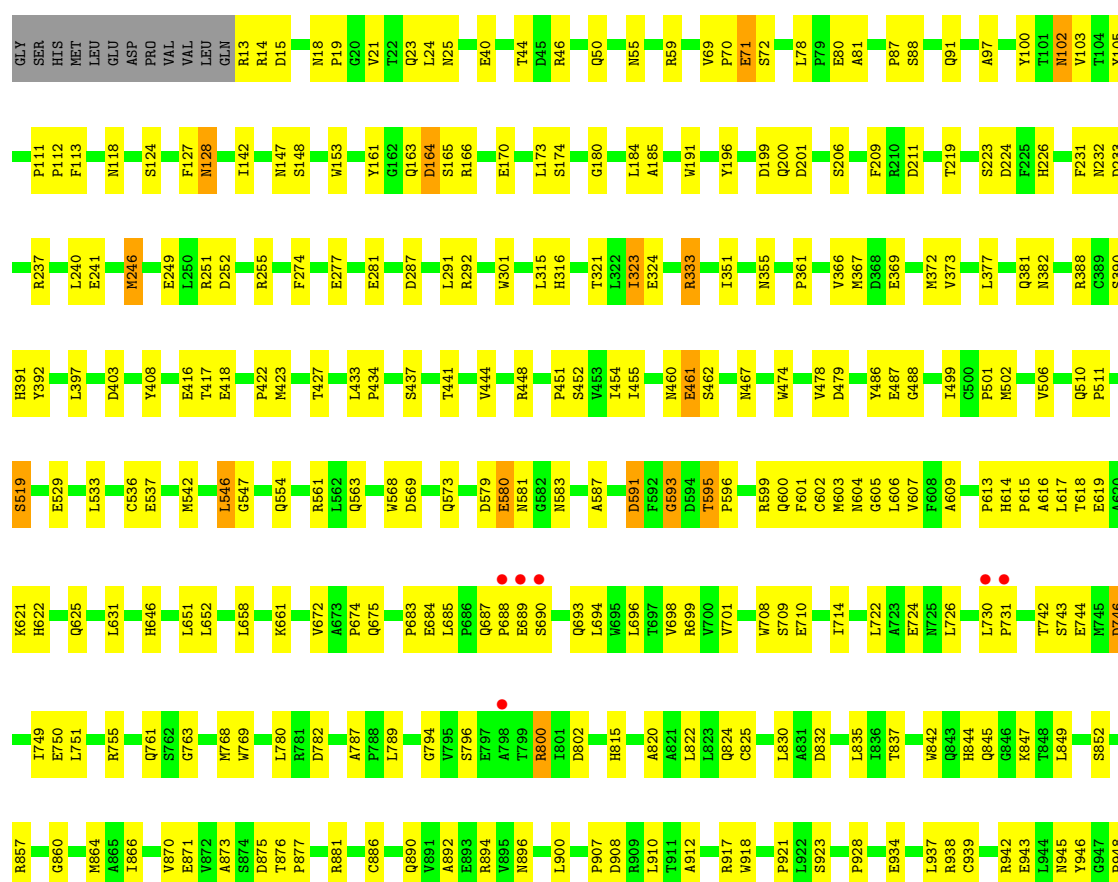


- Molecule 1: Beta-galactosidase



- Molecule 1: Beta-galactosidase





H949	Q950	G953	F957	S960	S963	Q966	R973	E979	Y983	T986	D987	G988	F989	H990	T993	G994	G995	D996	D997	S1000	P1001	E1006	F1007	A1011	Y1014	Q1017	L1018	V1019	W1020	C1021	Q1022	K1023
------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	-------	-------	-------	-------	-------	-------	-------	-------	-------	-------	-------	-------	-------

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	128.37Å 152.87Å 132.12Å 90.00° 102.68° 90.00°	Depositor
Resolution (Å)	98.54 – 3.00 98.54 – 3.00	Depositor EDS
% Data completeness (in resolution range)	91.1 (98.54-3.00) 91.1 (98.54-3.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.81 (at 3.01Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.218 , 0.249 0.208 , 0.238	Depositor DCC
R_{free} test set	4529 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	43.5	Xtriage
Anisotropy	0.423	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 52.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.022 for l,-k,h	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	33235	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.63% 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: DMS, GAL, MG, NA

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	1	0.40	0/8365	0.94	27/11412 (0.2%)
1	2	0.43	0/8365	0.94	30/11412 (0.3%)
1	3	0.44	0/8365	0.93	24/11412 (0.2%)
1	4	0.45	0/8365	0.95	28/11412 (0.2%)
All	All	0.43	0/33460	0.94	109/45648 (0.2%)

There are no bond length outliers.

All (109) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	1	157	ARG	NE-CZ-NH2	12.12	130.11	119.20
1	4	800	ARG	NE-CZ-NH2	11.93	129.93	119.20
1	2	809	ARG	NE-CZ-NH2	11.60	129.64	119.20
1	1	157	ARG	NE-CZ-NH1	-11.20	110.30	121.50
1	4	800	ARG	NE-CZ-NH1	-10.99	110.51	121.50
1	2	809	ARG	NE-CZ-NH1	-10.53	110.97	121.50
1	4	800	ARG	CD-NE-CZ	9.06	137.08	124.40
1	1	157	ARG	CD-NE-CZ	9.05	137.07	124.40
1	2	809	ARG	CD-NE-CZ	7.92	135.49	124.40
1	3	744	GLU	N-CA-C	-7.84	103.85	113.50
1	4	744	GLU	N-CA-C	-7.43	104.36	113.50
1	2	744	GLU	N-CA-C	-7.36	104.33	113.38
1	2	614	HIS	N-CA-C	-7.36	101.48	110.31
1	1	744	GLU	N-CA-C	-7.29	104.54	113.50
1	1	614	HIS	N-CA-C	-7.11	101.78	110.31
1	4	614	HIS	N-CA-C	-7.04	101.86	110.31
1	3	614	HIS	N-CA-C	-6.74	102.22	110.31
1	2	97	ALA	N-CA-C	6.65	118.18	109.93
1	1	506	VAL	N-CA-C	6.64	116.77	110.53
1	3	97	ALA	N-CA-C	6.64	118.16	109.93

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	2	506	VAL	N-CA-C	6.62	116.76	110.53
1	1	563	GLN	N-CA-C	6.58	120.30	112.93
1	3	323	ILE	N-CA-C	-6.56	106.17	111.81
1	4	97	ALA	N-CA-C	6.47	117.95	109.93
1	1	763	GLY	N-CA-C	-6.44	106.03	115.63
1	3	563	GLN	N-CA-C	6.44	120.14	112.93
1	4	323	ILE	N-CA-C	-6.42	106.29	111.81
1	2	763	GLY	N-CA-C	-6.38	106.12	115.63
1	1	97	ALA	N-CA-C	6.34	117.79	109.93
1	4	763	GLY	N-CA-C	-6.33	106.19	115.63
1	3	506	VAL	N-CA-C	6.28	116.44	110.53
1	4	506	VAL	N-CA-C	6.20	116.35	110.53
1	4	742	THR	N-CA-C	6.18	118.47	108.34
1	1	323	ILE	N-CA-C	-6.17	106.51	111.81
1	2	323	ILE	N-CA-C	-6.12	106.55	111.81
1	3	606	LEU	N-CA-C	-6.11	105.83	113.28
1	3	479	ASP	CA-C-N	6.08	125.70	119.56
1	3	479	ASP	C-N-CA	6.08	125.70	119.56
1	3	742	THR	N-CA-C	6.06	118.28	108.34
1	2	606	LEU	N-CA-C	-6.02	105.93	113.28
1	3	763	GLY	N-CA-C	-5.94	106.78	115.63
1	2	479	ASP	CA-C-N	5.92	125.54	119.56
1	2	479	ASP	C-N-CA	5.92	125.54	119.56
1	4	563	GLN	N-CA-C	5.91	119.54	112.93
1	2	554	GLN	N-CA-C	-5.89	104.76	111.07
1	1	949	HIS	N-CA-C	5.86	118.46	110.55
1	4	949	HIS	N-CA-C	5.83	118.42	110.55
1	1	742	THR	N-CA-C	5.82	117.88	108.34
1	1	606	LEU	N-CA-C	-5.80	106.20	113.28
1	2	742	THR	N-CA-C	5.79	117.83	108.34
1	2	949	HIS	N-CA-C	5.77	118.34	110.55
1	1	351	ILE	N-CA-C	5.77	115.85	108.12
1	1	730	LEU	CA-C-N	5.74	127.01	119.84
1	1	730	LEU	C-N-CA	5.74	127.01	119.84
1	3	730	LEU	CA-C-N	5.71	126.98	119.84
1	3	730	LEU	C-N-CA	5.71	126.98	119.84
1	2	563	GLN	N-CA-C	5.69	119.30	112.93
1	3	949	HIS	N-CA-C	5.66	118.19	110.55
1	2	730	LEU	CA-C-N	5.64	126.89	119.84
1	2	730	LEU	C-N-CA	5.64	126.89	119.84
1	4	479	ASP	CA-C-N	5.64	125.25	119.56
1	4	479	ASP	C-N-CA	5.64	125.25	119.56

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	4	554	GLN	N-CA-C	-5.61	105.07	111.07
1	4	730	LEU	CA-C-N	5.59	126.83	119.84
1	4	730	LEU	C-N-CA	5.59	126.83	119.84
1	3	351	ILE	N-CA-C	5.58	115.60	108.12
1	4	606	LEU	N-CA-C	-5.57	106.49	113.28
1	3	554	GLN	N-CA-C	-5.54	104.46	111.11
1	4	510	GLN	CA-C-N	5.52	126.74	119.84
1	4	510	GLN	C-N-CA	5.52	126.74	119.84
1	3	510	GLN	CA-C-N	5.51	126.73	119.84
1	3	510	GLN	C-N-CA	5.51	126.73	119.84
1	4	351	ILE	N-CA-C	5.50	115.50	108.12
1	1	510	GLN	CA-C-N	5.50	126.72	119.84
1	1	510	GLN	C-N-CA	5.50	126.72	119.84
1	4	646	HIS	N-CA-C	-5.48	101.64	110.14
1	2	519	SER	N-CA-C	-5.48	101.94	110.10
1	1	519	SER	N-CA-C	-5.46	101.96	110.10
1	4	605	GLY	N-CA-C	5.43	121.60	112.85
1	1	605	GLY	N-CA-C	5.41	121.56	112.85
1	3	142	ILE	N-CA-C	5.39	115.62	107.75
1	1	479	ASP	CA-C-N	5.38	125.00	119.56
1	1	479	ASP	C-N-CA	5.38	125.00	119.56
1	2	605	GLY	N-CA-C	5.38	121.51	112.85
1	3	746	ASP	N-CA-C	5.38	117.24	109.07
1	1	646	HIS	N-CA-C	-5.37	101.83	110.14
1	3	91	GLN	N-CA-C	-5.34	105.54	111.36
1	3	519	SER	N-CA-C	-5.34	102.15	110.10
1	1	746	ASP	N-CA-C	5.33	117.18	109.07
1	2	447	ASP	N-CA-C	5.30	119.74	113.28
1	1	91	GLN	N-CA-C	-5.22	105.67	111.36
1	4	519	SER	N-CA-C	-5.20	102.35	110.10
1	2	351	ILE	N-CA-C	5.19	115.07	108.12
1	4	746	ASP	N-CA-C	5.17	116.92	109.07
1	2	746	ASP	N-CA-C	5.14	116.88	109.07
1	4	113	PHE	N-CA-C	5.12	118.29	110.20
1	1	113	PHE	N-CA-C	5.10	118.26	110.20
1	2	91	GLN	N-CA-C	-5.09	105.82	111.36
1	2	510	GLN	CA-C-N	5.07	126.18	119.84
1	2	510	GLN	C-N-CA	5.07	126.18	119.84
1	3	646	HIS	N-CA-C	-5.07	102.28	110.14
1	2	646	HIS	N-CA-C	-5.06	102.29	110.14
1	1	499	ILE	N-CA-C	-5.06	100.34	107.98
1	2	113	PHE	N-CA-C	5.05	118.18	110.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	3	605	GLY	N-CA-C	5.05	120.87	112.89
1	4	91	GLN	N-CA-C	-5.04	105.87	111.36
1	4	593	GLY	N-CA-C	-5.03	108.67	115.36
1	2	142	ILE	N-CA-C	5.03	115.69	107.24
1	2	166	ARG	N-CA-C	5.00	120.00	113.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	1	8124	0	7714	248	0
1	2	8124	0	7715	233	0
1	3	8124	0	7714	236	0
1	4	8124	0	7714	220	0
2	1	3	0	0	0	0
2	2	3	0	0	0	0
2	3	2	0	0	0	0
2	4	2	0	0	0	0
3	1	3	0	0	0	0
3	2	4	0	0	0	0
3	3	4	0	0	0	0
3	4	2	0	0	0	0
4	1	96	0	144	10	0
4	2	80	0	120	11	0
4	3	92	0	138	11	0
4	4	24	0	36	2	0
5	1	24	0	24	1	0
5	2	24	0	24	1	0
5	3	24	0	23	0	0
5	4	24	0	24	0	0
6	1	56	0	0	2	0
6	2	59	0	0	0	0
6	3	109	0	0	2	0
6	4	104	0	0	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	33235	0	31390	898	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 14.

All (898) 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)
1:2:276:GLY:HA2	4:2:8001:DMS:H11	1.41	0.99
1:1:84:VAL:HG12	4:1:8414:DMS:H22	1.44	0.97
1:2:142:ILE:HG12	1:2:170:GLU:HG2	1.53	0.91
1:1:652:LEU:HD11	1:1:698:VAL:HB	1.51	0.90
1:2:289:VAL:HG23	4:2:8001:DMS:H23	1.52	0.90
1:3:142:ILE:HG12	1:3:170:GLU:HG2	1.52	0.90
1:3:652:LEU:HD11	1:3:698:VAL:HB	1.52	0.90
1:1:142:ILE:HG12	1:1:170:GLU:HG2	1.54	0.90
1:4:652:LEU:HD11	1:4:698:VAL:HB	1.53	0.89
1:2:652:LEU:HD11	1:2:698:VAL:HB	1.52	0.88
1:4:142:ILE:HG12	1:4:170:GLU:HG2	1.54	0.86
1:2:103:VAL:HG13	1:2:418:GLU:HG2	1.58	0.84
1:2:377:LEU:HD22	1:2:708:TRP:HA	1.60	0.84
1:4:103:VAL:HG13	1:4:418:GLU:HG2	1.60	0.83
1:3:377:LEU:HD22	1:3:708:TRP:HA	1.60	0.82
1:4:377:LEU:HD22	1:4:708:TRP:HA	1.59	0.82
1:1:103:VAL:HG13	1:1:418:GLU:HG2	1.61	0.81
1:3:103:VAL:HG13	1:3:418:GLU:HG2	1.61	0.80
1:1:15:ASP:CG	1:4:13:ARG:HH22	1.89	0.80
1:4:615:PRO:O	1:4:618:THR:HG22	1.83	0.79
1:1:615:PRO:O	1:1:618:THR:HG22	1.83	0.79
1:1:377:LEU:HD22	1:1:708:TRP:HA	1.64	0.78
1:1:13:ARG:HH22	1:4:15:ASP:CG	1.93	0.76
1:2:615:PRO:O	1:2:618:THR:HG22	1.86	0.76
1:2:687:GLN:OE1	1:2:688:PRO:HD2	1.87	0.75
1:3:615:PRO:O	1:3:618:THR:HG22	1.87	0.74
1:1:599:ARG:HB3	1:1:600:GLN:OE1	1.87	0.74
1:3:599:ARG:HB3	1:3:600:GLN:OE1	1.87	0.74
1:4:147:ASN:HB3	1:4:206:SER:HA	1.71	0.73
1:1:687:GLN:OE1	1:1:688:PRO:HD2	1.89	0.73
1:4:599:ARG:HB3	1:4:600:GLN:OE1	1.89	0.73
1:1:13:ARG:HD3	1:4:13:ARG:HD3	1.71	0.72
1:4:699:ARG:HD3	1:4:714:ILE:HD13	1.70	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:3:687:GLN:OE1	1:3:688:PRO:HD2	1.89	0.72
1:4:687:GLN:OE1	1:4:688:PRO:HD2	1.88	0.72
1:2:147:ASN:HB3	1:2:206:SER:HA	1.71	0.72
1:2:699:ARG:HD3	1:2:714:ILE:HD13	1.72	0.71
1:3:147:ASN:HB3	1:3:206:SER:HA	1.71	0.71
1:1:699:ARG:HD3	1:1:714:ILE:HD13	1.71	0.71
1:1:125:LEU:HA	4:1:8502:DMS:H13	1.72	0.71
1:2:599:ARG:HB3	1:2:600:GLN:OE1	1.89	0.71
1:4:945:ASN:HB3	1:4:1023:LYS:HE3	1.72	0.70
1:1:542:MET:CE	1:1:601:PHE:HA	2.21	0.70
1:2:542:MET:CE	1:2:601:PHE:HA	2.22	0.70
1:4:153:TRP:HB2	1:4:185:ALA:HB3	1.74	0.70
1:1:147:ASN:HB3	1:1:206:SER:HA	1.73	0.70
1:4:355:ASN:OD1	1:4:388:ARG:HD3	1.91	0.70
1:3:699:ARG:HD3	1:3:714:ILE:HD13	1.74	0.70
1:1:945:ASN:HB3	1:1:1023:LYS:HE3	1.74	0.69
1:1:153:TRP:HB2	1:1:185:ALA:HB3	1.74	0.69
1:2:153:TRP:HB2	1:2:185:ALA:HB3	1.73	0.69
1:3:542:MET:CE	1:3:601:PHE:HA	2.22	0.69
1:2:945:ASN:HB3	1:2:1023:LYS:HE3	1.74	0.69
1:3:622:HIS:O	1:3:625:GLN:HG3	1.92	0.69
1:2:249:GLU:HG2	1:2:251:ARG:NH2	2.08	0.68
1:3:153:TRP:HB2	1:3:185:ALA:HB3	1.74	0.68
1:1:355:ASN:OD1	1:1:388:ARG:HD3	1.94	0.68
1:3:355:ASN:OD1	1:3:388:ARG:HD3	1.93	0.68
1:3:249:GLU:HG2	1:3:251:ARG:NH2	2.08	0.68
1:4:542:MET:CE	1:4:601:PHE:HA	2.23	0.68
1:3:55:ASN:ND2	1:3:87:PRO:HD3	2.08	0.68
1:3:361:PRO:HB3	1:3:609:ALA:HB1	1.76	0.68
1:2:241:GLU:HG3	1:2:292:ARG:HG2	1.77	0.67
1:2:361:PRO:HB3	1:2:609:ALA:HB1	1.77	0.67
1:1:249:GLU:HG2	1:1:251:ARG:NH2	2.09	0.67
1:1:361:PRO:HB3	1:1:609:ALA:HB1	1.77	0.67
1:4:622:HIS:O	1:4:625:GLN:HG3	1.95	0.67
1:1:622:HIS:O	1:1:625:GLN:HG3	1.95	0.66
1:4:249:GLU:HG2	1:4:251:ARG:NH2	2.10	0.66
1:3:945:ASN:HB3	1:3:1023:LYS:HE3	1.77	0.66
1:2:658:LEU:O	1:2:661:LYS:HG3	1.96	0.66
1:1:986:ILE:HG21	1:1:1018:LEU:HD11	1.77	0.65
1:4:55:ASN:ND2	1:4:87:PRO:HD3	2.10	0.65
1:2:986:ILE:HG21	1:2:1018:LEU:HD11	1.77	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:55:ASN:ND2	1:2:87:PRO:HD3	2.09	0.65
1:1:125:LEU:HD11	4:1:8408:DMS:H23	1.78	0.65
1:1:241:GLU:HG3	1:1:292:ARG:HG2	1.78	0.65
1:3:290:THR:H	4:3:8001:DMS:H22	1.59	0.65
1:4:166:ARG:HG3	1:4:392:TYR:HB2	1.78	0.65
1:4:361:PRO:HB3	1:4:609:ALA:HB1	1.77	0.65
1:1:166:ARG:HG3	1:1:392:TYR:HB2	1.79	0.64
1:3:573:GLN:HB2	1:3:602:CYS:O	1.97	0.64
1:2:166:ARG:HG3	1:2:392:TYR:HB2	1.80	0.64
1:3:580:GLU:H	1:3:580:GLU:CD	2.06	0.64
1:4:923:SER:HA	4:4:8006:DMS:H12	1.80	0.64
1:2:355:ASN:OD1	1:2:388:ARG:HD3	1.96	0.64
1:4:928:PRO:HB2	1:4:973:ARG:HH11	1.63	0.64
1:4:986:ILE:HG21	1:4:1018:LEU:HD11	1.80	0.64
1:3:986:ILE:HG21	1:3:1018:LEU:HD11	1.80	0.64
1:1:928:PRO:HB2	1:1:973:ARG:HH11	1.62	0.63
1:4:658:LEU:O	1:4:661:LYS:HG3	1.98	0.63
1:1:55:ASN:ND2	1:1:87:PRO:HD3	2.13	0.63
1:3:546:LEU:HA	6:3:8807:HOH:O	1.97	0.63
1:2:945:ASN:OD1	1:2:950:GLN:HG3	1.98	0.63
1:1:580:GLU:CD	1:1:580:GLU:H	2.06	0.63
1:4:580:GLU:CD	1:4:580:GLU:H	2.07	0.63
1:2:782:ASP:HB2	1:2:842:TRP:CH2	2.34	0.63
1:3:557:ARG:NH1	4:3:8402:DMS:H13	2.14	0.62
1:3:658:LEU:O	1:3:661:LYS:HG3	1.99	0.62
1:2:580:GLU:CD	1:2:580:GLU:H	2.06	0.62
1:4:945:ASN:OD1	1:4:950:GLN:HG3	2.00	0.62
1:1:782:ASP:HB2	1:1:842:TRP:CH2	2.35	0.62
1:1:292:ARG:HH12	4:1:8001:DMS:H22	1.64	0.62
1:2:88:SER:HA	1:2:366:VAL:HG21	1.82	0.62
1:3:241:GLU:HG3	1:3:292:ARG:HG2	1.80	0.62
1:2:894:ARG:NH1	1:2:921:PRO:HD3	2.15	0.62
1:3:928:PRO:HB2	1:3:973:ARG:HH11	1.65	0.62
1:3:166:ARG:HG3	1:3:392:TYR:HB2	1.81	0.61
1:1:945:ASN:OD1	1:1:950:GLN:HG3	2.00	0.61
1:2:622:HIS:O	1:2:625:GLN:HG3	1.99	0.61
1:1:163:GLN:O	1:1:164:ASP:HB3	2.00	0.61
1:2:275:GLY:O	4:2:8001:DMS:H21	2.00	0.61
1:2:973:ARG:NH2	4:2:8411:DMS:H13	2.15	0.61
1:4:241:GLU:HG3	1:4:292:ARG:HG2	1.80	0.61
1:1:454:ILE:HG13	1:1:455:ILE:HG13	1.83	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:4:163:GLN:O	1:4:164:ASP:HB3	2.00	0.61
1:1:282:ARG:HG2	1:4:423:MET:HE2	1.81	0.61
1:1:658:LEU:O	1:1:661:LYS:HG3	2.00	0.61
1:4:232:ASN:ND2	1:4:237:ARG:HB2	2.16	0.61
1:1:59:ARG:HG2	4:1:8502:DMS:H12	1.83	0.60
1:1:433:LEU:HD23	1:4:437:SER:OG	1.99	0.60
1:1:894:ARG:NH1	1:1:921:PRO:HD3	2.15	0.60
1:2:163:GLN:O	1:2:164:ASP:HB3	2.01	0.60
1:2:928:PRO:HB2	1:2:973:ARG:HH11	1.65	0.60
1:4:200:GLN:HG2	1:4:391:HIS:HB2	1.83	0.60
1:2:290:THR:H	4:2:8001:DMS:H12	1.66	0.60
1:3:782:ASP:HB2	1:3:842:TRP:CH2	2.35	0.60
1:2:573:GLN:HB2	1:2:602:CYS:O	2.01	0.60
1:1:15:ASP:OD1	1:4:13:ARG:NH2	2.34	0.60
1:1:573:GLN:HB2	1:1:602:CYS:O	2.01	0.60
1:4:894:ARG:NH1	1:4:921:PRO:HD3	2.15	0.60
1:1:88:SER:HA	1:1:366:VAL:HG21	1.83	0.60
1:3:945:ASN:OD1	1:3:950:GLN:HG3	2.01	0.60
1:4:547:GLY:HA3	1:4:908:ASP:OD2	2.01	0.60
1:3:892:ALA:HB3	1:3:946:TYR:CE1	2.35	0.60
1:1:437:SER:OG	1:4:433:LEU:HD23	2.02	0.60
1:4:782:ASP:HB2	1:4:842:TRP:CH2	2.36	0.59
1:4:892:ALA:HB3	1:4:946:TYR:CE1	2.36	0.59
1:1:542:MET:HE1	1:1:601:PHE:HA	1.84	0.59
1:1:830:LEU:HD21	1:2:830:LEU:HD21	1.85	0.59
1:2:331:GLY:O	4:2:8401:DMS:H13	2.01	0.59
1:3:88:SER:HA	1:3:366:VAL:HG21	1.82	0.59
1:3:542:MET:HE1	1:3:601:PHE:HA	1.85	0.59
1:1:285:TYR:CE1	1:4:422:PRO:HG3	2.37	0.58
1:3:557:ARG:CZ	4:3:8402:DMS:H13	2.33	0.58
1:2:780:LEU:HA	1:2:886:CYS:HB3	1.86	0.58
1:1:200:GLN:HG2	1:1:391:HIS:HB2	1.85	0.58
1:1:13:ARG:NH2	1:4:15:ASP:OD1	2.36	0.58
1:2:100:TYR:CE1	1:2:602:CYS:HB3	2.38	0.58
1:3:894:ARG:NH1	1:3:921:PRO:HD3	2.19	0.58
1:4:88:SER:HA	1:4:366:VAL:HG21	1.83	0.58
1:1:85:VAL:HG23	4:1:8414:DMS:O	2.03	0.58
1:1:547:GLY:HA3	1:1:908:ASP:OD2	2.04	0.57
1:3:163:GLN:O	1:3:164:ASP:HB3	2.04	0.57
1:2:547:GLY:HA3	1:2:908:ASP:OD2	2.04	0.57
1:3:693:GLN:OE1	1:3:724:GLU:HB2	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:651:LEU:HD21	1:2:701:VAL:HB	1.87	0.57
1:3:200:GLN:HG2	1:3:391:HIS:HB2	1.86	0.57
1:1:38:ASN:HA	6:1:8613:HOH:O	2.04	0.57
1:2:892:ALA:HB3	1:2:946:TYR:CE1	2.40	0.57
1:3:750:GLU:HG2	1:3:755:ARG:HD3	1.87	0.57
1:3:369:GLU:HG3	1:3:397:LEU:HD21	1.86	0.57
1:4:573:GLN:HB2	1:4:602:CYS:O	2.05	0.57
1:2:246:MET:HE3	1:2:246:MET:C	2.30	0.57
1:2:542:MET:HE1	1:2:601:PHE:HA	1.86	0.56
1:1:892:ALA:HB3	1:1:946:TYR:CE1	2.39	0.56
1:2:125:LEU:HD11	4:2:8000:DMS:H23	1.87	0.56
1:3:547:GLY:HA3	1:3:908:ASP:OD2	2.04	0.56
1:4:454:ILE:HG13	1:4:455:ILE:HG13	1.85	0.56
1:4:750:GLU:HG2	1:4:755:ARG:HD3	1.86	0.56
1:1:44:THR:OG1	1:1:46:ARG:HG3	2.06	0.56
1:1:613:PRO:HB3	1:1:617:LEU:HD23	1.87	0.56
1:1:630:ARG:HA	4:1:8503:DMS:S	2.45	0.56
1:3:59:ARG:HB2	1:3:124:SER:OG	2.05	0.56
1:3:830:LEU:HD21	1:4:830:LEU:HD21	1.88	0.56
1:2:200:GLN:HG2	1:2:391:HIS:HB2	1.86	0.56
1:1:651:LEU:HD21	1:1:701:VAL:HB	1.88	0.56
1:1:780:LEU:HA	1:1:886:CYS:HB3	1.88	0.56
1:2:282:ARG:O	1:3:421:VAL:HG13	2.05	0.56
1:4:595:THR:HA	1:4:596:PRO:C	2.31	0.56
1:3:100:TYR:CE1	1:3:602:CYS:HB3	2.41	0.56
1:4:315:LEU:O	1:4:323:ILE:HB	2.06	0.56
1:3:780:LEU:HA	1:3:886:CYS:HB3	1.88	0.56
1:3:651:LEU:HD21	1:3:701:VAL:HB	1.87	0.55
1:2:454:ILE:HG13	1:2:455:ILE:HG13	1.89	0.55
1:4:822:LEU:HD11	1:4:824:GLN:O	2.06	0.55
1:4:651:LEU:HD21	1:4:701:VAL:HB	1.87	0.55
1:1:724:GLU:O	1:2:847:LYS:NZ	2.40	0.55
1:4:693:GLN:OE1	1:4:724:GLU:HB2	2.06	0.55
1:1:232:ASN:ND2	1:1:237:ARG:HB2	2.21	0.55
1:2:118:ASN:ND2	1:2:191:TRP:HB2	2.22	0.55
1:3:822:LEU:HD11	1:3:824:GLN:O	2.06	0.55
1:2:693:GLN:OE1	1:2:724:GLU:HB2	2.07	0.55
1:1:689:GLU:HG3	1:1:690:SER:H	1.72	0.55
1:3:246:MET:C	1:3:246:MET:HE3	2.31	0.55
1:3:246:MET:HE3	1:3:246:MET:O	2.06	0.55
1:3:499:ILE:HD11	1:3:529:GLU:CD	2.32	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:822:LEU:HD11	1:2:824:GLN:O	2.07	0.55
1:2:830:LEU:HD11	1:2:835:LEU:HD22	1.89	0.55
1:3:689:GLU:HG3	1:3:690:SER:H	1.72	0.55
1:2:557:ARG:NH1	4:2:8402:DMS:H13	2.22	0.55
1:4:403:ASP:OD2	1:4:451:PRO:HD2	2.07	0.55
1:1:59:ARG:HB2	1:1:124:SER:OG	2.07	0.55
1:3:316:HIS:HB2	1:3:321:THR:O	2.08	0.55
1:1:100:TYR:CE1	1:1:602:CYS:HB3	2.42	0.54
1:1:118:ASN:ND2	1:1:191:TRP:HB2	2.22	0.54
1:1:537:GLU:OE1	5:1:2001:GAL:H1	2.07	0.54
1:1:822:LEU:HD11	1:1:824:GLN:O	2.07	0.54
1:4:105:TYR:CE1	1:4:199:ASP:HB2	2.42	0.54
1:4:246:MET:HE3	1:4:246:MET:C	2.31	0.54
1:4:579:ASP:HB2	1:4:580:GLU:OE2	2.07	0.54
1:4:780:LEU:HA	1:4:886:CYS:HB3	1.88	0.54
1:1:561:ARG:HD3	1:2:525:SER:O	2.08	0.54
1:2:369:GLU:HG3	1:2:397:LEU:HD21	1.89	0.54
1:4:166:ARG:HG3	1:4:392:TYR:CB	2.37	0.54
1:4:377:LEU:HD22	1:4:708:TRP:CA	2.35	0.54
1:1:693:GLN:OE1	1:1:724:GLU:HB2	2.08	0.54
1:1:750:GLU:HG2	1:1:755:ARG:HD3	1.89	0.54
1:2:750:GLU:HG2	1:2:755:ARG:HD3	1.90	0.54
1:4:118:ASN:ND2	1:4:191:TRP:HB2	2.22	0.54
1:4:960:SER:HB2	6:4:8104:HOH:O	2.07	0.54
1:2:105:TYR:CE1	1:2:199:ASP:HB2	2.42	0.54
1:2:890:GLN:HE21	1:2:948:PRO:HG3	1.73	0.54
1:1:246:MET:C	1:1:246:MET:HE3	2.33	0.54
1:4:499:ILE:HD11	1:4:529:GLU:CD	2.32	0.54
1:1:755:ARG:HB3	1:1:769:TRP:HB2	1.90	0.54
1:2:232:ASN:ND2	1:2:237:ARG:HB2	2.23	0.54
1:3:377:LEU:HD22	1:3:708:TRP:CA	2.37	0.54
1:4:59:ARG:HB2	1:4:124:SER:OG	2.08	0.54
1:1:369:GLU:HG3	1:1:397:LEU:HD21	1.90	0.54
1:1:433:LEU:N	1:1:434:PRO:CD	2.71	0.54
1:2:44:THR:OG1	1:2:46:ARG:HG3	2.08	0.54
1:2:499:ILE:HD11	1:2:529:GLU:CD	2.33	0.54
1:1:166:ARG:HG3	1:1:392:TYR:CB	2.38	0.53
1:1:246:MET:HE2	1:1:287:ASP:HB2	1.90	0.53
1:1:595:THR:HA	1:1:596:PRO:C	2.34	0.53
1:1:683:PRO:O	1:1:685:LEU:N	2.42	0.53
1:3:830:LEU:HD11	1:3:835:LEU:HD22	1.89	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:166:ARG:HG3	1:2:392:TYR:CB	2.38	0.53
1:3:890:GLN:HE21	1:3:948:PRO:HG3	1.73	0.53
1:1:316:HIS:HB2	1:1:321:THR:O	2.07	0.53
1:3:87:PRO:HB2	1:3:209:PHE:C	2.34	0.53
1:3:595:THR:HA	1:3:596:PRO:C	2.32	0.53
1:3:613:PRO:HB3	1:3:617:LEU:HD23	1.89	0.53
1:4:369:GLU:HG3	1:4:397:LEU:HD21	1.90	0.53
1:4:542:MET:HE1	1:4:601:PHE:HA	1.88	0.53
1:2:246:MET:HE3	1:2:246:MET:O	2.08	0.53
1:2:579:ASP:HB2	1:2:580:GLU:OE2	2.08	0.53
1:3:232:ASN:ND2	1:3:237:ARG:HB2	2.24	0.53
1:4:689:GLU:HG3	1:4:690:SER:H	1.74	0.53
1:2:595:THR:HA	1:2:596:PRO:C	2.33	0.53
1:2:857:ARG:HG2	1:2:857:ARG:HH11	1.73	0.53
1:3:579:ASP:HB2	1:3:580:GLU:OE2	2.09	0.53
1:1:105:TYR:CE1	1:1:199:ASP:HB2	2.44	0.53
1:4:613:PRO:HB3	1:4:617:LEU:HD23	1.91	0.53
1:1:499:ILE:HD11	1:1:529:GLU:CD	2.34	0.53
1:4:163:GLN:O	1:4:164:ASP:CB	2.57	0.53
1:2:59:ARG:HB2	1:2:124:SER:OG	2.09	0.52
1:2:246:MET:HE2	1:2:287:ASP:HB2	1.90	0.52
1:2:755:ARG:HB3	1:2:769:TRP:HB2	1.91	0.52
1:3:166:ARG:HG3	1:3:392:TYR:CB	2.39	0.52
1:3:454:ILE:HG13	1:3:455:ILE:HG13	1.91	0.52
1:3:820:ALA:HB2	1:3:842:TRP:NE1	2.24	0.52
1:4:794:GLY:HA3	6:4:8105:HOH:O	2.08	0.52
1:1:749:ILE:N	1:1:749:ILE:HD12	2.24	0.52
1:2:749:ILE:N	1:2:749:ILE:HD12	2.24	0.52
1:3:844:HIS:CE1	1:3:845:GLN:HG3	2.44	0.52
1:4:100:TYR:CE1	1:4:602:CYS:HB3	2.44	0.52
1:2:408:TYR:HB3	1:2:454:ILE:HD13	1.91	0.52
1:3:486:TYR:CE2	1:3:488:GLY:HA3	2.44	0.52
1:3:749:ILE:HD12	1:3:749:ILE:N	2.25	0.52
1:1:780:LEU:HD22	1:1:1020:TRP:HZ3	1.74	0.52
1:3:105:TYR:CE1	1:3:199:ASP:HB2	2.45	0.52
1:4:830:LEU:HD11	1:4:835:LEU:HD22	1.90	0.52
1:1:87:PRO:HB2	1:1:209:PHE:C	2.34	0.52
1:1:249:GLU:HG2	1:1:251:ARG:CZ	2.40	0.52
1:1:689:GLU:CG	1:1:690:SER:H	2.23	0.52
1:2:24:LEU:HD11	1:3:13:ARG:NH2	2.25	0.52
1:4:44:THR:OG1	1:4:46:ARG:HG3	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:78:LEU:HB3	1:2:80:GLU:HG2	1.92	0.52
1:2:87:PRO:HB2	1:2:209:PHE:C	2.35	0.52
1:2:161:TYR:HE2	1:2:163:GLN:HE21	1.57	0.52
1:3:44:THR:OG1	1:3:46:ARG:HG3	2.08	0.52
1:3:743:SER:HB2	1:3:746:ASP:H	1.73	0.52
1:2:427:THR:O	1:2:467:ASN:HB2	2.10	0.52
1:1:579:ASP:HB2	1:1:580:GLU:OE2	2.09	0.52
1:2:100:TYR:CZ	1:2:602:CYS:HB3	2.45	0.52
1:3:780:LEU:HD22	1:3:1020:TRP:HZ3	1.75	0.52
1:4:755:ARG:HB3	1:4:769:TRP:HB2	1.90	0.52
1:4:890:GLN:HE21	1:4:948:PRO:HG3	1.75	0.52
1:2:844:HIS:CE1	1:2:845:GLN:HG3	2.45	0.52
1:3:755:ARG:HB3	1:3:769:TRP:HB2	1.90	0.52
1:1:13:ARG:HH21	1:4:24:LEU:HD11	1.74	0.51
1:2:296:GLU:HG2	4:2:8601:DMS:O	2.10	0.51
1:2:683:PRO:O	1:2:685:LEU:N	2.42	0.51
1:4:689:GLU:CG	1:4:690:SER:H	2.23	0.51
1:4:749:ILE:N	1:4:749:ILE:HD12	2.25	0.51
1:1:688:PRO:HD3	1:1:694:LEU:HD11	1.91	0.51
1:2:433:LEU:HD23	1:3:437:SER:OG	2.09	0.51
1:2:689:GLU:HG3	1:2:690:SER:H	1.74	0.51
1:2:780:LEU:HD22	1:2:1020:TRP:HZ3	1.74	0.51
1:3:118:ASN:ND2	1:3:191:TRP:HB2	2.25	0.51
1:3:249:GLU:HG2	1:3:251:ARG:CZ	2.40	0.51
1:4:246:MET:HE2	1:4:287:ASP:HB2	1.90	0.51
1:1:830:LEU:HD11	1:1:835:LEU:HD22	1.92	0.51
1:1:890:GLN:HE21	1:1:948:PRO:HG3	1.75	0.51
1:2:688:PRO:HD3	1:2:694:LEU:HD11	1.93	0.51
1:4:87:PRO:HB2	1:4:209:PHE:C	2.34	0.51
1:4:377:LEU:CD2	1:4:708:TRP:HA	2.37	0.51
1:2:128:ASN:HB3	1:2:180:GLY:C	2.36	0.51
1:3:876:THR:O	1:3:877:PRO:C	2.53	0.51
1:2:546:LEU:HD22	1:2:616:ALA:HB1	1.92	0.51
1:2:900:LEU:HB2	1:2:939:CYS:O	2.11	0.51
1:2:163:GLN:O	1:2:164:ASP:CB	2.59	0.51
1:3:427:THR:O	1:3:467:ASN:HB2	2.10	0.51
1:3:600:GLN:HB2	1:3:603:MET:HE2	1.92	0.51
1:3:688:PRO:HD3	1:3:694:LEU:HD11	1.92	0.51
1:4:249:GLU:HG2	1:4:251:ARG:CZ	2.41	0.51
1:1:163:GLN:O	1:1:164:ASP:CB	2.57	0.51
1:1:857:ARG:HG2	1:1:857:ARG:HH11	1.76	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:963:SER:HB3	1:2:983:TRP:CE2	2.46	0.51
1:4:486:TYR:CE2	1:4:488:GLY:HA3	2.46	0.51
1:4:857:ARG:HG2	1:4:857:ARG:HH11	1.76	0.51
1:1:226:HIS:ND1	1:1:448:ARG:HD3	2.26	0.51
1:1:687:GLN:OE1	1:1:687:GLN:HA	2.11	0.51
1:2:422:PRO:HG3	1:3:285:TYR:CE1	2.45	0.51
1:2:820:ALA:HB2	1:2:842:TRP:NE1	2.25	0.50
1:3:100:TYR:CZ	1:3:602:CYS:HB3	2.46	0.50
1:3:408:TYR:HB3	1:3:454:ILE:HD13	1.93	0.50
1:3:857:ARG:HH11	1:3:857:ARG:HG2	1.75	0.50
1:4:683:PRO:O	1:4:685:LEU:N	2.44	0.50
1:1:24:LEU:HD11	1:4:13:ARG:HH21	1.75	0.50
1:1:600:GLN:HB2	1:1:603:MET:HE2	1.94	0.50
1:2:743:SER:HB2	1:2:746:ASP:H	1.76	0.50
1:3:718:GLN:NE2	4:3:8503:DMS:H23	2.26	0.50
1:4:600:GLN:HB2	1:4:603:MET:HE2	1.94	0.50
1:4:687:GLN:OE1	1:4:687:GLN:HA	2.10	0.50
1:4:780:LEU:HD22	1:4:1020:TRP:HZ3	1.76	0.50
1:1:315:LEU:O	1:1:323:ILE:HB	2.12	0.50
1:1:486:TYR:CE2	1:1:488:GLY:HA3	2.46	0.50
1:2:689:GLU:CG	1:2:690:SER:H	2.25	0.50
1:4:743:SER:HB2	1:4:746:ASP:H	1.75	0.50
1:2:249:GLU:HG2	1:2:251:ARG:CZ	2.40	0.50
1:2:651:LEU:CD2	1:2:701:VAL:HB	2.41	0.50
1:2:408:TYR:HB3	1:2:454:ILE:CD1	2.42	0.50
1:3:525:SER:O	1:4:561:ARG:HD3	2.11	0.50
1:4:546:LEU:HD22	1:4:616:ALA:HB1	1.93	0.50
1:4:710:GLU:CD	1:4:710:GLU:H	2.19	0.50
1:1:246:MET:HG2	1:1:274:PHE:CE1	2.46	0.50
1:1:651:LEU:CD2	1:1:701:VAL:HB	2.42	0.50
1:1:844:HIS:CE1	1:1:845:GLN:HG3	2.47	0.50
1:1:876:THR:O	1:1:877:PRO:C	2.54	0.50
1:3:718:GLN:HE21	4:3:8503:DMS:H23	1.74	0.50
1:2:226:HIS:ND1	1:2:448:ARG:HD3	2.26	0.50
1:2:710:GLU:CD	1:2:710:GLU:H	2.20	0.50
1:1:246:MET:HE3	1:1:246:MET:O	2.11	0.50
1:1:1000:SER:O	1:1:1001:PRO:C	2.55	0.50
1:3:78:LEU:HB3	1:3:80:GLU:HG2	1.93	0.50
1:3:161:TYR:HE2	1:3:163:GLN:HE21	1.58	0.50
1:4:316:HIS:HB2	1:4:321:THR:O	2.11	0.50
1:4:844:HIS:CE1	1:4:845:GLN:HG3	2.47	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:315:LEU:O	1:2:323:ILE:HB	2.12	0.50
1:2:316:HIS:HB2	1:2:321:THR:O	2.12	0.50
1:2:377:LEU:HD22	1:2:708:TRP:CA	2.37	0.50
1:3:710:GLU:CD	1:3:710:GLU:H	2.19	0.50
1:1:403:ASP:OD2	1:1:451:PRO:HD2	2.11	0.49
1:2:533:LEU:C	1:2:533:LEU:HD23	2.37	0.49
1:3:689:GLU:CG	1:3:690:SER:H	2.24	0.49
1:3:900:LEU:HB2	1:3:939:CYS:O	2.12	0.49
1:4:963:SER:HB3	1:4:983:TRP:CE2	2.47	0.49
1:2:787:ALA:HB3	1:2:934:GLU:N	2.27	0.49
1:4:751:LEU:HD21	1:4:860:GLY:O	2.12	0.49
1:1:631:LEU:HD22	1:1:696:LEU:HD23	1.93	0.49
1:4:246:MET:HE3	1:4:246:MET:O	2.12	0.49
1:2:423:MET:HE1	1:3:281:GLU:HB2	1.94	0.49
1:3:246:MET:HE2	1:3:287:ASP:HB2	1.93	0.49
1:3:433:LEU:N	1:3:434:PRO:CD	2.75	0.49
1:3:651:LEU:CD2	1:3:701:VAL:HB	2.41	0.49
1:1:1011:ALA:HB3	1:1:1014:TYR:CZ	2.47	0.49
1:2:223:SER:O	1:2:224:ASP:HB2	2.12	0.49
1:3:233:ASP:HB3	4:3:8417:DMS:H23	1.94	0.49
1:3:751:LEU:HD21	1:3:860:GLY:O	2.12	0.49
1:4:631:LEU:HD22	1:4:696:LEU:HD23	1.93	0.49
1:4:1000:SER:O	1:4:1001:PRO:C	2.56	0.49
1:2:613:PRO:HB3	1:2:617:LEU:HD23	1.94	0.49
1:4:688:PRO:HD3	1:4:694:LEU:HD11	1.93	0.49
1:4:900:LEU:HB2	1:4:939:CYS:O	2.12	0.49
1:1:377:LEU:HD22	1:1:708:TRP:CA	2.39	0.49
1:3:226:HIS:ND1	1:3:448:ARG:HD3	2.28	0.49
1:3:408:TYR:HB3	1:3:454:ILE:CD1	2.42	0.49
1:3:683:PRO:O	1:3:685:LEU:N	2.45	0.49
1:3:687:GLN:OE1	1:3:687:GLN:HA	2.12	0.49
1:4:651:LEU:CD2	1:4:701:VAL:HB	2.43	0.49
1:4:820:ALA:HB2	1:4:842:TRP:NE1	2.28	0.49
1:1:751:LEU:HD21	1:1:860:GLY:O	2.13	0.49
1:1:820:ALA:HB2	1:1:842:TRP:NE1	2.27	0.49
1:4:78:LEU:HB3	1:4:80:GLU:HG2	1.94	0.49
1:4:246:MET:HG2	1:4:274:PHE:CE1	2.47	0.49
1:4:599:ARG:HG3	6:4:8028:HOH:O	2.12	0.49
1:1:223:SER:O	1:1:224:ASP:HB2	2.13	0.48
1:2:1000:SER:O	1:2:1001:PRO:C	2.54	0.48
1:1:100:TYR:CZ	1:1:602:CYS:HB3	2.48	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:333:ARG:HG2	1:2:333:ARG:HH11	1.78	0.48
1:2:631:LEU:HD22	1:2:696:LEU:HD23	1.94	0.48
1:2:815:HIS:HE1	1:2:877:PRO:O	1.96	0.48
1:4:333:ARG:HG2	1:4:333:ARG:HH11	1.77	0.48
1:1:251:ARG:HG3	1:1:251:ARG:HH11	1.78	0.48
1:4:876:THR:O	1:4:877:PRO:C	2.55	0.48
1:1:13:ARG:NH2	1:4:24:LEU:HD11	2.29	0.48
1:2:687:GLN:OE1	1:2:687:GLN:HA	2.12	0.48
1:1:710:GLU:CD	1:1:710:GLU:H	2.20	0.48
1:2:441:THR:HG22	1:2:474:TRP:CZ2	2.49	0.48
1:3:377:LEU:CD2	1:3:708:TRP:HA	2.38	0.48
1:3:581:ASN:HB2	1:3:583:ASN:OD1	2.14	0.48
1:4:441:THR:HG22	1:4:474:TRP:CZ2	2.48	0.48
1:1:78:LEU:HB3	1:1:80:GLU:HG2	1.96	0.48
1:1:743:SER:HB2	1:1:746:ASP:H	1.78	0.48
1:1:900:LEU:HB2	1:1:939:CYS:O	2.13	0.48
1:3:18:ASN:ND2	1:3:21:VAL:HG23	2.29	0.48
1:3:963:SER:HB3	1:3:983:TRP:CE2	2.48	0.48
1:1:282:ARG:HG2	1:4:423:MET:CE	2.43	0.48
1:1:474:TRP:CE2	1:1:478:VAL:HG21	2.48	0.48
1:2:19:PRO:HD3	1:2:112:PRO:CB	2.44	0.48
1:3:196:TYR:O	1:3:417:THR:HG22	2.13	0.48
1:3:403:ASP:OD2	1:3:451:PRO:HD2	2.13	0.48
1:4:533:LEU:C	1:4:533:LEU:HD23	2.39	0.48
1:3:128:ASN:HB3	1:3:180:GLY:C	2.38	0.48
1:1:546:LEU:HD22	1:1:616:ALA:HB1	1.96	0.48
1:1:815:HIS:HE1	1:1:877:PRO:O	1.96	0.48
1:3:24:LEU:O	1:3:25:ASN:HB2	2.14	0.48
1:1:408:TYR:HB3	1:1:454:ILE:HD13	1.96	0.47
1:1:515:VAL:HG21	1:4:281:GLU:CD	2.39	0.47
1:2:486:TYR:CE2	1:2:488:GLY:HA3	2.49	0.47
1:3:844:HIS:ND1	1:3:845:GLN:HG3	2.29	0.47
1:1:124:SER:HA	1:1:184:LEU:O	2.14	0.47
1:1:963:SER:HB3	1:1:983:TRP:CE2	2.50	0.47
1:3:546:LEU:HD22	1:3:616:ALA:HB1	1.96	0.47
1:4:128:ASN:HB3	1:4:180:GLY:C	2.39	0.47
1:2:600:GLN:HB2	1:2:603:MET:HE2	1.95	0.47
1:3:815:HIS:HE1	1:3:877:PRO:O	1.98	0.47
1:1:433:LEU:HB3	1:1:434:PRO:HD3	1.96	0.47
1:2:276:GLY:CA	4:2:8001:DMS:H11	2.30	0.47
1:2:433:LEU:N	1:2:434:PRO:CD	2.77	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:4:100:TYR:CZ	1:4:602:CYS:HB3	2.49	0.47
1:1:441:THR:HG22	1:1:474:TRP:CZ2	2.48	0.47
1:1:536:CYS:O	1:1:537:GLU:HG3	2.15	0.47
1:2:876:THR:O	1:2:877:PRO:C	2.56	0.47
1:4:474:TRP:CE2	1:4:478:VAL:HG21	2.49	0.47
1:1:126:THR:H	4:1:8502:DMS:C1	2.27	0.47
1:1:128:ASN:HB3	1:1:180:GLY:C	2.40	0.47
1:1:949:HIS:O	1:1:1023:LYS:HD2	2.15	0.47
1:3:441:THR:HG22	1:3:474:TRP:CZ2	2.50	0.47
1:4:226:HIS:ND1	1:4:448:ARG:HD3	2.29	0.47
1:4:255:ARG:HB2	1:4:316:HIS:CE1	2.49	0.47
1:1:18:ASN:ND2	1:1:21:VAL:HG23	2.30	0.47
1:1:19:PRO:HD3	1:1:112:PRO:CB	2.45	0.47
1:1:499:ILE:HG22	1:1:501:PRO:HD3	1.97	0.47
1:2:40:GLU:HA	1:2:40:GLU:OE1	2.15	0.47
1:2:474:TRP:CE2	1:2:478:VAL:HG21	2.50	0.47
1:3:533:LEU:C	1:3:533:LEU:HD23	2.40	0.47
1:3:1000:SER:O	1:3:1001:PRO:C	2.56	0.47
1:4:223:SER:O	1:4:224:ASP:HB2	2.13	0.47
1:4:815:HIS:HE1	1:4:877:PRO:O	1.98	0.47
1:2:937:LEU:O	1:2:938:ARG:HD2	2.15	0.47
1:2:949:HIS:O	1:2:1023:LYS:HD2	2.15	0.47
1:3:40:GLU:OE1	1:3:40:GLU:HA	2.14	0.47
1:3:1011:ALA:HB3	1:3:1014:TYR:CZ	2.50	0.47
1:4:127:PHE:CE1	1:4:184:LEU:HG	2.50	0.47
1:4:433:LEU:N	1:4:434:PRO:CD	2.77	0.47
1:2:418:GLU:HB2	1:2:461:GLU:HB3	1.96	0.47
1:3:902:PRO:HD2	4:3:8415:DMS:H21	1.97	0.47
1:2:124:SER:HA	1:2:184:LEU:O	2.15	0.47
1:2:988:GLY:C	1:2:989:PHE:CD1	2.93	0.47
1:3:19:PRO:HD3	1:3:112:PRO:CB	2.44	0.47
1:3:486:TYR:CZ	1:3:488:GLY:HA3	2.50	0.47
1:3:631:LEU:HD22	1:3:696:LEU:HD23	1.97	0.47
1:1:474:TRP:CZ2	1:1:478:VAL:HG21	2.50	0.46
1:2:708:TRP:CE3	1:2:709:SER:HB3	2.51	0.46
1:3:163:GLN:O	1:3:164:ASP:CB	2.62	0.46
1:4:486:TYR:CZ	1:4:488:GLY:HA3	2.50	0.46
1:4:603:MET:C	1:4:604:ASN:HD22	2.23	0.46
1:4:1011:ALA:HB3	1:4:1014:TYR:CZ	2.50	0.46
1:2:603:MET:C	1:2:604:ASN:HD22	2.23	0.46
1:2:820:ALA:HB2	1:2:842:TRP:CE2	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:59:ARG:CG	4:1:8502:DMS:H12	2.46	0.46
1:1:533:LEU:HD23	1:1:533:LEU:C	2.40	0.46
1:1:581:ASN:HB2	1:1:583:ASN:OD1	2.16	0.46
1:1:593:GLY:O	1:1:595:THR:HG22	2.15	0.46
1:2:433:LEU:HB3	1:2:434:PRO:HD3	1.97	0.46
1:2:1006:GLU:HG2	1:2:1007:PHE:CD2	2.51	0.46
1:3:59:ARG:HB2	1:3:124:SER:HG	1.79	0.46
1:3:127:PHE:CE1	1:3:184:LEU:HG	2.50	0.46
1:3:367:MET:HB3	1:3:372:MET:HE2	1.97	0.46
1:3:460:ASN:O	1:3:461:GLU:C	2.58	0.46
1:4:40:GLU:OE1	1:4:40:GLU:HA	2.14	0.46
1:1:787:ALA:HB3	1:1:934:GLU:N	2.30	0.46
1:2:18:ASN:ND2	1:2:21:VAL:HG23	2.31	0.46
1:3:290:THR:H	4:3:8001:DMS:C2	2.28	0.46
1:3:820:ALA:HB2	1:3:842:TRP:CE2	2.50	0.46
1:4:942:ARG:HA	1:4:953:GLY:O	2.15	0.46
1:2:537:GLU:OE1	5:2:2001:GAL:H1	2.16	0.46
1:4:408:TYR:HB3	1:4:454:ILE:HD13	1.97	0.46
1:4:427:THR:O	1:4:467:ASN:HB2	2.15	0.46
1:2:423:MET:CE	1:3:281:GLU:HB2	2.45	0.46
1:3:802:ASP:O	1:3:808:GLU:HG3	2.16	0.46
1:4:161:TYR:HE2	1:4:163:GLN:HE21	1.61	0.46
1:4:460:ASN:O	1:4:461:GLU:C	2.59	0.46
1:2:844:HIS:ND1	1:2:845:GLN:HG3	2.31	0.46
1:3:474:TRP:CE2	1:3:478:VAL:HG21	2.50	0.46
1:4:147:ASN:HA	1:4:148:SER:HA	1.71	0.46
1:1:161:TYR:HE2	1:1:163:GLN:HE21	1.60	0.46
1:1:255:ARG:HB2	1:1:316:HIS:CE1	2.51	0.46
1:2:751:LEU:HD21	1:2:860:GLY:O	2.16	0.46
1:2:1011:ALA:HB3	1:2:1014:TYR:CZ	2.50	0.46
1:4:433:LEU:HB3	1:4:434:PRO:HD3	1.98	0.46
1:1:333:ARG:HH11	1:1:333:ARG:HG2	1.81	0.46
1:2:50:GLN:CD	1:2:50:GLN:N	2.74	0.46
1:2:377:LEU:CD2	1:2:708:TRP:HA	2.39	0.46
1:2:499:ILE:HG22	1:2:501:PRO:HD3	1.98	0.46
1:3:50:GLN:N	1:3:50:GLN:CD	2.74	0.46
1:3:311:ALA:HB2	1:3:330:VAL:HG21	1.97	0.46
1:1:24:LEU:HD11	1:4:13:ARG:NH2	2.31	0.46
1:1:50:GLN:N	1:1:50:GLN:CD	2.74	0.46
1:2:593:GLY:O	1:2:595:THR:HG22	2.16	0.46
1:3:768:MET:HE1	1:3:1020:TRP:CZ2	2.52	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:3:499:ILE:HG22	1:3:501:PRO:HD3	1.97	0.45
1:3:593:GLY:O	1:3:595:THR:HG22	2.16	0.45
1:4:499:ILE:HG22	1:4:501:PRO:HD3	1.97	0.45
1:4:536:CYS:O	1:4:537:GLU:HG3	2.16	0.45
1:1:942:ARG:HA	1:1:953:GLY:O	2.17	0.45
1:3:474:TRP:CZ2	1:3:478:VAL:HG21	2.51	0.45
1:3:890:GLN:HE21	1:3:948:PRO:CG	2.29	0.45
1:4:787:ALA:HB3	1:4:934:GLU:N	2.30	0.45
1:1:408:TYR:HB3	1:1:454:ILE:CD1	2.46	0.45
1:1:418:GLU:HB2	1:1:461:GLU:HB3	1.99	0.45
1:3:127:PHE:HE1	1:3:184:LEU:HG	1.82	0.45
1:3:536:CYS:O	1:3:537:GLU:HG3	2.16	0.45
1:1:708:TRP:CE3	1:1:709:SER:HB3	2.52	0.45
1:1:917:ARG:HH12	1:1:943:GLU:CD	2.24	0.45
1:2:857:ARG:HG2	1:2:857:ARG:NH1	2.30	0.45
1:3:1020:TRP:HD1	1:3:1021:CYS:N	2.14	0.45
1:4:232:ASN:HD22	1:4:237:ARG:HB2	1.80	0.45
1:4:487:GLU:OE1	1:4:502:MET:HE3	2.16	0.45
1:4:768:MET:HE1	1:4:1020:TRP:CZ2	2.52	0.45
1:4:988:GLY:C	1:4:989:PHE:CD1	2.95	0.45
1:1:1006:GLU:HG2	1:1:1007:PHE:CD2	2.52	0.45
1:2:403:ASP:OD2	1:2:451:PRO:HD2	2.16	0.45
1:2:581:ASN:HB2	1:2:583:ASN:OD1	2.17	0.45
1:3:433:LEU:HB3	1:3:434:PRO:HD3	1.97	0.45
1:4:910:LEU:C	1:4:910:LEU:HD12	2.41	0.45
1:4:937:LEU:O	1:4:938:ARG:HD2	2.16	0.45
1:4:963:SER:HB3	1:4:983:TRP:NE1	2.32	0.45
1:1:127:PHE:CE1	1:1:184:LEU:HG	2.51	0.45
1:1:796:SER:HB2	1:1:802:ASP:HB3	1.99	0.45
1:2:231:PHE:CD1	1:2:231:PHE:N	2.85	0.45
1:2:910:LEU:C	1:2:910:LEU:HD12	2.41	0.45
1:3:124:SER:HA	1:3:184:LEU:O	2.16	0.45
1:3:787:ALA:HB3	1:3:934:GLU:N	2.31	0.45
1:3:847:LYS:HG3	1:3:849:LEU:HD23	1.98	0.45
1:4:19:PRO:HD3	1:4:112:PRO:CB	2.46	0.45
1:4:124:SER:HA	1:4:184:LEU:O	2.17	0.45
1:1:988:GLY:C	1:1:989:PHE:CD1	2.95	0.45
1:2:24:LEU:O	1:2:25:ASN:HB2	2.17	0.45
1:2:474:TRP:CZ2	1:2:478:VAL:HG21	2.52	0.45
1:2:890:GLN:HE21	1:2:948:PRO:CG	2.29	0.45
1:3:223:SER:O	1:3:224:ASP:HB2	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:3:315:LEU:O	1:3:323:ILE:HB	2.16	0.45
1:3:1006:GLU:HG2	1:3:1007:PHE:CD2	2.52	0.45
1:4:251:ARG:HG3	1:4:251:ARG:HH11	1.81	0.45
1:4:418:GLU:HB2	1:4:461:GLU:HB3	1.99	0.45
1:4:474:TRP:CZ2	1:4:478:VAL:HG21	2.51	0.45
1:4:487:GLU:CD	1:4:502:MET:HE3	2.42	0.45
1:1:196:TYR:O	1:1:417:THR:HG22	2.17	0.45
1:1:246:MET:HG2	1:1:274:PHE:CZ	2.51	0.45
1:2:196:TYR:O	1:2:417:THR:HG22	2.16	0.45
1:2:423:MET:HE2	1:3:282:ARG:HG2	1.99	0.45
1:3:949:HIS:O	1:3:1023:LYS:HD2	2.16	0.45
1:4:50:GLN:CD	1:4:50:GLN:N	2.75	0.45
1:4:581:ASN:HB2	1:4:583:ASN:OD1	2.17	0.45
1:1:857:ARG:HG2	1:1:857:ARG:NH1	2.32	0.45
1:2:311:ALA:HB2	1:2:330:VAL:HG21	1.98	0.45
1:2:542:MET:HE1	1:2:601:PHE:CG	2.52	0.45
1:3:147:ASN:HA	1:3:148:SER:HA	1.71	0.45
1:4:542:MET:HE1	1:4:601:PHE:CG	2.52	0.45
1:4:937:LEU:HD12	1:4:957:PHE:C	2.42	0.45
1:1:40:GLU:OE1	1:1:40:GLU:HA	2.17	0.45
1:1:231:PHE:CD1	1:1:231:PHE:N	2.84	0.45
1:1:542:MET:HE3	1:1:604:ASN:HD21	1.82	0.45
1:1:631:LEU:HD12	1:1:635:THR:O	2.17	0.45
1:2:127:PHE:CE1	1:2:184:LEU:HG	2.51	0.45
1:2:279:ILE:HD11	1:3:422:PRO:CG	2.47	0.45
1:3:251:ARG:HG3	1:3:251:ARG:HH11	1.81	0.45
1:1:285:TYR:CD1	1:4:422:PRO:HG3	2.52	0.44
1:1:487:GLU:CD	1:1:502:MET:HE3	2.42	0.44
1:1:820:ALA:HB2	1:1:842:TRP:CE2	2.52	0.44
1:1:910:LEU:C	1:1:910:LEU:HD12	2.42	0.44
1:2:246:MET:HG2	1:2:274:PHE:CE1	2.51	0.44
1:2:301:TRP:CH2	1:2:452:SER:HA	2.52	0.44
1:2:917:ARG:HH12	1:2:943:GLU:CD	2.25	0.44
1:3:369:GLU:O	1:3:373:VAL:HG23	2.17	0.44
1:3:603:MET:C	1:3:604:ASN:HD22	2.25	0.44
1:4:164:ASP:HB3	6:4:8033:HOH:O	2.16	0.44
1:4:820:ALA:HB2	1:4:842:TRP:CE2	2.53	0.44
1:4:890:GLN:HE21	1:4:948:PRO:CG	2.30	0.44
1:1:807:VAL:HG13	1:1:808:GLU:N	2.32	0.44
1:3:260:LEU:O	1:3:267:VAL:HG22	2.17	0.44
1:4:196:TYR:O	1:4:417:THR:HG22	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:301:TRP:CH2	1:1:452:SER:HA	2.51	0.44
1:2:251:ARG:HG3	1:2:251:ARG:HH11	1.81	0.44
1:2:674:PRO:O	1:2:675:GLN:HB2	2.17	0.44
1:2:870:VAL:HG12	1:2:871:GLU:N	2.33	0.44
1:3:301:TRP:CH2	1:3:452:SER:HA	2.52	0.44
1:3:674:PRO:O	1:3:675:GLN:HB2	2.18	0.44
1:4:231:PHE:N	1:4:231:PHE:CD1	2.85	0.44
1:4:796:SER:HB2	1:4:802:ASP:HB3	2.00	0.44
1:1:649:ASN:O	1:1:650:GLU:HG3	2.17	0.44
1:2:416:GLU:HG3	1:2:460:ASN:O	2.17	0.44
1:2:460:ASN:O	1:2:461:GLU:C	2.60	0.44
1:3:390:SER:HA	1:3:391:HIS:HA	1.81	0.44
1:3:478:VAL:HG13	4:3:8420:DMS:H22	1.99	0.44
1:3:600:GLN:HB2	1:3:603:MET:CE	2.47	0.44
1:4:966:GLN:NE2	1:4:979:GLU:OE2	2.51	0.44
1:4:1020:TRP:HD1	1:4:1021:CYS:N	2.14	0.44
1:1:768:MET:HE1	1:1:1020:TRP:CZ2	2.53	0.44
1:1:937:LEU:O	1:1:938:ARG:HD2	2.18	0.44
1:2:200:GLN:N	1:2:200:GLN:OE1	2.51	0.44
1:3:796:SER:HB2	1:3:802:ASP:HB3	1.98	0.44
1:3:988:GLY:C	1:3:989:PHE:CD1	2.96	0.44
1:4:127:PHE:HE1	1:4:184:LEU:HG	1.82	0.44
1:4:708:TRP:CE3	1:4:709:SER:HB3	2.53	0.44
1:2:423:MET:H	1:3:282:ARG:HB2	1.83	0.44
1:2:536:CYS:O	1:2:537:GLU:HG3	2.18	0.44
1:3:857:ARG:HG2	1:3:857:ARG:NH1	2.32	0.44
1:4:408:TYR:HB3	1:4:454:ILE:CD1	2.48	0.44
1:1:896:ASN:HA	1:1:918:TRP:O	2.18	0.44
1:1:937:LEU:HD12	1:1:957:PHE:C	2.43	0.44
1:2:587:ALA:HB1	1:2:591:ASP:CB	2.48	0.44
1:3:870:VAL:HG12	1:3:871:GLU:N	2.32	0.44
1:3:942:ARG:HA	1:3:953:GLY:O	2.17	0.44
1:4:233:ASP:HA	4:4:8004:DMS:S	2.58	0.44
1:1:487:GLU:OE1	1:1:502:MET:HE3	2.18	0.44
1:1:662:PRO:O	1:1:663:LEU:HD23	2.18	0.44
1:1:847:LYS:HG3	1:1:849:LEU:HD23	2.00	0.44
1:3:937:LEU:O	1:3:938:ARG:HD2	2.17	0.44
1:1:41:GLU:OE2	1:1:46:ARG:NH2	2.48	0.44
1:1:870:VAL:HG12	1:1:871:GLU:N	2.33	0.44
1:3:842:TRP:HZ3	1:3:852:SER:HB3	1.82	0.44
1:4:369:GLU:O	1:4:373:VAL:HG23	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:147:ASN:HA	1:1:148:SER:HA	1.71	0.43
1:1:603:MET:C	1:1:604:ASN:HD22	2.26	0.43
1:2:255:ARG:HB2	1:2:316:HIS:CE1	2.53	0.43
1:3:708:TRP:CH2	4:3:8403:DMS:H23	2.53	0.43
1:3:910:LEU:C	1:3:910:LEU:HD12	2.43	0.43
1:4:118:ASN:HD21	1:4:191:TRP:HB2	1.83	0.43
1:4:864:MET:HE3	1:4:866:ILE:HD11	2.00	0.43
1:4:870:VAL:HG12	1:4:871:GLU:N	2.31	0.43
1:1:844:HIS:ND1	1:1:845:GLN:HG3	2.32	0.43
1:3:255:ARG:HB2	1:3:316:HIS:CE1	2.53	0.43
1:3:662:PRO:O	1:3:663:LEU:HD23	2.17	0.43
1:4:873:ALA:C	1:4:875:ASP:H	2.27	0.43
1:1:200:GLN:N	1:1:200:GLN:OE1	2.51	0.43
1:1:377:LEU:CD2	1:1:708:TRP:HA	2.41	0.43
1:1:460:ASN:O	1:1:461:GLU:C	2.62	0.43
1:1:890:GLN:HE21	1:1:948:PRO:CG	2.30	0.43
1:1:1020:TRP:HD1	1:1:1021:CYS:N	2.16	0.43
1:2:240:LEU:C	1:2:240:LEU:HD23	2.42	0.43
1:2:282:ARG:HB2	1:3:423:MET:N	2.33	0.43
1:2:1020:TRP:HD1	1:2:1021:CYS:N	2.15	0.43
1:3:427:THR:HG21	1:3:462:SER:HB3	2.00	0.43
1:4:416:GLU:HG3	1:4:460:ASN:O	2.18	0.43
1:1:418:GLU:OE1	1:1:461:GLU:OE1	2.36	0.43
1:1:587:ALA:HB1	1:1:591:ASP:CB	2.49	0.43
1:2:802:ASP:O	1:2:808:GLU:HG3	2.19	0.43
1:2:847:LYS:HG3	1:2:849:LEU:HD23	1.99	0.43
1:2:873:ALA:C	1:2:875:ASP:H	2.26	0.43
1:3:587:ALA:HB1	1:3:591:ASP:CB	2.48	0.43
1:3:631:LEU:HD12	1:3:635:THR:O	2.18	0.43
1:3:649:ASN:O	1:3:650:GLU:HG3	2.19	0.43
1:4:857:ARG:HG2	1:4:857:ARG:NH1	2.33	0.43
1:1:542:MET:HE1	1:1:601:PHE:CG	2.54	0.43
1:2:502:MET:HB2	1:2:537:GLU:HB2	2.00	0.43
1:2:513:PRO:O	1:2:514:ALA:HB3	2.19	0.43
1:3:966:GLN:NE2	1:3:979:GLU:OE2	2.50	0.43
1:4:24:LEU:O	1:4:25:ASN:HB2	2.18	0.43
1:4:291:LEU:N	1:4:291:LEU:HD22	2.33	0.43
1:1:127:PHE:HE1	1:1:184:LEU:HG	1.84	0.43
1:1:311:ALA:HB2	1:1:330:VAL:HG21	1.99	0.43
1:3:876:THR:O	1:3:876:THR:HG23	2.18	0.43
1:4:568:TRP:CD2	1:4:569:ASP:HB3	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:35:SER:HB2	1:1:217:LYS:HD2	2.01	0.43
1:1:486:TYR:CZ	1:1:488:GLY:HA3	2.54	0.43
1:2:963:SER:HB3	1:2:983:TRP:NE1	2.33	0.43
1:3:367:MET:HB3	1:3:372:MET:CE	2.48	0.43
1:3:917:ARG:HH12	1:3:943:GLU:CD	2.24	0.43
1:4:674:PRO:O	1:4:675:GLN:HB2	2.19	0.43
1:3:231:PHE:N	1:3:231:PHE:CD1	2.87	0.43
1:1:883:GLY:HA3	1:1:987:ASP:HA	2.01	0.43
1:2:437:SER:OG	1:3:433:LEU:HD23	2.19	0.43
1:3:23:GLN:C	1:3:23:GLN:CD	2.86	0.43
1:3:246:MET:HG2	1:3:274:PHE:CE1	2.54	0.43
1:3:333:ARG:HH11	1:3:333:ARG:HG2	1.82	0.43
1:4:111:PRO:HA	1:4:112:PRO:HA	1.83	0.43
1:4:917:ARG:HH12	1:4:943:GLU:CD	2.26	0.43
1:1:600:GLN:HB2	1:1:603:MET:CE	2.49	0.43
1:2:942:ARG:HA	1:2:953:GLY:O	2.17	0.43
1:3:502:MET:HB2	1:3:537:GLU:HB2	2.00	0.43
1:3:658:LEU:O	1:3:659:ASP:C	2.61	0.43
1:3:1008:GLN:O	1:3:1010:SER:N	2.52	0.43
1:4:1006:GLU:HG2	1:4:1007:PHE:CD2	2.54	0.43
1:2:13:ARG:NH2	1:3:24:LEU:HD11	2.34	0.42
1:2:24:LEU:HD11	1:3:13:ARG:HH21	1.83	0.42
1:2:639:THR:HG23	1:2:677:LYS:HG2	2.01	0.42
1:2:662:PRO:O	1:2:663:LEU:HD23	2.19	0.42
1:2:768:MET:HE1	1:2:1020:TRP:CZ2	2.54	0.42
1:2:887:GLN:OE1	1:2:980:GLU:O	2.37	0.42
1:3:200:GLN:OE1	1:3:200:GLN:N	2.51	0.42
1:3:639:THR:HG23	1:3:677:LYS:HG2	2.00	0.42
1:4:427:THR:HG21	1:4:462:SER:HB3	2.00	0.42
1:1:842:TRP:HZ3	1:1:852:SER:HB3	1.83	0.42
1:2:842:TRP:HZ3	1:2:852:SER:HB3	1.84	0.42
1:2:864:MET:HE3	1:2:866:ILE:HD11	2.01	0.42
1:4:59:ARG:NH2	1:4:81:ALA:HB3	2.34	0.42
1:2:631:LEU:HD12	1:2:635:THR:O	2.19	0.42
1:3:240:LEU:HD23	1:3:240:LEU:C	2.45	0.42
1:3:418:GLU:HB2	1:3:461:GLU:HB3	2.01	0.42
1:3:542:MET:HE3	1:3:604:ASN:HD21	1.85	0.42
1:4:949:HIS:O	1:4:1023:LYS:HD2	2.19	0.42
1:1:995:GLY:C	1:1:997:ASP:N	2.77	0.42
1:2:807:VAL:HG13	1:2:808:GLU:N	2.34	0.42
1:3:111:PRO:HA	1:3:112:PRO:HA	1.84	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:118:ASN:HD21	1:1:191:TRP:HB2	1.84	0.42
1:1:367:MET:HB3	1:1:372:MET:HE2	2.01	0.42
1:1:427:THR:O	1:1:467:ASN:HB2	2.19	0.42
1:1:873:ALA:C	1:1:875:ASP:H	2.26	0.42
1:2:796:SER:HB2	1:2:802:ASP:HB3	2.00	0.42
1:4:301:TRP:CH2	1:4:452:SER:HA	2.54	0.42
1:4:593:GLY:O	1:4:595:THR:HG22	2.19	0.42
1:1:173:LEU:O	1:1:174:SER:C	2.62	0.42
1:2:502:MET:HG3	1:2:502:MET:O	2.20	0.42
1:3:825:CYS:HA	1:3:837:THR:O	2.20	0.42
1:4:842:TRP:HZ3	1:4:852:SER:HB3	1.85	0.42
1:1:291:LEU:N	1:1:291:LEU:HD22	2.35	0.42
1:1:674:PRO:O	1:1:675:GLN:HB2	2.19	0.42
1:2:418:GLU:OE1	1:2:461:GLU:OE1	2.37	0.42
1:2:542:MET:HE3	1:2:604:ASN:HD21	1.84	0.42
1:2:607:VAL:HG12	1:2:613:PRO:HA	2.02	0.42
1:2:896:ASN:HA	1:2:918:TRP:O	2.19	0.42
1:3:902:PRO:HD2	4:3:8415:DMS:S	2.59	0.42
1:4:71:GLU:O	1:4:72:SER:C	2.63	0.42
1:4:876:THR:O	1:4:876:THR:HG23	2.19	0.42
1:4:894:ARG:HH12	1:4:921:PRO:HD3	1.83	0.42
1:1:367:MET:HB3	1:1:372:MET:CE	2.49	0.42
1:1:526:LEU:HD23	1:1:526:LEU:HA	1.86	0.42
1:1:689:GLU:HG3	1:1:690:SER:N	2.35	0.42
1:2:59:ARG:HB2	1:2:124:SER:HG	1.85	0.42
1:2:290:THR:N	4:2:8001:DMS:H12	2.32	0.42
1:2:486:TYR:CZ	1:2:488:GLY:HA3	2.55	0.42
1:3:542:MET:HE1	1:3:601:PHE:CG	2.55	0.42
1:3:708:TRP:CE3	1:3:709:SER:HB3	2.54	0.42
1:1:240:LEU:C	1:1:240:LEU:HD23	2.45	0.42
1:1:252:ASP:OD1	4:1:8416:DMS:S	2.77	0.42
1:2:165:SER:O	1:2:209:PHE:HZ	2.03	0.42
1:2:568:TRP:CD2	1:2:569:ASP:HB3	2.55	0.42
1:3:379:MET:HE1	1:3:387:VAL:HB	2.02	0.42
1:3:568:TRP:HA	1:3:569:ASP:HA	1.77	0.42
1:3:807:VAL:HG13	1:3:808:GLU:N	2.34	0.42
1:4:18:ASN:ND2	1:4:21:VAL:HG23	2.34	0.42
1:4:102:ASN:HA	1:4:201:ASP:OD1	2.20	0.42
1:4:367:MET:HB3	1:4:372:MET:CE	2.50	0.42
1:4:844:HIS:ND1	1:4:845:GLN:HG3	2.35	0.42
1:4:847:LYS:HG3	1:4:849:LEU:HD23	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:379:MET:HE1	1:1:387:VAL:HB	2.02	0.41
1:1:658:LEU:O	1:1:659:ASP:C	2.63	0.41
1:2:418:GLU:HA	1:2:423:MET:HG3	2.02	0.41
1:2:421:VAL:HG13	1:3:282:ARG:O	2.20	0.41
1:3:416:GLU:HG3	1:3:460:ASN:O	2.19	0.41
1:3:568:TRP:CD2	1:3:569:ASP:HB3	2.55	0.41
1:1:23:GLN:C	1:1:23:GLN:CD	2.88	0.41
1:1:148:SER:HA	1:1:165:SER:OG	2.20	0.41
1:1:802:ASP:O	1:1:808:GLU:HG3	2.19	0.41
1:2:231:PHE:CD2	1:2:238:ALA:HB2	2.56	0.41
1:2:423:MET:N	1:3:282:ARG:HB2	2.35	0.41
1:2:526:LEU:HD23	1:2:526:LEU:HA	1.87	0.41
1:2:883:GLY:HA3	1:2:987:ASP:HA	2.02	0.41
1:2:986:ILE:CG2	1:2:1018:LEU:HD11	2.48	0.41
1:3:148:SER:HA	1:3:165:SER:OG	2.20	0.41
1:3:444:VAL:O	1:3:448:ARG:HB3	2.21	0.41
1:3:513:PRO:O	1:3:514:ALA:HB3	2.21	0.41
1:4:240:LEU:C	1:4:240:LEU:HD23	2.46	0.41
1:4:825:CYS:HA	1:4:837:THR:O	2.21	0.41
1:4:995:GLY:C	1:4:997:ASP:N	2.78	0.41
1:1:277:GLU:CD	1:1:277:GLU:H	2.27	0.41
1:1:730:LEU:HB3	1:1:731:PRO:HD2	2.02	0.41
1:2:361:PRO:HB3	1:2:609:ALA:CB	2.47	0.41
1:2:746:ASP:HA	1:2:760:ARG:HG3	2.02	0.41
1:2:907:PRO:HG2	1:2:990:HIS:O	2.20	0.41
1:3:487:GLU:CD	1:3:502:MET:HE3	2.45	0.41
1:3:607:VAL:HG12	1:3:613:PRO:HA	2.01	0.41
1:3:847:LYS:NZ	1:4:724:GLU:O	2.53	0.41
1:3:883:GLY:HA3	1:3:987:ASP:HA	2.03	0.41
1:3:937:LEU:HD12	1:3:957:PHE:C	2.45	0.41
1:1:361:PRO:HB3	1:1:609:ALA:CB	2.47	0.41
1:2:35:SER:HB2	1:2:217:LYS:HD2	2.02	0.41
1:2:127:PHE:HE1	1:2:184:LEU:HG	1.83	0.41
1:2:937:LEU:HD12	1:2:957:PHE:C	2.46	0.41
1:3:482:ARG:HA	1:3:483:PRO:HD3	1.88	0.41
1:4:23:GLN:CD	1:4:23:GLN:C	2.87	0.41
1:4:277:GLU:CD	1:4:277:GLU:H	2.28	0.41
1:4:367:MET:HB3	1:4:372:MET:HE2	2.01	0.41
1:4:789:LEU:HD11	1:4:993:ILE:HG22	2.02	0.41
1:1:24:LEU:O	1:1:25:ASN:HB2	2.19	0.41
1:1:463:GLY:HA2	6:1:8649:HOH:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:683:PRO:O	1:1:684:GLU:C	2.63	0.41
1:2:600:GLN:HB2	1:2:603:MET:CE	2.51	0.41
1:1:502:MET:HB2	1:1:537:GLU:HB2	2.01	0.41
1:1:607:VAL:HG12	1:1:613:PRO:HA	2.03	0.41
1:1:825:CYS:HA	1:1:837:THR:O	2.21	0.41
1:2:96:ASP:OD1	1:2:96:ASP:C	2.64	0.41
1:3:652:LEU:HD23	1:3:680:ILE:HD12	2.02	0.41
1:3:873:ALA:C	1:3:875:ASP:H	2.28	0.41
1:4:103:VAL:HG22	1:4:418:GLU:CD	2.46	0.41
1:4:246:MET:HG2	1:4:274:PHE:CZ	2.55	0.41
1:4:896:ASN:HA	1:4:918:TRP:O	2.21	0.41
1:1:369:GLU:O	1:1:373:VAL:HG23	2.21	0.41
1:1:689:GLU:CG	1:1:690:SER:N	2.84	0.41
1:2:69:VAL:HA	1:2:70:PRO:HD2	1.91	0.41
1:4:907:PRO:HG2	1:4:990:HIS:O	2.21	0.41
1:1:153:TRP:CD1	1:1:158:TRP:HA	2.56	0.41
1:1:232:ASN:HD22	1:1:237:ARG:HB2	1.85	0.41
1:1:513:PRO:O	1:1:514:ALA:HB3	2.21	0.41
1:1:652:LEU:HD23	1:1:680:ILE:HD12	2.03	0.41
1:1:864:MET:HE3	1:1:866:ILE:HD11	2.03	0.41
1:2:118:ASN:HD21	1:2:191:TRP:HB2	1.84	0.41
1:2:658:LEU:O	1:2:659:ASP:C	2.63	0.41
1:2:825:CYS:HA	1:2:837:THR:O	2.21	0.41
1:3:361:PRO:HB3	1:3:609:ALA:CB	2.48	0.41
1:3:963:SER:HB3	1:3:983:TRP:NE1	2.36	0.41
1:4:444:VAL:O	1:4:448:ARG:HB3	2.21	0.41
1:4:502:MET:HB2	1:4:537:GLU:HB2	2.03	0.41
1:4:881:ARG:HE	1:4:987:ASP:CG	2.27	0.41
1:1:165:SER:O	1:1:209:PHE:HZ	2.04	0.41
1:1:246:MET:HE2	1:1:287:ASP:CB	2.51	0.41
1:1:772:ASP:OD2	1:1:773:LYS:HE2	2.21	0.41
1:2:92:MET:HE3	1:2:205:MET:CE	2.51	0.41
1:2:127:PHE:CD1	1:2:127:PHE:N	2.89	0.41
1:2:382:ASN:HA	1:2:621:LYS:HG3	2.03	0.41
1:3:71:GLU:O	1:3:72:SER:C	2.64	0.41
1:3:352:ARG:HG2	1:3:553:TRP:CH2	2.56	0.41
1:3:944:LEU:O	1:3:950:GLN:HA	2.21	0.41
1:4:381:GLN:O	1:4:621:LYS:HE3	2.21	0.41
1:1:414:ASN:O	1:1:439:ARG:HD3	2.21	0.40
1:1:726:LEU:HD22	1:2:871:GLU:OE1	2.21	0.40
1:1:756:TRP:CD2	1:1:858:ILE:HD13	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:873:ALA:O	1:1:876:THR:HG22	2.21	0.40
1:2:966:GLN:NE2	1:2:979:GLU:OE2	2.52	0.40
1:3:756:TRP:CD2	1:3:858:ILE:HD13	2.56	0.40
1:4:607:VAL:HG12	1:4:613:PRO:HA	2.02	0.40
1:1:789:LEU:HD11	1:1:993:ILE:HG22	2.04	0.40
1:1:853:ARG:HG3	1:1:853:ARG:HH11	1.87	0.40
1:1:876:THR:O	1:1:876:THR:HG23	2.21	0.40
1:1:963:SER:HB3	1:1:983:TRP:NE1	2.36	0.40
1:2:246:MET:HE2	1:2:287:ASP:CB	2.51	0.40
1:2:959:ILE:HG23	1:2:959:ILE:O	2.21	0.40
1:3:277:GLU:H	1:3:277:GLU:CD	2.29	0.40
1:3:871:GLU:OE1	1:4:726:LEU:HD22	2.21	0.40
1:4:69:VAL:HA	1:4:70:PRO:HD2	1.91	0.40
1:4:382:ASN:HA	1:4:621:LYS:HG3	2.03	0.40
1:4:390:SER:HA	1:4:391:HIS:HA	1.82	0.40
1:4:587:ALA:HB1	1:4:591:ASP:CB	2.51	0.40
1:1:71:GLU:O	1:1:72:SER:C	2.64	0.40
1:1:427:THR:HG21	1:1:462:SER:HB3	2.03	0.40
1:1:542:MET:HE3	1:1:604:ASN:ND2	2.36	0.40
1:1:887:GLN:OE1	1:1:980:GLU:O	2.39	0.40
1:2:265:THR:HG22	1:2:266:GLN:N	2.37	0.40
1:2:756:TRP:CD2	1:2:858:ILE:HD13	2.56	0.40
1:3:291:LEU:HD22	1:3:291:LEU:N	2.37	0.40
1:4:173:LEU:O	1:4:174:SER:C	2.64	0.40
1:1:944:LEU:O	1:1:950:GLN:HA	2.22	0.40
1:1:959:ILE:O	1:1:959:ILE:HG23	2.21	0.40
1:2:708:TRP:CZ3	1:2:709:SER:HB3	2.57	0.40
1:2:869:ASP:CG	1:2:1015:HIS:HD1	2.29	0.40
1:2:927:THR:HA	1:2:928:PRO:HD2	1.95	0.40
1:3:246:MET:HG2	1:3:274:PHE:CZ	2.57	0.40
1:3:615:PRO:HD2	6:3:8808:HOH:O	2.20	0.40
1:3:782:ASP:OD1	1:3:842:TRP:HH2	2.04	0.40
1:3:853:ARG:HH11	1:3:853:ARG:HG3	1.86	0.40
1:1:102:ASN:HA	1:1:201:ASP:OD1	2.22	0.40
1:1:416:GLU:HG3	1:1:460:ASN:O	2.21	0.40
1:1:907:PRO:HG2	1:1:990:HIS:O	2.22	0.40
1:2:147:ASN:HA	1:2:148:SER:HA	1.72	0.40
1:2:333:ARG:HG2	1:2:333:ARG:NH1	2.36	0.40
1:2:427:THR:HG21	1:2:462:SER:HB3	2.03	0.40
1:3:526:LEU:HA	1:3:526:LEU:HD23	1.86	0.40
1:3:746:ASP:HA	1:3:760:ARG:HG3	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:3:894:ARG:HH12	1:3:921:PRO:HD3	1.87	0.40
1:4:165:SER:O	1:4:209:PHE:HZ	2.04	0.40
1:4:333:ARG:HG2	1:4:333:ARG:NH1	2.36	0.40
1:4:619:GLU:HA	1:4:912:ALA:HB2	2.03	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	1	1009/1023 (99%)	912 (90%)	85 (8%)	12 (1%)	10	40
1	2	1009/1023 (99%)	907 (90%)	91 (9%)	11 (1%)	11	43
1	3	1009/1023 (99%)	911 (90%)	88 (9%)	10 (1%)	12	45
1	4	1009/1023 (99%)	911 (90%)	87 (9%)	11 (1%)	11	43
All	All	4036/4092 (99%)	3641 (90%)	351 (9%)	44 (1%)	11	43

All (44) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	1	722	LEU
1	2	722	LEU
1	3	722	LEU
1	4	722	LEU
1	1	164	ASP
1	1	684	GLU
1	2	164	ASP
1	2	684	GLU
1	3	684	GLU
1	4	164	ASP
1	4	684	GLU

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Mol	Chain	Res	Type
1	1	211	ASP
1	1	461	GLU
1	1	832	ASP
1	2	211	ASP
1	2	461	GLU
1	2	591	ASP
1	2	832	ASP
1	3	164	ASP
1	3	211	ASP
1	3	461	GLU
1	3	832	ASP
1	3	1009	LEU
1	4	211	ASP
1	4	461	GLU
1	4	832	ASP
1	1	102	ASN
1	1	591	ASP
1	2	1009	LEU
1	4	591	ASP
1	1	1009	LEU
1	2	102	ASN
1	4	14	ARG
1	4	102	ASN
1	1	14	ARG
1	3	102	ASN
1	1	511	PRO
1	1	891	VAL
1	2	891	VAL
1	4	511	PRO
1	2	511	PRO
1	3	891	VAL
1	3	511	PRO
1	4	731	PRO

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	1	864/875 (99%)	849 (98%)	15 (2%)	53	78
1	2	864/875 (99%)	851 (98%)	13 (2%)	57	80
1	3	864/875 (99%)	850 (98%)	14 (2%)	55	79
1	4	864/875 (99%)	849 (98%)	15 (2%)	53	78
All	All	3456/3500 (99%)	3399 (98%)	57 (2%)	55	79

All (57) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	1	71	GLU
1	1	128	ASN
1	1	219	THR
1	1	231	PHE
1	1	246	MET
1	1	252	ASP
1	1	324	GLU
1	1	333	ARG
1	1	519	SER
1	1	546	LEU
1	1	580	GLU
1	1	595	THR
1	1	672	VAL
1	1	761	GLN
1	1	1017	GLN
1	2	71	GLU
1	2	128	ASN
1	2	219	THR
1	2	246	MET
1	2	252	ASP
1	2	324	GLU
1	2	333	ARG
1	2	519	SER
1	2	546	LEU
1	2	580	GLU
1	2	595	THR
1	2	672	VAL
1	2	761	GLN
1	3	71	GLU
1	3	128	ASN
1	3	219	THR
1	3	246	MET
1	3	252	ASP

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Mol	Chain	Res	Type
1	3	324	GLU
1	3	333	ARG
1	3	519	SER
1	3	546	LEU
1	3	580	GLU
1	3	595	THR
1	3	672	VAL
1	3	761	GLN
1	3	1017	GLN
1	4	71	GLU
1	4	128	ASN
1	4	219	THR
1	4	246	MET
1	4	252	ASP
1	4	324	GLU
1	4	333	ARG
1	4	519	SER
1	4	546	LEU
1	4	580	GLU
1	4	595	THR
1	4	672	VAL
1	4	761	GLN
1	4	800	ARG
1	4	1017	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (41) such sidechains are listed below:

Mol	Chain	Res	Type
1	1	38	ASN
1	1	50	GLN
1	1	93	HIS
1	1	485	GLN
1	1	702	GLN
1	1	878	HIS
1	1	885	ASN
1	1	887	GLN
1	1	1008	GLN
1	2	50	GLN
1	2	93	HIS
1	2	163	GLN
1	2	262	GLN
1	2	485	GLN

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Mol	Chain	Res	Type
1	2	702	GLN
1	2	878	HIS
1	2	885	ASN
1	2	887	GLN
1	2	1008	GLN
1	3	38	ASN
1	3	49	GLN
1	3	50	GLN
1	3	93	HIS
1	3	485	GLN
1	3	702	GLN
1	3	718	GLN
1	3	878	HIS
1	3	885	ASN
1	3	887	GLN
1	3	1008	GLN
1	4	38	ASN
1	4	50	GLN
1	4	245	GLN
1	4	294	ASN
1	4	485	GLN
1	4	702	GLN
1	4	713	HIS
1	4	878	HIS
1	4	885	ASN
1	4	887	GLN
1	4	1008	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 104 ligands modelled in this entry, 23 are monoatomic - leaving 81 for Mogul analysis.

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	$\# Z > 2$
4	DMS	1	8003	-	3,3,3	0.26	0	3,3,3	0.67	0
4	DMS	1	8503	-	3,3,3	0.28	0	3,3,3	0.67	0
4	DMS	4	8002	-	3,3,3	0.30	0	3,3,3	0.67	0
4	DMS	2	8503	-	3,3,3	0.26	0	3,3,3	0.68	0
4	DMS	3	8000	-	3,3,3	0.26	0	3,3,3	0.68	0
4	DMS	2	8423	-	3,3,3	0.27	0	3,3,3	0.71	0
4	DMS	3	8419	-	3,3,3	0.29	0	3,3,3	0.68	0
4	DMS	2	8001	-	3,3,3	0.29	0	3,3,3	0.68	0
4	DMS	3	8501	-	3,3,3	0.25	0	3,3,3	0.65	0
4	DMS	1	8402	-	3,3,3	0.25	0	3,3,3	0.64	0
4	DMS	3	8401	-	3,3,3	0.25	0	3,3,3	0.63	0
4	DMS	3	8411	-	3,3,3	0.26	0	3,3,3	0.68	0
4	DMS	1	8423	-	3,3,3	0.25	0	3,3,3	0.65	0
4	DMS	1	8421	-	3,3,3	0.28	0	3,3,3	0.69	0
4	DMS	1	8401	-	3,3,3	0.20	0	3,3,3	0.57	0
4	DMS	2	8003	-	3,3,3	0.24	0	3,3,3	0.68	0
4	DMS	3	8402	-	3,3,3	0.28	0	3,3,3	0.66	0
4	DMS	1	8411	-	3,3,3	0.22	0	3,3,3	0.71	0
4	DMS	3	8405	-	3,3,3	0.26	0	3,3,3	0.70	0
5	GAL	3	2002	-	12,12,12	0.87	0	17,17,17	0.83	0
4	DMS	1	8406	-	3,3,3	0.20	0	3,3,3	0.62	0
5	GAL	1	2002	-	12,12,12	1.00	0	17,17,17	0.82	0
4	DMS	4	8001	-	3,3,3	0.22	0	3,3,3	0.60	0
5	GAL	2	2001	3	12,12,12	0.82	0	17,17,17	0.80	0
4	DMS	2	8601	-	3,3,3	0.31	0	3,3,3	0.66	0
4	DMS	1	8425	-	3,3,3	0.25	0	3,3,3	0.65	0
5	GAL	1	2001	2,3	12,12,12	0.77	0	17,17,17	0.86	0
4	DMS	2	8419	-	3,3,3	0.27	0	3,3,3	0.67	0
4	DMS	3	8425	-	3,3,3	0.21	0	3,3,3	0.67	0
4	DMS	3	8406	-	3,3,3	0.26	0	3,3,3	0.67	0
4	DMS	1	8410	-	3,3,3	0.23	0	3,3,3	0.66	0
4	DMS	1	8501	-	3,3,3	0.28	0	3,3,3	0.60	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	GAL	2	2002	-	12,12,12	1.00	0	17,17,17	0.80	0
4	DMS	3	8001	-	3,3,3	0.23	0	3,3,3	0.65	0
4	DMS	4	8005	-	3,3,3	0.30	0	3,3,3	0.71	0
4	DMS	1	8602	-	3,3,3	0.27	0	3,3,3	0.66	0
4	DMS	4	8004	-	3,3,3	0.27	0	3,3,3	0.64	0
4	DMS	3	8403	-	3,3,3	0.18	0	3,3,3	0.62	0
4	DMS	2	8414	-	3,3,3	0.25	0	3,3,3	0.65	0
4	DMS	2	8002	-	3,3,3	0.19	0	3,3,3	0.63	0
4	DMS	1	8403	-	3,3,3	0.28	0	3,3,3	0.70	0
4	DMS	1	8416	-	3,3,3	0.25	0	3,3,3	0.68	0
4	DMS	1	8409	-	3,3,3	0.25	0	3,3,3	0.68	0
4	DMS	1	8002	-	3,3,3	0.22	0	3,3,3	0.64	0
4	DMS	2	8409	-	3,3,3	0.24	0	3,3,3	0.67	0
4	DMS	2	8504	-	3,3,3	0.26	0	3,3,3	0.59	0
4	DMS	3	8416	-	3,3,3	0.24	0	3,3,3	0.70	0
4	DMS	1	8502	-	3,3,3	0.26	0	3,3,3	0.60	0
4	DMS	1	8405	-	3,3,3	0.31	0	3,3,3	0.68	0
4	DMS	3	8407	-	3,3,3	0.25	0	3,3,3	0.71	0
4	DMS	2	8421	-	3,3,3	0.26	0	3,3,3	0.66	0
4	DMS	4	8006	-	3,3,3	0.29	0	3,3,3	0.74	0
4	DMS	2	8405	-	3,3,3	0.28	0	3,3,3	0.69	0
4	DMS	3	8421	-	3,3,3	0.24	0	3,3,3	0.66	0
5	GAL	3	2001	3	12,12,12	0.73	0	17,17,17	0.79	0
4	DMS	3	8414	-	3,3,3	0.24	0	3,3,3	0.67	0
4	DMS	2	8427	-	3,3,3	0.22	0	3,3,3	0.66	0
4	DMS	2	8401	-	3,3,3	0.27	0	3,3,3	0.60	0
4	DMS	3	8404	-	3,3,3	0.28	0	3,3,3	0.72	0
4	DMS	4	8003	-	3,3,3	0.31	0	3,3,3	0.69	0
4	DMS	2	8000	-	3,3,3	0.32	0	3,3,3	0.67	0
4	DMS	2	8411	-	3,3,3	0.25	0	3,3,3	0.63	0
4	DMS	1	8415	-	3,3,3	0.30	0	3,3,3	0.64	0
4	DMS	2	8402	-	3,3,3	0.24	0	3,3,3	0.63	0
4	DMS	3	8503	-	3,3,3	0.27	0	3,3,3	0.68	0
4	DMS	1	8417	-	3,3,3	0.23	0	3,3,3	0.66	0
4	DMS	3	8420	-	3,3,3	0.26	0	3,3,3	0.62	0
4	DMS	1	8414	-	3,3,3	0.32	0	3,3,3	0.68	0
4	DMS	1	8404	-	3,3,3	0.23	0	3,3,3	0.71	0
4	DMS	2	8404	-	3,3,3	0.23	0	3,3,3	0.63	0
5	GAL	4	2001	3	12,12,12	1.15	1 (8%)	17,17,17	0.77	0
5	GAL	4	2002	-	12,12,12	0.81	0	17,17,17	0.82	0
4	DMS	2	8501	-	3,3,3	0.28	0	3,3,3	0.65	0
4	DMS	1	8408	-	3,3,3	0.31	0	3,3,3	0.66	0
4	DMS	3	8415	-	3,3,3	0.27	0	3,3,3	0.64	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	DMS	1	8001	-	3,3,3	0.23	0	3,3,3	0.67	0
4	DMS	3	8417	-	3,3,3	0.26	0	3,3,3	0.70	0
4	DMS	3	8705	-	3,3,3	0.23	0	3,3,3	0.66	0
4	DMS	3	8408	-	3,3,3	0.24	0	3,3,3	0.70	0
4	DMS	3	8409	-	3,3,3	0.30	0	3,3,3	0.69	0
4	DMS	2	8403	-	3,3,3	0.32	0	3,3,3	0.68	0

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
5	GAL	3	2002	-	-	2/2/22/22	0/1/1/1
5	GAL	1	2001	2,3	-	2/2/22/22	0/1/1/1
5	GAL	1	2002	-	-	2/2/22/22	0/1/1/1
5	GAL	4	2001	3	-	1/2/22/22	0/1/1/1
5	GAL	4	2002	-	-	2/2/22/22	0/1/1/1
5	GAL	2	2002	-	-	2/2/22/22	0/1/1/1
5	GAL	3	2001	3	-	1/2/22/22	0/1/1/1
5	GAL	2	2001	3	-	1/2/22/22	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	4	2001	GAL	O5-C1	2.36	1.48	1.42

There are no bond angle outliers.

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	2	2002	GAL	O5-C5-C6-O6
5	2	2002	GAL	C4-C5-C6-O6
5	1	2001	GAL	O5-C5-C6-O6
5	1	2001	GAL	C4-C5-C6-O6
5	1	2002	GAL	O5-C5-C6-O6
5	1	2002	GAL	C4-C5-C6-O6
5	4	2002	GAL	O5-C5-C6-O6
5	4	2002	GAL	C4-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
5	2	2001	GAL	O5-C5-C6-O6
5	4	2001	GAL	O5-C5-C6-O6
5	3	2001	GAL	O5-C5-C6-O6
5	3	2002	GAL	C4-C5-C6-O6
5	3	2002	GAL	O5-C5-C6-O6

There are no ring outliers.

23 monomers are involved in 36 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	1	8503	DMS	1	0
4	2	8001	DMS	6	0
4	3	8402	DMS	2	0
5	2	2001	GAL	1	0
4	2	8601	DMS	1	0
5	1	2001	GAL	1	0
4	3	8001	DMS	2	0
4	4	8004	DMS	1	0
4	3	8403	DMS	1	0
4	1	8416	DMS	1	0
4	1	8502	DMS	4	0
4	4	8006	DMS	1	0
4	2	8401	DMS	1	0
4	2	8000	DMS	1	0
4	2	8411	DMS	1	0
4	2	8402	DMS	1	0
4	3	8503	DMS	2	0
4	3	8420	DMS	1	0
4	1	8414	DMS	2	0
4	1	8408	DMS	1	0
4	3	8415	DMS	2	0
4	1	8001	DMS	1	0
4	3	8417	DMS	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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	1	1011/1023 (98%)	0.25	18 (1%) 67 44	18, 51, 89, 108	0
1	2	1011/1023 (98%)	0.06	29 (2%) 53 31	15, 39, 91, 114	0
1	3	1011/1023 (98%)	-0.18	6 (0%) 85 69	16, 35, 57, 89	0
1	4	1011/1023 (98%)	-0.11	6 (0%) 85 69	13, 36, 59, 86	0
All	All	4044/4092 (98%)	0.01	59 (1%) 72 49	13, 39, 79, 114	0

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	2	735	HIS	4.2
1	2	734	SER	4.1
1	1	686	PRO	3.4
1	3	688	PRO	3.4
1	1	733	ALA	3.4
1	2	730	LEU	3.3
1	2	763	GLY	3.2
1	2	653	HIS	3.0
1	2	764	PHE	3.0
1	2	768	MET	3.0
1	4	730	LEU	2.9
1	2	745	MET	2.9
1	2	729	THR	2.8
1	4	688	PRO	2.8
1	2	736	ALA	2.8
1	1	730	LEU	2.8
1	2	737	ILE	2.8
1	1	319	ASP	2.8
1	2	821	ALA	2.7
1	2	733	ALA	2.7
1	4	731	PRO	2.6

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Mol	Chain	Res	Type	RSRZ
1	2	739	HIS	2.6
1	1	890	GLN	2.6
1	2	49	GLN	2.6
1	2	765	LEU	2.5
1	1	861	SER	2.5
1	4	798	ALA	2.5
1	2	823	LEU	2.5
1	3	755	ARG	2.5
1	2	1022	GLN	2.5
1	2	685	LEU	2.4
1	1	830	LEU	2.4
1	4	689	GLU	2.4
1	2	688	PRO	2.3
1	2	732	ALA	2.3
1	1	798	ALA	2.3
1	2	856	TYR	2.3
1	2	771	GLY	2.2
1	2	766	SER	2.2
1	1	863	GLN	2.2
1	2	862	GLY	2.2
1	3	690	SER	2.2
1	4	690	SER	2.2
1	2	799	THR	2.2
1	1	831	ALA	2.2
1	2	767	GLN	2.1
1	3	80	GLU	2.1
1	2	758	PHE	2.1
1	1	728	VAL	2.1
1	1	763	GLY	2.1
1	1	688	PRO	2.1
1	1	726	LEU	2.1
1	1	732	ALA	2.1
1	1	751	LEU	2.0
1	3	730	LEU	2.0
1	2	1020	TRP	2.0
1	3	735	HIS	2.0
1	1	737	ILE	2.0
1	1	770	ILE	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	DMS	1	8416	4/4	0.32	0.35	115,115,115,117	0
4	DMS	3	8421	4/4	0.33	0.31	120,121,121,121	0
4	DMS	1	8417	4/4	0.45	0.39	106,106,106,108	0
4	DMS	3	8414	4/4	0.46	0.28	114,114,115,115	0
4	DMS	1	8425	4/4	0.47	0.35	124,125,125,126	0
4	DMS	1	8410	4/4	0.48	0.30	110,110,110,111	0
4	DMS	1	8503	4/4	0.54	0.39	115,116,116,116	0
4	DMS	2	8503	4/4	0.55	0.34	105,106,106,107	0
4	DMS	1	8602	4/4	0.56	0.28	115,115,115,115	0
4	DMS	3	8416	4/4	0.58	0.28	106,106,107,108	0
3	NA	2	3102	1/1	0.59	0.12	49,49,49,49	0
4	DMS	3	8503	4/4	0.59	0.31	105,105,106,106	0
4	DMS	3	8415	4/4	0.70	0.45	133,133,133,134	0
4	DMS	2	8423	4/4	0.72	0.30	84,85,85,86	0
4	DMS	1	8415	4/4	0.75	0.28	84,84,85,85	0
4	DMS	1	8421	4/4	0.75	0.25	88,88,89,90	0
4	DMS	2	8421	4/4	0.77	0.27	116,116,116,117	0
4	DMS	2	8427	4/4	0.78	0.27	78,79,79,80	0
4	DMS	1	8502	4/4	0.79	0.25	91,91,92,92	0
4	DMS	3	8407	4/4	0.80	0.23	79,79,80,81	0
3	NA	3	3102	1/1	0.80	0.09	44,44,44,44	0
4	DMS	3	8417	4/4	0.81	0.27	82,82,83,84	0
3	NA	1	3103	1/1	0.82	0.07	43,43,43,43	0
4	DMS	3	8705	4/4	0.82	0.28	107,107,107,107	0
4	DMS	4	8004	4/4	0.82	0.27	80,81,81,81	0
2	MG	2	3002	1/1	0.84	0.18	26,26,26,26	1
4	DMS	2	8003	4/4	0.84	0.23	75,75,75,76	0
4	DMS	3	8406	4/4	0.84	0.25	78,79,79,80	0
4	DMS	3	8419	4/4	0.85	0.16	85,86,86,86	0
4	DMS	3	8411	4/4	0.86	0.23	67,67,68,69	0
4	DMS	4	8003	4/4	0.86	0.20	56,56,57,58	0
4	DMS	1	8414	4/4	0.86	0.18	54,54,56,56	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	DMS	4	8005	4/4	0.86	0.21	74,74,76,76	0
5	GAL	2	2002	12/12	0.86	0.16	81,82,83,84	0
2	MG	1	3002	1/1	0.88	0.23	63,63,63,63	0
4	DMS	1	8409	4/4	0.88	0.21	65,66,67,67	0
3	NA	2	3103	1/1	0.89	0.10	30,30,30,30	0
4	DMS	1	8001	4/4	0.89	0.20	73,73,74,75	0
4	DMS	2	8419	4/4	0.89	0.15	69,69,70,70	0
4	DMS	1	8406	4/4	0.89	0.22	82,82,83,84	0
5	GAL	1	2002	12/12	0.89	0.14	76,79,80,80	0
4	DMS	3	8420	4/4	0.89	0.22	65,65,66,66	0
4	DMS	2	8000	4/4	0.90	0.16	56,57,57,58	0
3	NA	2	3100	1/1	0.90	0.12	36,36,36,36	0
4	DMS	4	8006	4/4	0.90	0.16	57,59,59,60	0
4	DMS	1	8423	4/4	0.90	0.18	82,83,84,84	0
4	DMS	3	8403	4/4	0.90	0.14	48,48,50,50	0
5	GAL	3	2002	12/12	0.90	0.15	76,78,79,80	0
4	DMS	2	8501	4/4	0.91	0.19	62,62,62,63	0
4	DMS	2	8414	4/4	0.91	0.17	74,74,74,74	0
4	DMS	1	8405	4/4	0.91	0.16	62,62,63,64	0
4	DMS	1	8501	4/4	0.91	0.15	58,58,58,59	0
2	MG	1	3001	1/1	0.91	0.13	131,131,131,131	0
4	DMS	2	8409	4/4	0.91	0.14	42,43,43,43	0
5	GAL	4	2002	12/12	0.91	0.15	79,80,81,82	0
4	DMS	2	8402	4/4	0.92	0.17	71,71,72,72	0
4	DMS	2	8403	4/4	0.92	0.12	40,41,42,43	0
3	NA	3	3103	1/1	0.92	0.17	48,48,48,48	0
4	DMS	1	8403	4/4	0.92	0.12	52,52,54,54	0
4	DMS	2	8504	4/4	0.92	0.17	53,53,54,54	0
3	NA	1	3100	1/1	0.92	0.07	44,44,44,44	0
4	DMS	1	8003	4/4	0.92	0.17	79,79,79,80	0
4	DMS	2	8411	4/4	0.93	0.14	53,53,54,54	0
4	DMS	2	8601	4/4	0.93	0.15	63,63,63,64	0
4	DMS	1	8408	4/4	0.93	0.13	62,62,62,63	0
4	DMS	2	8001	4/4	0.93	0.13	46,46,46,47	0
4	DMS	3	8425	4/4	0.93	0.23	5,6,6,7	4
5	GAL	2	2001	12/12	0.93	0.10	39,41,42,42	0
4	DMS	1	8002	4/4	0.93	0.14	72,73,73,74	0
3	NA	4	3100	1/1	0.93	0.30	74,74,74,74	0
5	GAL	4	2001	12/12	0.93	0.09	21,29,30,31	0
4	DMS	3	8001	4/4	0.93	0.14	55,55,55,56	0
4	DMS	3	8409	4/4	0.94	0.14	46,47,47,48	0
5	GAL	1	2001	12/12	0.94	0.09	37,40,43,44	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	NA	4	3101	1/1	0.94	0.11	31,31,31,31	0
4	DMS	1	8402	4/4	0.94	0.13	66,66,67,67	0
4	DMS	4	8002	4/4	0.95	0.12	48,49,50,51	0
4	DMS	3	8405	4/4	0.95	0.12	51,52,52,53	0
4	DMS	2	8002	4/4	0.95	0.13	67,67,68,68	0
5	GAL	3	2001	12/12	0.95	0.08	30,32,33,35	0
4	DMS	3	8402	4/4	0.95	0.12	45,46,47,47	0
4	DMS	1	8411	4/4	0.95	0.11	47,48,48,49	0
4	DMS	3	8404	4/4	0.95	0.14	39,40,40,41	0
4	DMS	1	8404	4/4	0.96	0.11	50,50,50,51	0
4	DMS	3	8000	4/4	0.96	0.10	55,55,56,57	0
4	DMS	2	8401	4/4	0.96	0.12	45,45,46,46	0
2	MG	2	3001	1/1	0.96	0.09	57,57,57,57	0
4	DMS	3	8408	4/4	0.96	0.12	58,59,59,60	0
3	NA	1	3101	1/1	0.96	0.06	30,30,30,30	0
4	DMS	2	8405	4/4	0.96	0.11	45,46,46,47	0
3	NA	3	3101	1/1	0.96	0.06	38,38,38,38	0
4	DMS	3	8501	4/4	0.97	0.09	44,44,45,45	0
3	NA	2	3101	1/1	0.97	0.07	25,25,25,25	0
3	NA	3	3100	1/1	0.97	0.11	16,16,16,16	0
4	DMS	3	8401	4/4	0.98	0.08	36,38,38,38	0
4	DMS	2	8404	4/4	0.98	0.09	44,44,45,46	0
2	MG	3	3001	1/1	0.98	0.04	48,48,48,48	0
2	MG	4	3000	1/1	0.98	0.06	40,40,40,40	0
2	MG	1	3000	1/1	0.98	0.07	18,18,18,18	0
4	DMS	1	8401	4/4	0.98	0.07	39,40,41,41	0
2	MG	2	3000	1/1	0.98	0.04	30,30,30,30	0
4	DMS	4	8001	4/4	0.99	0.05	30,30,31,32	0
2	MG	3	3000	1/1	0.99	0.03	9,9,9,9	0
2	MG	4	3004	1/1	1.00	0.03	1,1,1,1	0

6.5 Other polymers ⓘ

There are no such residues in this entry.