



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 01:45 PM UTC

PDB ID : 3MUY / pdb_00003muy
Title : E. coli (lacZ) beta-galactosidase (R599A)
Authors : Dugdale, M.L.; Vance, M.; Driedger, M.L.; Nibber, A.; Tran, A.; Huber, R.E.
Deposited on : 2010-05-03
Resolution : 2.50 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

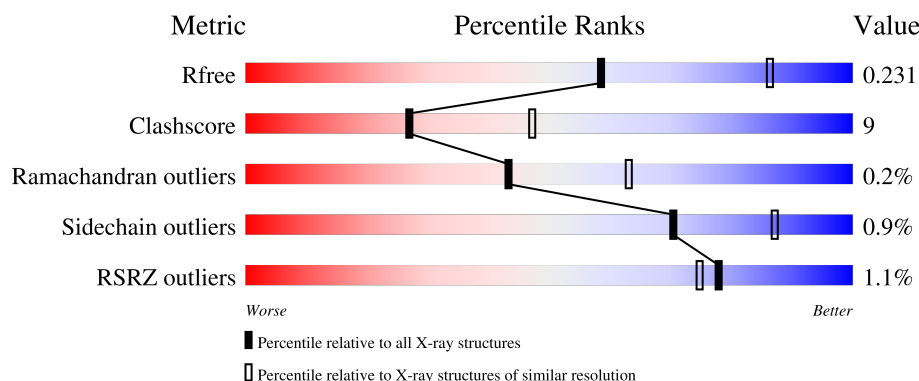
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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% 77% 21% ..
1	2	1023	 % 74% 24% ..
1	3	1023	 % 74% 24% ..
1	4	1023	 % 75% 22% ..

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 36360 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-D-galactosidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	1	1011	Total	C	N	O	S	0	0	0
			8119	5135	1437	1509	38			
1	2	1011	Total	C	N	O	S	0	0	0
			8119	5135	1437	1509	38			
1	3	1011	Total	C	N	O	S	0	0	0
			8119	5135	1437	1509	38			
1	4	1011	Total	C	N	O	S	0	0	0
			8119	5135	1437	1509	38			

There are 36 discrepancies between the modelled and reference sequences:

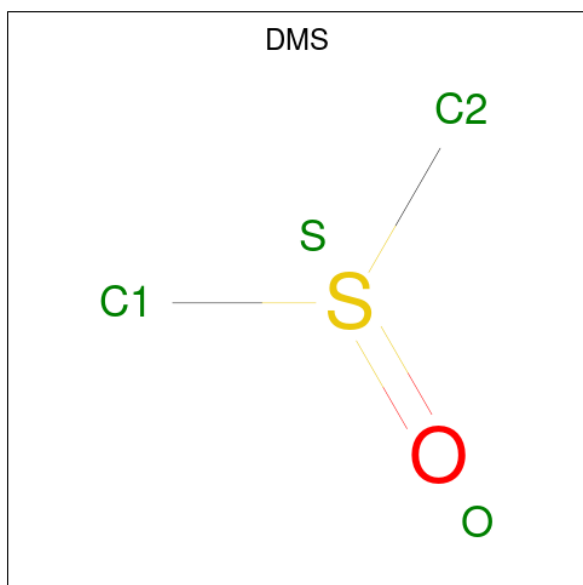
Chain	Residue	Modelled	Actual	Comment	Reference
1	1	GLY	-	expression tag	UNP C6UB28
1	2	SER	-	expression tag	UNP C6UB28
1	3	HIS	-	expression tag	UNP C6UB28
1	4	MET	-	expression tag	UNP C6UB28
1	5	LEU	-	expression tag	UNP C6UB28
1	6	GLU	-	expression tag	UNP C6UB28
1	7	ASP	-	expression tag	UNP C6UB28
1	8	PRO	-	expression tag	UNP C6UB28
1	599	ALA	ARG	engineered mutation	UNP C6UB28
2	1	GLY	-	expression tag	UNP C6UB28
2	2	SER	-	expression tag	UNP C6UB28
2	3	HIS	-	expression tag	UNP C6UB28
2	4	MET	-	expression tag	UNP C6UB28
2	5	LEU	-	expression tag	UNP C6UB28
2	6	GLU	-	expression tag	UNP C6UB28
2	7	ASP	-	expression tag	UNP C6UB28
2	8	PRO	-	expression tag	UNP C6UB28
2	599	ALA	ARG	engineered mutation	UNP C6UB28
3	1	GLY	-	expression tag	UNP C6UB28
3	2	SER	-	expression tag	UNP C6UB28
3	3	HIS	-	expression tag	UNP C6UB28

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Chain	Residue	Modelled	Actual	Comment	Reference
3	4	MET	-	expression tag	UNP C6UB28
3	5	LEU	-	expression tag	UNP C6UB28
3	6	GLU	-	expression tag	UNP C6UB28
3	7	ASP	-	expression tag	UNP C6UB28
3	8	PRO	-	expression tag	UNP C6UB28
3	599	ALA	ARG	engineered mutation	UNP C6UB28
4	1	GLY	-	expression tag	UNP C6UB28
4	2	SER	-	expression tag	UNP C6UB28
4	3	HIS	-	expression tag	UNP C6UB28
4	4	MET	-	expression tag	UNP C6UB28
4	5	LEU	-	expression tag	UNP C6UB28
4	6	GLU	-	expression tag	UNP C6UB28
4	7	ASP	-	expression tag	UNP C6UB28
4	8	PRO	-	expression tag	UNP C6UB28
4	599	ALA	ARG	engineered mutation	UNP C6UB28

- Molecule 2 is DIMETHYL SULFOXIDE (CCD ID: DMS) (formula: C_2H_6OS).



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

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

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

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- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

- Molecule 4 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	1	3	Total 3	Na 3	0	0
4	2	4	Total 4	Na 4	0	0
4	3	5	Total 5	Na 5	0	0
4	4	4	Total 4	Na 4	0	0

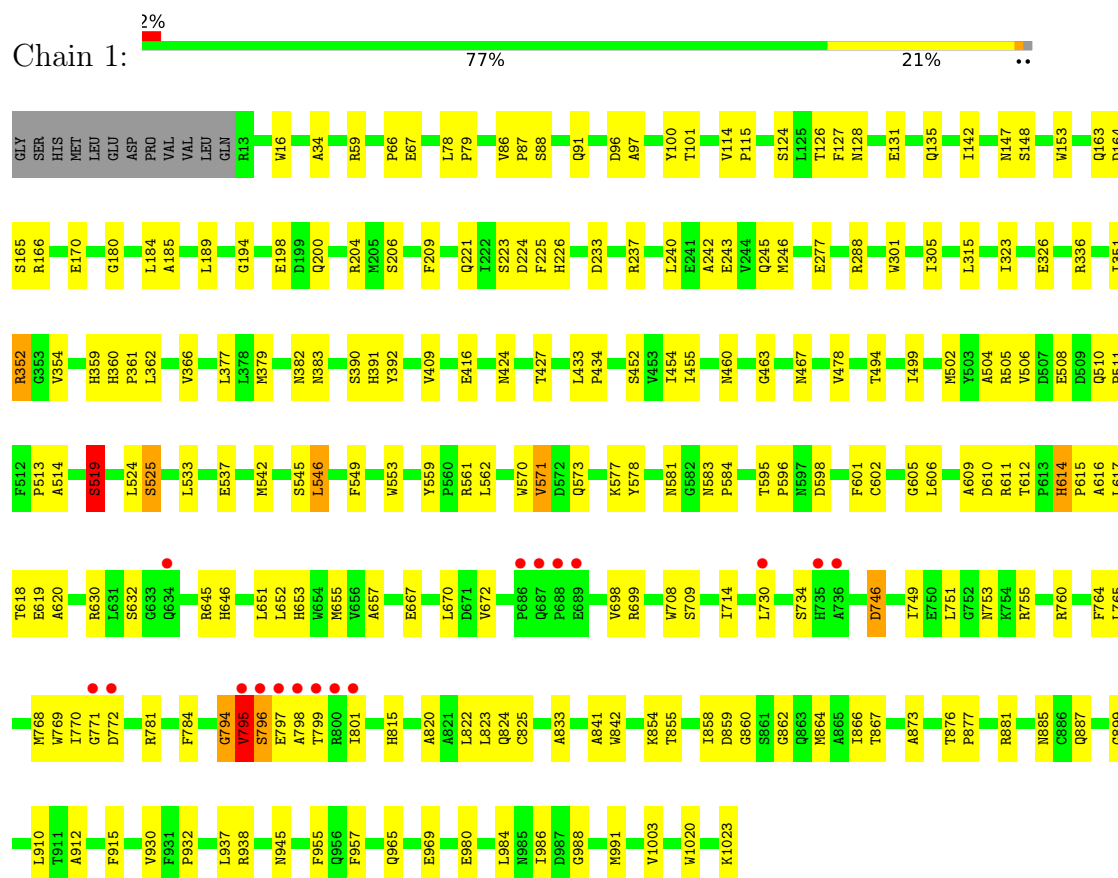
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	1	846	Total 846	O 846	0	0
5	2	821	Total 821	O 821	0	0
5	3	838	Total 838	O 838	0	0
5	4	833	Total 833	O 833	0	0

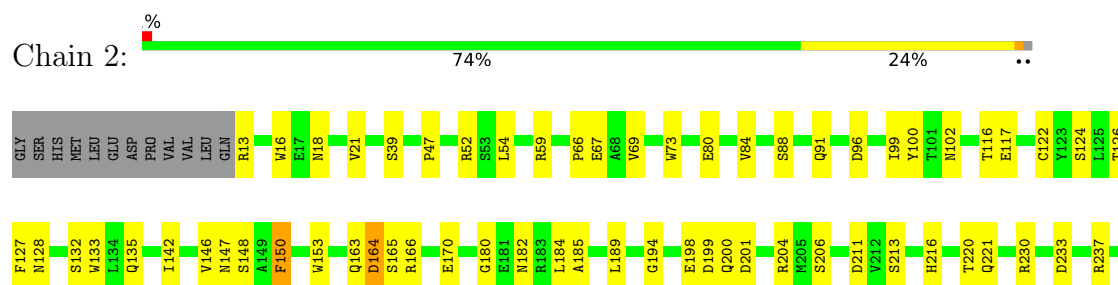
3 Residue-property plots

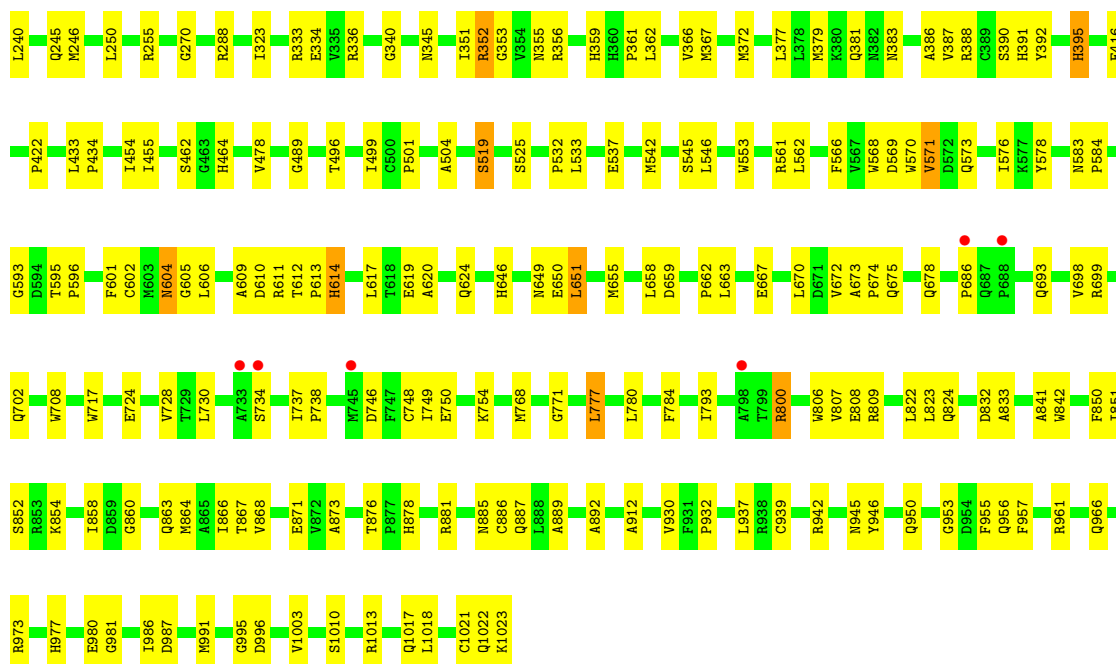
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density.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

• Molecule 1: Beta-D-galactosidase

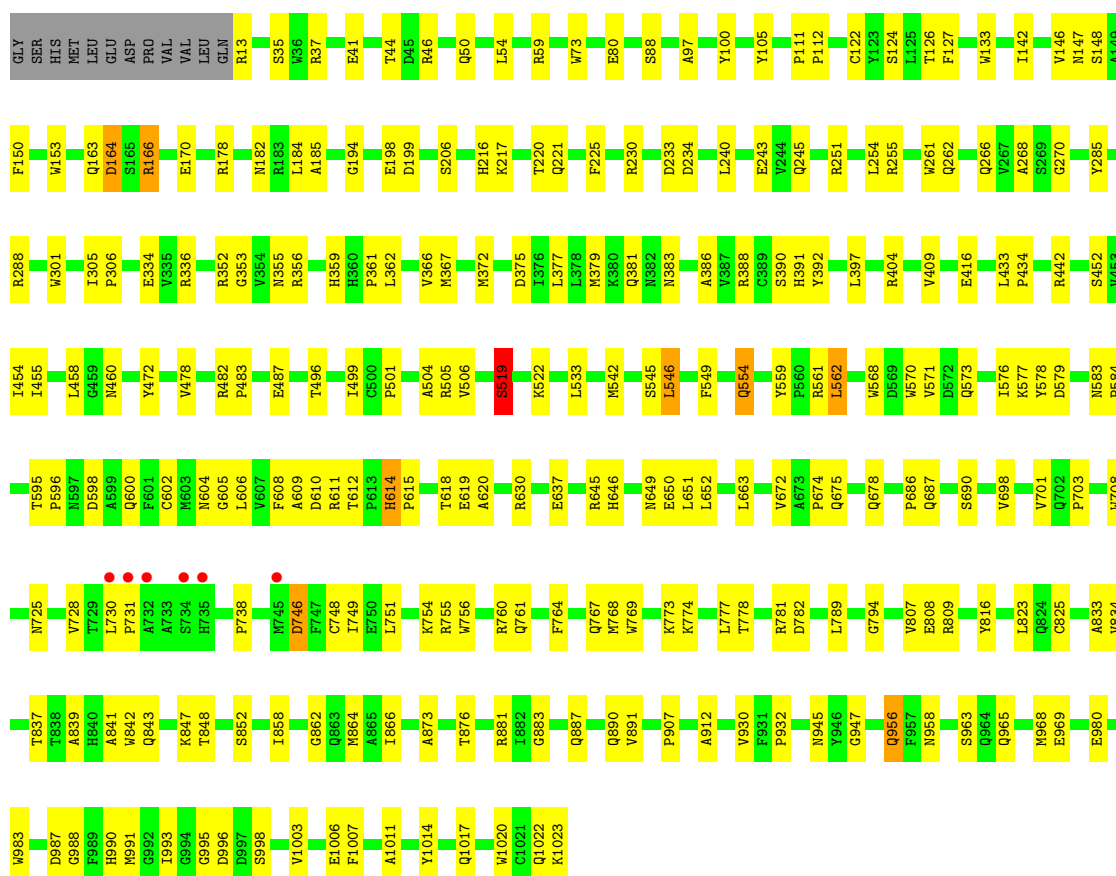
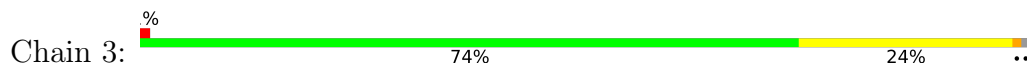


• Molecule 1: Beta-D-galactosidase

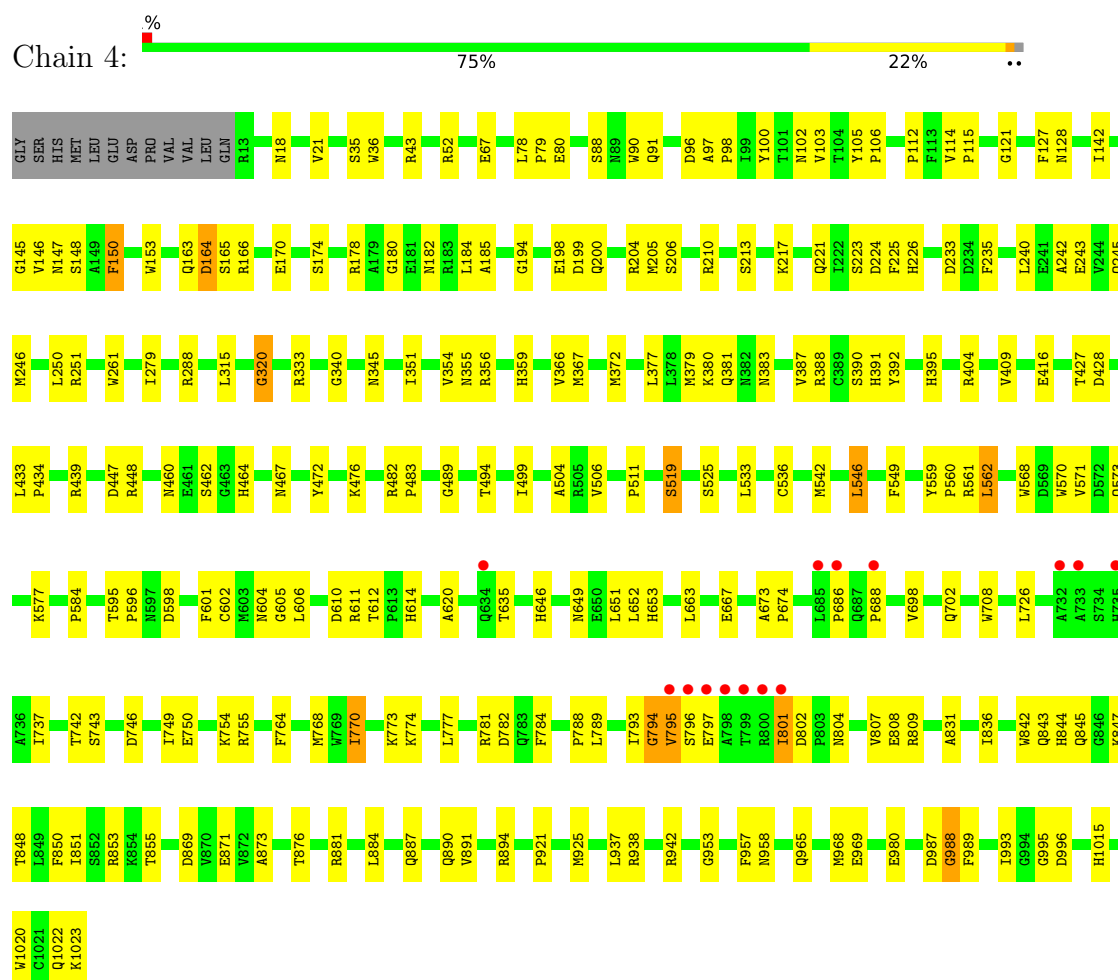




• Molecule 1: Beta-D-galactosidase



• Molecule 1: Beta-D-galactosidase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	149.27Å 167.24Å 200.61Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	86.02 – 2.50 86.02 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.7 (86.02-2.50) 99.8 (86.02-2.50)	Depositor EDS
R_{merge}	0.18	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.82 (at 2.51Å)	Xtrriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.179 , 0.235 0.175 , 0.231	Depositor DCC
R_{free} test set	2456 reflections (1.42%)	wwPDB-VP
Wilson B-factor (Å ²)	20.7	Xtrriage
Anisotropy	0.718	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 58.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	36360	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 37.76 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 4.0857e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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, 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	0.36	0/8361	0.89	26/11408 (0.2%)
1	2	0.36	0/8361	0.89	30/11408 (0.3%)
1	3	0.36	0/8361	0.89	24/11408 (0.2%)
1	4	0.37	0/8361	0.90	31/11408 (0.3%)
All	All	0.36	0/33444	0.89	111/45632 (0.2%)

There are no bond length outliers.

All (111) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	2	771	GLY	N-CA-C	-10.36	102.14	111.67
1	2	570	TRP	N-CA-C	8.88	120.66	110.97
1	4	614	HIS	N-CA-C	-8.46	99.32	110.40
1	3	570	TRP	N-CA-C	7.79	119.67	111.03
1	1	614	HIS	N-CA-C	-7.76	99.97	110.36
1	2	614	HIS	N-CA-C	-7.47	100.35	110.36
1	4	988	GLY	N-CA-C	-7.35	102.83	113.86
1	3	614	HIS	N-CA-C	-7.21	100.96	110.40
1	1	988	GLY	N-CA-C	-7.00	103.35	113.86
1	1	570	TRP	N-CA-C	6.98	118.78	111.03
1	1	746	ASP	N-CA-C	6.92	119.46	109.14
1	4	98	PRO	N-CA-C	-6.88	100.18	111.68
1	4	571	VAL	N-CA-C	6.81	118.61	108.46
1	1	97	ALA	N-CA-C	6.66	119.50	110.31
1	1	519	SER	N-CA-C	-6.63	100.22	110.10
1	4	570	TRP	N-CA-C	6.61	118.36	111.03
1	4	770	ILE	N-CA-C	-6.60	96.23	106.72
1	4	97	ALA	N-CA-C	6.60	117.97	109.72
1	4	519	SER	N-CA-C	-6.53	100.76	110.23
1	4	612	THR	N-CA-C	-6.48	101.48	109.65

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	3	988	GLY	N-CA-C	-6.43	104.21	113.86
1	4	605	GLY	N-CA-C	6.42	121.69	112.82
1	3	519	SER	N-CA-C	-6.41	100.94	110.23
1	1	506	VAL	N-CA-C	6.37	116.54	110.42
1	2	777	LEU	N-CA-C	-6.37	106.10	114.31
1	1	323	ILE	N-CA-C	-6.35	105.54	111.45
1	3	97	ALA	N-CA-C	6.35	118.07	110.07
1	4	746	ASP	N-CA-C	6.20	118.73	108.99
1	2	395	HIS	N-CA-C	-6.17	102.50	109.60
1	4	150	PHE	N-CA-C	6.15	118.20	108.79
1	3	166	ARG	N-CA-C	6.14	120.80	112.88
1	1	612	THR	N-CA-C	-6.11	101.96	109.65
1	3	605	GLY	N-CA-C	6.09	122.51	112.89
1	4	320	GLY	N-CA-C	6.09	124.02	115.43
1	4	606	LEU	N-CA-C	-6.07	105.83	113.18
1	2	593	GLY	N-CA-C	-5.98	107.43	115.21
1	4	504	ALA	N-CA-C	-5.98	100.75	110.20
1	2	519	SER	N-CA-C	-5.95	101.23	110.10
1	3	746	ASP	N-CA-C	5.95	118.01	109.14
1	1	646	HIS	N-CA-C	-5.94	101.72	110.52
1	1	571	VAL	N-CA-C	5.94	117.31	108.46
1	3	506	VAL	N-CA-C	5.91	116.09	110.53
1	1	34	ALA	N-CA-C	-5.91	107.30	114.75
1	4	777	LEU	N-CA-C	-5.88	106.45	113.97
1	3	504	ALA	N-CA-C	-5.86	101.60	110.28
1	1	352	ARG	N-CA-C	-5.85	101.45	109.71
1	2	211	ASP	N-CA-C	5.83	117.82	110.24
1	2	352	ARG	N-CA-C	-5.82	100.85	109.11
1	2	793	ILE	N-CA-C	5.80	116.45	110.82
1	1	605	GLY	N-CA-C	5.80	122.05	112.89
1	4	506	VAL	N-CA-C	5.78	115.96	110.53
1	2	955	PHE	N-CA-C	5.77	116.99	108.86
1	3	352	ARG	N-CA-C	-5.75	100.94	109.11
1	2	220	THR	N-CA-C	-5.75	100.46	109.25
1	2	116	THR	N-CA-C	-5.73	105.11	111.36
1	2	383	ASN	N-CA-C	5.71	119.45	111.74
1	2	323	ILE	N-CA-C	-5.71	106.25	111.67
1	2	233	ASP	N-CA-C	5.68	118.21	111.33
1	1	770	ILE	N-CA-C	-5.68	99.57	107.80
1	3	545	SER	N-CA-C	5.67	115.89	108.07
1	1	955	PHE	N-CA-C	5.64	116.82	108.86
1	3	404	ARG	N-CA-C	5.64	117.23	111.14

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	1	525	SER	N-CA-C	5.56	119.06	112.72
1	3	221	GLN	N-CA-C	5.48	117.30	109.14
1	4	447	ASP	N-CA-C	5.47	120.63	113.30
1	2	545	SER	N-CA-C	5.46	115.60	108.07
1	2	728	VAL	N-CA-C	-5.45	107.49	113.43
1	2	504	ALA	N-CA-C	-5.42	101.63	110.20
1	4	925	MET	N-CA-C	-5.41	106.52	113.02
1	4	448	ARG	N-CA-C	5.41	118.98	112.38
1	4	536	CYS	N-CA-C	-5.37	106.23	112.89
1	1	383	ASN	N-CA-C	5.35	118.96	111.90
1	2	562	LEU	N-CA-C	-5.34	99.81	108.41
1	1	499	ILE	N-CA-C	-5.33	99.93	107.98
1	2	221	GLN	N-CA-C	5.30	117.13	109.07
1	3	571	VAL	N-CA-C	5.30	116.36	108.46
1	4	383	ASN	N-CA-C	5.30	119.02	112.24
1	3	220	THR	N-CA-C	-5.30	99.88	108.41
1	1	409	VAL	N-CA-C	5.29	116.11	108.48
1	2	868	VAL	N-CA-C	5.27	115.48	108.11
1	1	510	GLN	CA-C-N	5.26	126.42	119.84
1	1	510	GLN	C-N-CA	5.26	126.42	119.84
1	1	351	ILE	N-CA-C	5.26	115.16	108.12
1	4	499	ILE	N-CA-C	-5.24	100.07	107.98
1	3	383	ASN	N-CA-C	5.22	118.95	112.58
1	4	635	THR	N-CA-C	5.22	117.75	109.24
1	3	777	LEU	N-CA-C	-5.21	107.47	113.88
1	2	746	ASP	N-CA-C	5.21	117.16	108.99
1	4	351	ILE	N-CA-C	5.17	115.05	108.12
1	1	504	ALA	N-CA-C	-5.15	102.06	110.20
1	4	646	HIS	N-CA-C	-5.15	102.12	110.32
1	2	351	ILE	N-CA-C	5.15	115.02	108.12
1	3	409	VAL	N-CA-C	5.14	115.99	108.53
1	4	409	VAL	N-CA-C	5.14	115.88	108.48
1	4	404	ARG	N-CA-C	5.13	118.00	111.69
1	1	221	GLN	N-CA-C	5.13	116.78	109.14
1	1	315	LEU	N-CA-C	-5.13	100.88	109.24
1	4	315	LEU	N-CA-C	-5.12	99.44	108.20
1	2	462	SER	N-CA-C	-5.12	104.85	111.71
1	2	69	VAL	N-CA-C	5.11	112.92	107.76
1	3	562	LEU	N-CA-C	-5.09	99.04	107.99
1	2	39	SER	N-CA-C	5.08	116.82	111.28
1	2	571	VAL	N-CA-C	5.07	115.95	108.45
1	4	221	GLN	N-CA-C	5.05	117.09	109.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	3	554	GLN	N-CA-C	-5.04	105.78	111.28
1	4	562	LEU	N-CA-C	-5.04	99.12	107.99
1	2	612	THR	N-CA-C	-5.04	102.09	109.50
1	2	150	PHE	N-CA-C	5.03	116.16	108.42
1	3	773	LYS	N-CA-C	5.03	117.75	109.76
1	3	956	GLN	N-CA-C	-5.03	101.91	109.85
1	3	612	THR	N-CA-C	-5.02	103.32	109.65

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	1	8119	0	7708	146	0
1	2	8119	0	7707	156	0
1	3	8119	0	7707	157	0
1	4	8119	0	7707	144	0
2	1	108	0	162	4	0
2	2	136	0	204	11	0
2	3	140	0	210	11	0
2	4	136	0	204	11	0
3	1	3	0	0	0	0
3	2	2	0	0	0	0
3	3	2	0	0	0	0
3	4	3	0	0	0	0
4	1	3	0	0	0	0
4	2	4	0	0	0	0
4	3	5	0	0	0	0
4	4	4	0	0	0	0
5	1	846	0	0	2	0
5	2	821	0	0	4	0
5	3	838	0	0	8	0
5	4	833	0	0	4	0
All	All	36360	0	31609	593	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 9.

All (593) 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)
1:4:142:ILE:HG12	1:4:170:GLU:HG2	1.37	1.05
1:3:356:ARG:HD2	1:3:379:MET:HE1	1.49	0.94
1:2:966:GLN:HE22	1:2:977:HIS:H	1.04	0.93
1:1:142:ILE:HG12	1:1:170:GLU:HG2	1.52	0.92
1:2:142:ILE:HG12	1:2:170:GLU:HG2	1.54	0.87
1:4:804:ASN:ND2	1:4:809:ARG:HH21	1.73	0.86
1:4:804:ASN:HD22	1:4:809:ARG:HH21	1.24	0.83
1:1:794:GLY:O	1:1:795:VAL:HG13	1.77	0.83
1:3:44:THR:OG1	1:3:46:ARG:HD3	1.79	0.83
1:1:630:ARG:HH12	1:1:632:SER:HB2	1.43	0.81
1:1:245:GLN:HG2	1:1:288:ARG:HG2	1.63	0.80
1:3:142:ILE:HG12	1:3:170:GLU:HG2	1.64	0.79
1:3:375:ASP:HB3	1:3:379:MET:HE3	1.67	0.77
1:1:147:ASN:HB3	1:1:206:SER:HA	1.66	0.77
1:2:966:GLN:NE2	1:2:977:HIS:H	1.83	0.75
1:1:88:SER:HA	1:1:366:VAL:HG21	1.68	0.75
1:1:655:MET:HE3	1:1:657:ALA:HB2	1.69	0.75
1:2:873:ALA:O	1:2:876:THR:HG22	1.85	0.74
1:2:147:ASN:HB3	1:2:206:SER:HA	1.70	0.73
1:3:579:ASP:OD2	1:3:583:ASN:HB2	1.88	0.73
1:2:737:ILE:HD12	1:2:738:PRO:HD2	1.68	0.73
1:1:615:PRO:O	1:1:618:THR:HG22	1.89	0.73
1:4:568:TRP:HE1	1:4:604:ASN:HD22	1.38	0.72
1:2:245:GLN:HG2	1:2:288:ARG:HG2	1.72	0.71
1:4:245:GLN:HG2	1:4:288:ARG:HG2	1.72	0.71
1:1:546:LEU:HG	1:1:616:ALA:HB1	1.73	0.71
1:4:147:ASN:HB3	1:4:206:SER:HA	1.73	0.71
1:4:372:MET:HE1	1:4:395:HIS:HB3	1.72	0.69
1:1:730:LEU:HD21	1:2:823:LEU:O	1.91	0.69
1:1:965:GLN:O	1:1:969:GLU:HG3	1.93	0.69
1:3:416:GLU:HG3	1:3:460:ASN:O	1.93	0.69
1:3:755:ARG:HB3	1:3:769:TRP:HB2	1.75	0.69
1:4:250:LEU:O	1:4:251:ARG:HG2	1.93	0.69
1:2:966:GLN:HE22	1:2:977:HIS:N	1.87	0.68
1:2:748:CYS:C	1:2:749:ILE:HD12	2.18	0.68
1:2:379:MET:HE1	1:2:387:VAL:HB	1.76	0.68
1:2:991:MET:HE3	1:2:1003:VAL:HG21	1.76	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:3:652:LEU:HD11	1:3:698:VAL:HB	1.76	0.67
1:4:377:LEU:HD22	1:4:708:TRP:HA	1.76	0.67
1:3:88:SER:HA	1:3:366:VAL:HG21	1.75	0.67
1:3:147:ASN:HB3	1:3:206:SER:HA	1.75	0.67
1:2:270:GLY:HA3	2:2:8405:DMS:H23	1.77	0.67
1:3:615:PRO:O	1:3:618:THR:HG22	1.94	0.67
1:4:320:GLY:HA2	2:4:8406:DMS:H22	1.75	0.67
1:3:245:GLN:HG2	1:3:288:ARG:HG2	1.74	0.67
1:2:356:ARG:HH22	1:2:367:MET:HE2	1.60	0.67
1:3:268:ALA:HA	2:3:8603:DMS:H23	1.77	0.66
1:4:355:ASN:OD1	1:4:388:ARG:HD3	1.94	0.66
1:2:336:ARG:HB3	2:2:8409:DMS:H23	1.76	0.66
1:1:630:ARG:NH1	1:1:632:SER:HB2	2.10	0.66
1:2:737:ILE:HD13	1:2:832:ASP:HA	1.77	0.66
1:4:174:SER:HB3	2:4:8705:DMS:H12	1.78	0.65
1:4:598:ASP:OD1	1:4:797:GLU:HA	1.96	0.65
1:3:651:LEU:O	1:3:651:LEU:HD12	1.96	0.65
1:4:651:LEU:HD12	1:4:651:LEU:O	1.97	0.65
1:4:804:ASN:HD22	1:4:809:ARG:NH2	1.94	0.65
1:1:653:HIS:CD2	1:1:667:GLU:HG2	2.32	0.64
1:2:127:PHE:HE2	1:2:184:LEU:HG	1.61	0.64
1:3:454:ILE:HG13	1:3:455:ILE:HG13	1.79	0.64
1:4:595:THR:HA	1:4:596:PRO:C	2.23	0.64
1:1:598:ASP:OD1	1:1:797:GLU:HA	1.96	0.64
1:1:277:GLU:CD	1:1:277:GLU:H	2.06	0.63
1:2:372:MET:HE1	1:2:395:HIS:HB3	1.80	0.63
1:3:843:GLN:HG2	1:3:848:THR:HA	1.81	0.63
1:4:233:ASP:HA	2:4:8417:DMS:H11	1.79	0.63
1:3:240:LEU:HD23	1:3:240:LEU:C	2.24	0.62
1:4:88:SER:HA	1:4:366:VAL:HG21	1.81	0.62
1:1:542:MET:HE3	1:1:601:PHE:HA	1.81	0.62
1:3:991:MET:HE3	1:3:1003:VAL:HG21	1.80	0.62
1:3:890:GLN:HG2	1:3:891:VAL:N	2.14	0.62
1:2:620:ALA:O	1:2:624:GLN:HG3	2.00	0.62
1:2:91:GLN:HG3	1:2:96:ASP:OD1	1.99	0.62
1:2:699:ARG:HG2	1:2:717:TRP:HB3	1.82	0.61
1:4:801:ILE:N	1:4:801:ILE:HD12	2.13	0.61
1:4:36:TRP:HH2	2:4:8404:DMS:H23	1.65	0.61
1:3:568:TRP:HE1	1:3:604:ASN:HD22	1.48	0.61
1:4:427:THR:HG21	1:4:462:SER:HB3	1.82	0.61
1:4:793:ILE:HG22	1:4:795:VAL:HG22	1.83	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:3:251:ARG:HG3	1:3:251:ARG:HH11	1.66	0.61
1:2:124:SER:HA	1:2:184:LEU:O	2.01	0.61
1:4:794:GLY:O	1:4:795:VAL:HG13	2.01	0.60
1:1:382:ASN:ND2	1:1:617:LEU:HD21	2.17	0.60
1:1:945:ASN:HB3	1:1:1023:LYS:HE3	1.82	0.60
1:1:873:ALA:O	1:1:876:THR:HG22	2.01	0.60
1:3:864:MET:HE3	1:3:866:ILE:HD11	1.82	0.60
1:2:18:ASN:ND2	1:2:21:VAL:HG23	2.17	0.60
1:4:380:LYS:O	2:4:8403:DMS:H21	2.00	0.60
1:2:377:LEU:HD22	1:2:708:TRP:HA	1.84	0.60
1:1:765:LEU:HD21	1:1:768:MET:HE2	1.85	0.59
1:2:663:LEU:HD21	1:2:686:PRO:HG2	1.83	0.59
1:1:937:LEU:O	1:1:938:ARG:HD2	2.02	0.59
1:2:754:LYS:HE2	1:2:1022:GLN:NE2	2.18	0.59
1:2:777:LEU:HG	1:2:889:ALA:HA	1.84	0.59
1:2:942:ARG:HA	1:2:953:GLY:O	2.03	0.59
1:1:801:ILE:N	1:1:801:ILE:HD12	2.17	0.59
1:2:986:ILE:HD13	1:2:1018:LEU:HD13	1.85	0.59
1:3:356:ARG:HD2	1:3:379:MET:CE	2.30	0.59
1:3:749:ILE:HD12	1:3:749:ILE:N	2.17	0.59
1:4:142:ILE:HG12	1:4:170:GLU:CG	2.25	0.59
1:3:767:GLN:NE2	1:3:774:LYS:HD3	2.17	0.59
1:4:651:LEU:HD12	1:4:651:LEU:C	2.28	0.59
1:1:801:ILE:HD12	1:1:801:ILE:H	1.66	0.59
1:2:672:VAL:HG22	1:2:678:GLN:HB2	1.83	0.59
1:1:746:ASP:HA	1:1:760:ARG:HG3	1.83	0.58
1:1:377:LEU:HD22	1:1:708:TRP:HA	1.85	0.58
1:3:672:VAL:HG22	1:3:678:GLN:HB2	1.85	0.58
1:2:88:SER:HA	1:2:366:VAL:HG21	1.86	0.58
1:2:864:MET:HE3	1:2:866:ILE:HD11	1.85	0.58
1:1:545:SER:O	1:1:546:LEU:HB2	2.02	0.58
1:3:270:GLY:HA3	2:3:8405:DMS:H23	1.85	0.58
1:3:377:LEU:O	1:3:381:GLN:HG3	2.02	0.57
1:2:610:ASP:O	1:2:611:ARG:HB2	2.04	0.57
1:4:768:MET:HE1	1:4:1020:TRP:CZ2	2.39	0.57
2:2:8420:DMS:H13	1:3:478:VAL:HB	1.86	0.57
1:3:533:LEU:C	1:3:533:LEU:HD23	2.29	0.57
1:3:579:ASP:CG	1:3:583:ASN:HB2	2.29	0.57
1:1:887:GLN:NE2	1:1:980:GLU:O	2.38	0.57
1:2:595:THR:HA	1:2:596:PRO:C	2.28	0.57
1:1:494:THR:HA	2:1:8607:DMS:H12	1.85	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:3:41:GLU:HG2	1:3:46:ARG:NH2	2.19	0.57
1:2:194:GLY:O	1:2:198:GLU:HG3	2.05	0.57
1:1:749:ILE:N	1:1:749:ILE:HD12	2.19	0.57
1:1:764:PHE:CE1	1:1:781:ARG:NH1	2.72	0.57
1:2:127:PHE:CE2	1:2:184:LEU:HG	2.40	0.57
1:3:890:GLN:OE1	1:3:947:GLY:HA3	2.05	0.56
1:4:494:THR:HA	2:4:8615:DMS:H11	1.86	0.56
1:4:965:GLN:O	1:4:969:GLU:HG3	2.06	0.56
1:3:768:MET:HE1	1:3:1020:TRP:CZ2	2.41	0.56
1:4:416:GLU:HG3	1:4:460:ASN:O	2.05	0.56
1:1:508:GLU:OE2	2:1:8407:DMS:H11	2.06	0.56
1:4:153:TRP:HB2	1:4:185:ALA:HB3	1.88	0.56
1:4:788:PRO:HD2	1:4:968:MET:HG3	1.87	0.56
1:3:965:GLN:O	1:3:969:GLU:HG3	2.06	0.56
1:4:433:LEU:HB3	1:4:434:PRO:HD3	1.87	0.56
1:2:13:ARG:HG3	1:3:13:ARG:NH1	2.21	0.55
1:2:619:GLU:HA	1:2:912:ALA:HB2	1.88	0.55
1:3:823:LEU:HD11	1:3:841:ALA:HB2	1.89	0.55
1:2:153:TRP:HB2	1:2:185:ALA:HB3	1.87	0.55
1:3:748:CYS:C	1:3:749:ILE:HD12	2.31	0.55
1:4:546:LEU:HA	5:4:4130:HOH:O	2.07	0.55
1:2:573:GLN:HB2	1:2:602:CYS:O	2.06	0.55
1:4:942:ARG:HA	1:4:953:GLY:O	2.07	0.54
1:2:651:LEU:HD11	1:2:667:GLU:OE1	2.06	0.54
1:4:91:GLN:HG3	1:4:96:ASP:OD1	2.07	0.54
1:3:728:VAL:HG12	1:4:851:ILE:HD11	1.89	0.54
1:3:356:ARG:HH22	1:3:367:MET:HE2	1.73	0.54
1:3:573:GLN:HB2	1:3:602:CYS:O	2.07	0.54
1:4:894:ARG:NH1	1:4:921:PRO:HG3	2.23	0.54
1:2:47:PRO:HA	2:2:8606:DMS:H12	1.89	0.54
1:4:533:LEU:C	1:4:533:LEU:HD23	2.32	0.54
1:2:930:VAL:O	1:2:932:PRO:HD3	2.08	0.54
1:3:59:ARG:HB2	1:3:124:SER:OG	2.07	0.54
1:4:340:GLY:O	1:4:561:ARG:HG2	2.08	0.54
1:4:356:ARG:HH22	1:4:367:MET:HE2	1.73	0.54
1:2:784:PHE:HD1	2:2:8611:DMS:H13	1.73	0.53
1:1:652:LEU:HD11	1:1:698:VAL:HB	1.90	0.53
1:2:166:ARG:HG3	1:2:392:TYR:HB2	1.89	0.53
1:4:784:PHE:HA	1:4:881:ARG:O	2.08	0.53
1:1:237:ARG:HB3	1:1:237:ARG:NH1	2.24	0.53
1:2:356:ARG:NH2	1:2:367:MET:HE2	2.24	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:433:LEU:HB3	1:1:434:PRO:HD3	1.90	0.53
1:2:823:LEU:HD11	1:2:841:ALA:HB2	1.91	0.53
1:4:372:MET:HE2	5:4:4195:HOH:O	2.08	0.53
1:2:881:ARG:HE	1:2:987:ASP:CG	2.17	0.53
1:3:334:GLU:OE1	1:3:336:ARG:NH1	2.41	0.53
1:1:795:VAL:C	1:1:797:GLU:H	2.17	0.52
1:2:542:MET:HE3	1:2:601:PHE:HA	1.90	0.52
1:3:306:PRO:HD2	2:3:8612:DMS:H23	1.91	0.52
1:2:945:ASN:OD1	1:2:950:GLN:HG3	2.08	0.52
1:1:100:TYR:CE1	1:1:602:CYS:HB3	2.45	0.52
1:3:703:PRO:HG2	2:3:8425:DMS:H11	1.90	0.52
1:4:549:PHE:CE2	1:4:620:ALA:HA	2.44	0.52
1:4:802:ASP:O	1:4:808:GLU:HG3	2.09	0.52
1:2:863:GLN:HG2	1:2:1021:CYS:CB	2.39	0.52
1:2:100:TYR:CE1	1:2:602:CYS:HB3	2.44	0.52
1:4:770:ILE:HG21	1:4:1022:GLN:HE22	1.75	0.52
1:1:87:PRO:HB2	1:1:209:PHE:C	2.35	0.52
1:1:533:LEU:HD23	1:1:533:LEU:C	2.34	0.52
1:3:230:ARG:HD3	5:3:7199:HOH:O	2.10	0.52
1:4:749:ILE:N	1:4:749:ILE:HD12	2.25	0.52
1:2:454:ILE:HG13	1:2:455:ILE:HG13	1.91	0.52
1:4:148:SER:HA	1:4:165:SER:OG	2.10	0.52
1:2:734:SER:HB3	1:2:860:GLY:HA3	1.92	0.52
1:2:646:HIS:HB3	5:2:6063:HOH:O	2.09	0.51
1:1:784:PHE:HA	1:1:881:ARG:O	2.09	0.51
1:4:686:PRO:O	1:4:688:PRO:HD3	2.11	0.51
1:1:542:MET:CE	1:1:601:PHE:HA	2.40	0.51
1:3:251:ARG:HG3	1:3:251:ARG:NH1	2.24	0.51
1:3:839:ALA:HA	1:3:852:SER:O	2.10	0.51
1:1:301:TRP:CH2	1:1:452:SER:HA	2.45	0.51
1:4:937:LEU:O	1:4:938:ARG:HD2	2.10	0.51
1:3:362:LEU:HD23	1:3:576:ILE:HB	1.92	0.51
1:4:873:ALA:O	1:4:876:THR:HG22	2.10	0.51
1:1:153:TRP:HB2	1:1:185:ALA:HB3	1.93	0.51
1:1:336:ARG:HB3	2:1:8409:DMS:O	2.11	0.51
1:1:424:ASN:OD1	1:4:279:ILE:HD11	2.11	0.51
1:1:833:ALA:HB1	1:1:858:ILE:O	2.11	0.51
1:2:542:MET:HA	1:2:604:ASN:HA	1.92	0.51
1:3:499:ILE:HG22	1:3:501:PRO:HD3	1.93	0.51
1:4:250:LEU:C	1:4:251:ARG:HG2	2.36	0.51
1:2:533:LEU:C	1:2:533:LEU:HD23	2.36	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:768:MET:HE1	1:1:1020:TRP:CZ2	2.45	0.50
1:2:433:LEU:HB3	1:2:434:PRO:HD3	1.94	0.50
1:1:755:ARG:HB2	1:1:769:TRP:HB2	1.94	0.50
1:2:499:ILE:HG22	1:2:501:PRO:HD3	1.94	0.50
1:2:362:LEU:HD23	1:2:576:ILE:HB	1.93	0.50
1:2:478:VAL:HG12	2:3:8420:DMS:H13	1.94	0.50
1:3:458:LEU:HD11	1:3:472:TYR:HB2	1.93	0.50
1:2:377:LEU:O	1:2:381:GLN:HG3	2.11	0.50
1:2:863:GLN:HG2	1:2:1021:CYS:HB3	1.92	0.50
1:3:595:THR:HA	1:3:596:PRO:C	2.37	0.50
1:3:738:PRO:HB2	1:3:834:VAL:HG23	1.94	0.50
1:1:559:TYR:HB2	1:1:562:LEU:HD12	1.93	0.50
1:2:132:SER:HA	1:2:135:GLN:CD	2.37	0.50
1:2:146:VAL:HG11	1:2:150:PHE:CD2	2.46	0.50
1:3:610:ASP:O	1:3:611:ARG:HB2	2.12	0.50
1:1:127:PHE:CE2	1:1:184:LEU:HG	2.47	0.49
1:3:649:ASN:O	1:3:650:GLU:HG3	2.12	0.49
1:4:428:ASP:O	2:4:8611:DMS:H22	2.12	0.49
1:4:673:ALA:HB1	1:4:674:PRO:HD2	1.93	0.49
1:3:646:HIS:HB3	5:3:6063:HOH:O	2.13	0.49
1:4:225:PHE:HA	1:4:243:GLU:O	2.12	0.49
1:1:734:SER:HB2	1:1:860:GLY:HA3	1.95	0.49
1:2:353:GLY:HA2	1:2:386:ALA:O	2.13	0.49
1:2:464:HIS:HB2	1:2:489:GLY:HA3	1.93	0.49
1:4:568:TRP:HE1	1:4:604:ASN:ND2	2.08	0.49
1:4:958:ASN:C	1:4:958:ASN:OD1	2.55	0.49
1:1:578:TYR:CE2	1:1:584:PRO:HB3	2.47	0.49
1:2:100:TYR:CZ	1:2:602:CYS:HB3	2.48	0.49
1:2:237:ARG:HG3	1:2:237:ARG:HH11	1.77	0.49
1:2:655:MET:HE1	1:2:662:PRO:HB3	1.94	0.49
1:2:542:MET:CE	1:2:601:PHE:HA	2.43	0.49
1:3:126:THR:HA	1:3:182:ASN:O	2.12	0.49
1:3:1011:ALA:HB3	1:3:1014:TYR:CZ	2.48	0.49
1:4:573:GLN:HB2	1:4:602:CYS:O	2.11	0.49
1:4:651:LEU:HD11	1:4:653:HIS:CE1	2.48	0.49
1:4:200:GLN:HG2	1:4:391:HIS:HB2	1.94	0.48
1:4:843:GLN:HA	1:4:847:LYS:O	2.13	0.48
1:1:581:ASN:HB2	1:1:583:ASN:OD1	2.12	0.48
1:2:822:LEU:HD11	1:2:824:GLN:O	2.13	0.48
1:4:887:GLN:NE2	1:4:980:GLU:O	2.42	0.48
1:1:699:ARG:HD3	1:1:714:ILE:HD13	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:133:TRP:CE3	1:2:216:HIS:HB2	2.48	0.48
1:2:961:ARG:NE	1:2:981:GLY:O	2.43	0.48
1:4:764:PHE:CD1	1:4:781:ARG:HG2	2.48	0.48
1:2:352:ARG:HG2	1:2:553:TRP:CH2	2.48	0.48
1:2:708:TRP:CZ2	2:2:8403:DMS:H11	2.48	0.48
1:4:742:THR:HG22	1:4:743:SER:N	2.27	0.48
1:1:245:GLN:HG2	1:1:288:ARG:CG	2.38	0.48
1:4:18:ASN:ND2	1:4:21:VAL:HG23	2.29	0.48
1:4:737:ILE:HD12	1:4:831:ALA:O	2.14	0.48
1:1:708:TRP:CE3	1:1:709:SER:HB3	2.47	0.48
1:3:336:ARG:HB3	2:3:8409:DMS:H23	1.95	0.48
1:4:114:VAL:HB	1:4:115:PRO:HD2	1.94	0.48
1:4:128:ASN:HB3	1:4:180:GLY:O	2.13	0.48
1:2:255:ARG:HG2	1:2:255:ARG:HH11	1.79	0.48
1:2:663:LEU:C	1:2:663:LEU:HD23	2.39	0.48
1:4:782:ASP:HA	1:4:884:LEU:HD23	1.95	0.48
1:3:749:ILE:HB	1:3:756:TRP:HB2	1.95	0.48
1:3:768:MET:HE1	1:3:1020:TRP:CH2	2.48	0.48
1:3:608:PHE:CE1	1:3:614:HIS:HD2	2.32	0.47
1:4:145:GLY:N	1:4:210:ARG:HB2	2.29	0.47
1:1:755:ARG:HD2	5:1:6193:HOH:O	2.13	0.47
1:1:854:LYS:HA	1:1:867:THR:O	2.14	0.47
1:3:833:ALA:HB1	1:3:858:ILE:O	2.14	0.47
1:4:1022:GLN:O	1:4:1023:LYS:HB2	2.14	0.47
1:2:613:PRO:HB3	1:2:617:LEU:HD23	1.94	0.47
1:2:1013:ARG:O	2:2:8612:DMS:H22	2.14	0.47
1:2:355:ASN:OD1	1:2:388:ARG:HD3	2.14	0.47
1:3:251:ARG:HD3	2:3:8416:DMS:H22	1.96	0.47
1:3:674:PRO:O	1:3:675:GLN:HB2	2.13	0.47
1:3:746:ASP:HA	1:3:760:ARG:HG3	1.95	0.47
1:4:91:GLN:OE1	1:4:205:MET:HB3	2.13	0.47
1:1:549:PHE:CE2	1:1:620:ALA:HA	2.49	0.47
1:2:496:THR:O	2:2:8610:DMS:H13	2.15	0.47
1:3:651:LEU:HD12	1:3:651:LEU:C	2.40	0.47
1:1:795:VAL:C	1:1:797:GLU:N	2.72	0.47
1:4:194:GLY:O	1:4:198:GLU:HG3	2.14	0.47
1:4:843:GLN:HG2	1:4:848:THR:HA	1.96	0.47
1:1:454:ILE:HG13	1:1:455:ILE:HG13	1.97	0.47
1:4:146:VAL:HG11	1:4:150:PHE:CD2	2.49	0.47
1:1:240:LEU:C	1:1:240:LEU:HD23	2.38	0.47
1:2:651:LEU:C	1:2:651:LEU:HD23	2.40	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:698:VAL:O	1:2:717:TRP:HA	2.14	0.47
1:3:559:TYR:HB2	1:3:562:LEU:HD12	1.97	0.47
1:3:794:GLY:HA2	1:3:998:SER:O	2.15	0.47
1:1:502:MET:HB2	1:1:537:GLU:HB2	1.95	0.47
1:2:750:GLU:HB3	5:2:5177:HOH:O	2.13	0.47
1:3:166:ARG:HG3	1:3:392:TYR:HB2	1.97	0.47
1:3:809:ARG:HD2	5:3:4651:HOH:O	2.14	0.47
1:4:750:GLU:OE2	1:4:755:ARG:HD3	2.15	0.47
1:2:334:GLU:OE1	1:2:336:ARG:NH1	2.46	0.46
1:3:549:PHE:CE2	1:3:620:ALA:HA	2.49	0.46
1:3:816:TYR:CE2	1:3:968:MET:HE1	2.50	0.46
1:4:164:ASP:OD2	1:4:439:ARG:HG2	2.15	0.46
1:4:246:MET:SD	1:4:246:MET:C	2.99	0.46
1:4:890:GLN:HG3	1:4:891:VAL:N	2.30	0.46
1:3:377:LEU:HD22	1:3:708:TRP:HA	1.97	0.46
1:4:354:VAL:CG1	1:4:379:MET:HE3	2.46	0.46
1:1:200:GLN:HG2	1:1:391:HIS:HB2	1.97	0.46
1:1:573:GLN:HB2	1:1:602:CYS:O	2.16	0.46
1:2:986:ILE:HG23	1:2:986:ILE:O	2.15	0.46
1:3:73:TRP:CE2	1:3:122:CYS:HB3	2.51	0.46
1:2:390:SER:HA	1:2:391:HIS:HA	1.74	0.46
1:2:833:ALA:HB1	1:2:858:ILE:O	2.15	0.46
1:4:79:PRO:HG2	1:4:80:GLU:OE2	2.16	0.46
1:4:390:SER:HA	1:4:391:HIS:HA	1.76	0.46
1:1:114:VAL:HB	1:1:115:PRO:HD2	1.97	0.46
1:1:361:PRO:HB3	1:1:609:ALA:HB1	1.97	0.46
1:2:148:SER:HA	1:2:165:SER:OG	2.16	0.46
1:2:336:ARG:CB	2:2:8409:DMS:H23	2.45	0.46
1:2:754:LYS:HE2	1:2:1022:GLN:HE21	1.80	0.46
1:3:100:TYR:CZ	1:3:602:CYS:HB3	2.51	0.46
1:1:166:ARG:HG3	1:1:392:TYR:HB2	1.97	0.46
1:1:930:VAL:O	1:1:932:PRO:HD3	2.15	0.46
1:2:199:ASP:C	1:2:416:GLU:HG2	2.41	0.46
1:3:35:SER:HB2	1:3:217:LYS:HD2	1.97	0.46
1:4:333:ARG:HA	1:4:345:ASN:OD1	2.15	0.46
1:4:601:PHE:CD1	1:4:796:SER:HA	2.51	0.46
1:1:359:HIS:CD2	1:1:573:GLN:HA	2.51	0.46
1:1:768:MET:HE1	1:1:1020:TRP:CH2	2.51	0.46
1:2:99:ILE:O	1:2:204:ARG:HG2	2.16	0.46
1:2:230:ARG:HD3	5:2:7562:HOH:O	2.14	0.46
1:1:237:ARG:HB3	1:1:237:ARG:HH11	1.80	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:546:LEU:HD22	1:1:546:LEU:HA	1.80	0.45
1:3:372:MET:HE1	1:3:397:LEU:HB3	1.98	0.45
1:3:433:LEU:N	1:3:434:PRO:CD	2.78	0.45
1:3:554:GLN:HG3	5:3:7553:HOH:O	2.15	0.45
1:4:246:MET:HE3	1:4:250:LEU:HD23	1.98	0.45
1:1:797:GLU:C	1:1:799:THR:H	2.24	0.45
1:2:166:ARG:HG3	1:2:392:TYR:CB	2.46	0.45
1:3:598:ASP:OD1	2:3:8602:DMS:H21	2.16	0.45
1:1:101:THR:HG23	1:1:204:ARG:NH2	2.31	0.45
1:1:147:ASN:HA	1:1:148:SER:HA	1.63	0.45
1:1:194:GLY:O	1:1:198:GLU:HG3	2.16	0.45
1:1:799:THR:HG22	1:1:799:THR:O	2.16	0.45
1:3:245:GLN:HG2	1:3:288:ARG:CG	2.46	0.45
1:2:245:GLN:HG2	1:2:288:ARG:CG	2.43	0.45
1:2:887:GLN:NE2	1:2:980:GLU:O	2.49	0.45
1:2:892:ALA:HB3	1:2:946:TYR:CE1	2.51	0.45
1:3:496:THR:O	2:3:8606:DMS:H21	2.16	0.45
1:4:90:TRP:CZ3	1:4:121:GLY:HA3	2.52	0.45
1:4:163:GLN:O	1:4:164:ASP:HB3	2.17	0.45
1:4:235:PHE:CE2	1:4:345:ASN:HA	2.52	0.45
1:2:16:TRP:CG	1:2:189:LEU:HD13	2.52	0.45
1:2:66:PRO:HG2	1:2:67:GLU:OE1	2.17	0.45
1:4:881:ARG:HE	1:4:987:ASP:CG	2.24	0.45
1:1:66:PRO:HG2	1:1:67:GLU:OE1	2.15	0.45
1:2:163:GLN:O	1:2:164:ASP:HB3	2.17	0.45
1:3:54:LEU:O	2:3:8408:DMS:H12	2.17	0.45
1:3:233:ASP:OD1	1:3:234:ASP:N	2.49	0.45
1:1:305:ILE:HD11	1:1:645:ARG:HB3	1.98	0.45
1:1:822:LEU:HD11	1:1:824:GLN:O	2.17	0.45
1:2:52:ARG:O	1:2:213:SER:HB2	2.16	0.45
1:2:537:GLU:HA	1:2:566:PHE:O	2.17	0.45
1:2:780:LEU:HA	1:2:886:CYS:HB3	1.99	0.45
1:3:789:LEU:HD11	1:3:993:ILE:HG22	1.97	0.45
1:3:153:TRP:HB2	1:3:185:ALA:HB3	1.99	0.44
1:3:356:ARG:HH22	1:3:367:MET:CE	2.30	0.44
1:3:505:ARG:O	1:3:519:SER:HA	2.16	0.44
1:1:619:GLU:HA	1:1:912:ALA:HB2	1.99	0.44
1:1:390:SER:HA	1:1:391:HIS:HA	1.72	0.44
1:1:505:ARG:O	1:1:519:SER:HA	2.17	0.44
1:2:246:MET:HE3	1:2:250:LEU:HD23	1.98	0.44
1:4:649:ASN:HA	2:4:1024:DMS:H22	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:3:194:GLY:O	1:3:198:GLU:HG3	2.18	0.44
1:3:433:LEU:HB3	1:3:434:PRO:HD3	1.99	0.44
1:3:1006:GLU:HG2	1:3:1007:PHE:CD1	2.53	0.44
1:4:100:TYR:CE1	1:4:602:CYS:HB3	2.51	0.44
1:1:223:SER:O	1:1:224:ASP:HB2	2.17	0.44
1:1:478:VAL:HB	2:4:8611:DMS:H12	1.98	0.44
1:2:59:ARG:HB2	1:2:124:SER:OG	2.18	0.44
1:2:84:VAL:HG12	2:2:8414:DMS:H13	1.99	0.44
1:1:91:GLN:HG3	1:1:96:ASP:OD1	2.17	0.44
1:2:674:PRO:O	1:2:675:GLN:HB2	2.18	0.44
1:4:472:TYR:O	1:4:476:LYS:HG2	2.17	0.44
1:1:795:VAL:O	1:1:797:GLU:N	2.50	0.44
1:2:117:GLU:OE1	1:2:117:GLU:HA	2.18	0.44
1:4:379:MET:HE1	1:4:387:VAL:HB	2.00	0.44
1:1:525:SER:O	1:2:561:ARG:HD3	2.17	0.44
1:1:571:VAL:CG2	1:1:609:ALA:HA	2.48	0.44
1:3:147:ASN:HA	1:3:148:SER:HA	1.70	0.44
1:3:225:PHE:HA	1:3:243:GLU:O	2.18	0.44
1:3:353:GLY:HA2	1:3:386:ALA:O	2.18	0.44
1:3:390:SER:HA	1:3:391:HIS:HA	1.78	0.44
1:4:995:GLY:O	1:4:996:ASP:C	2.61	0.44
1:1:354:VAL:CG1	1:1:379:MET:HE3	2.48	0.44
1:2:102:ASN:OD1	1:2:201:ASP:HB2	2.18	0.44
1:2:606:LEU:O	1:2:614:HIS:HB2	2.17	0.44
1:3:690:SER:HB2	5:3:5032:HOH:O	2.18	0.44
1:4:223:SER:O	1:4:224:ASP:HB2	2.17	0.44
1:4:245:GLN:HG2	1:4:288:ARG:CG	2.46	0.44
1:1:128:ASN:HA	1:1:180:GLY:O	2.18	0.43
1:1:163:GLN:O	1:1:164:ASP:HB3	2.18	0.43
1:1:910:LEU:C	1:1:910:LEU:HD12	2.43	0.43
1:3:606:LEU:O	1:3:614:HIS:HB2	2.18	0.43
1:2:945:ASN:HB3	1:2:1023:LYS:HE2	1.99	0.43
1:4:773:LYS:HE3	1:4:774:LYS:O	2.19	0.43
1:1:226:HIS:O	1:1:242:ALA:HA	2.19	0.43
1:1:651:LEU:C	1:1:651:LEU:HD12	2.42	0.43
1:1:801:ILE:H	1:1:801:ILE:CD1	2.29	0.43
1:1:855:THR:OG1	1:1:867:THR:HB	2.19	0.43
1:1:991:MET:HE3	1:1:1003:VAL:HG21	2.00	0.43
1:2:73:TRP:CE2	1:2:122:CYS:HB3	2.52	0.43
1:2:245:GLN:CG	1:2:288:ARG:HG2	2.45	0.43
1:2:939:CYS:SG	1:2:956:GLN:HG3	2.58	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:3:359:HIS:CD2	1:3:573:GLN:HA	2.53	0.43
1:3:546:LEU:HA	5:3:4130:HOH:O	2.17	0.43
1:3:764:PHE:CE1	1:3:781:ARG:NH1	2.86	0.43
1:4:127:PHE:CE2	1:4:184:LEU:HG	2.53	0.43
1:1:277:GLU:CD	1:1:277:GLU:N	2.74	0.43
1:4:67:GLU:OE1	1:4:67:GLU:N	2.45	0.43
1:4:359:HIS:CD2	1:4:573:GLN:HA	2.53	0.43
1:1:815:HIS:HE1	1:1:877:PRO:O	2.00	0.43
1:1:864:MET:HE3	1:1:866:ILE:HD11	2.00	0.43
1:4:427:THR:O	1:4:467:ASN:HB2	2.17	0.43
1:4:807:VAL:HG13	1:4:808:GLU:N	2.33	0.43
1:2:422:PRO:HG3	1:3:285:TYR:CE1	2.52	0.43
1:2:800:ARG:HH11	1:2:800:ARG:HB3	1.83	0.43
1:4:651:LEU:HD22	1:4:667:GLU:OE1	2.18	0.43
1:4:750:GLU:HG3	1:4:755:ARG:HG2	1.99	0.43
1:1:610:ASP:O	1:1:611:ARG:HB2	2.18	0.43
1:3:301:TRP:CH2	1:3:452:SER:HA	2.53	0.43
1:3:663:LEU:HD21	1:3:686:PRO:HG2	2.01	0.43
1:3:963:SER:HB3	1:3:983:TRP:NE1	2.33	0.43
1:2:240:LEU:C	1:2:240:LEU:HD23	2.43	0.43
1:2:569:ASP:O	1:2:605:GLY:HA2	2.19	0.43
1:3:577:LYS:O	1:3:584:PRO:HA	2.19	0.43
1:4:102:ASN:ND2	1:4:103:VAL:HG23	2.34	0.43
1:4:937:LEU:HD12	1:4:957:PHE:C	2.43	0.43
1:1:59:ARG:HB2	1:1:124:SER:OG	2.18	0.43
1:1:352:ARG:HG2	1:1:553:TRP:CH2	2.54	0.43
1:1:708:TRP:CZ3	1:1:709:SER:HB3	2.54	0.43
1:3:807:VAL:HG13	1:3:808:GLU:N	2.32	0.43
1:3:843:GLN:HA	1:3:847:LYS:O	2.19	0.43
1:3:958:ASN:OD1	1:3:958:ASN:C	2.62	0.43
1:4:166:ARG:HG3	1:4:392:TYR:HB2	2.01	0.43
1:4:869:ASP:OD1	1:4:1015:HIS:ND1	2.42	0.43
1:1:606:LEU:O	1:1:614:HIS:HB2	2.19	0.43
1:1:984:LEU:HD21	1:1:986:ILE:HD11	2.00	0.43
1:2:126:THR:HA	1:2:182:ASN:O	2.19	0.43
1:2:128:ASN:HA	1:2:180:GLY:O	2.19	0.43
1:2:237:ARG:HG3	1:2:237:ARG:NH1	2.34	0.43
1:2:878:HIS:NE2	1:2:1010:SER:HB3	2.34	0.43
1:3:305:ILE:HD11	1:3:645:ARG:HB3	2.01	0.43
1:3:619:GLU:HA	1:3:912:ALA:HB2	2.01	0.43
1:3:663:LEU:HD23	1:3:663:LEU:C	2.44	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:4:464:HIS:HB2	1:4:489:GLY:HA3	2.01	0.43
1:4:542:MET:HE3	1:4:601:PHE:HA	2.01	0.43
1:4:737:ILE:HG23	1:4:737:ILE:O	2.18	0.43
1:1:246:MET:C	1:1:246:MET:SD	3.02	0.42
1:1:823:LEU:HD11	1:1:841:ALA:HB2	2.01	0.42
1:3:37:ARG:HG2	1:3:50:GLN:CD	2.44	0.42
1:3:930:VAL:O	1:3:932:PRO:HD3	2.19	0.42
1:1:427:THR:O	1:1:467:ASN:HB2	2.19	0.42
1:3:163:GLN:O	1:3:164:ASP:HB3	2.19	0.42
1:3:608:PHE:CE1	1:3:614:HIS:CD2	3.07	0.42
1:3:887:GLN:NE2	1:3:980:GLU:O	2.52	0.42
1:4:652:LEU:HD11	1:4:698:VAL:HB	2.00	0.42
1:1:433:LEU:N	1:1:434:PRO:CD	2.82	0.42
1:1:751:LEU:HD23	1:1:862:GLY:HA2	2.01	0.42
1:2:147:ASN:HA	1:2:148:SER:HA	1.67	0.42
1:3:111:PRO:HA	1:3:112:PRO:HA	1.85	0.42
1:3:127:PHE:CE2	1:3:184:LEU:HG	2.55	0.42
1:3:442:ARG:HD3	5:3:4049:HOH:O	2.19	0.42
1:4:853:ARG:NH1	1:4:871:GLU:OE2	2.42	0.42
1:2:54:LEU:O	2:2:8408:DMS:H12	2.20	0.42
1:3:754:LYS:HE3	1:3:1022:GLN:OE1	2.20	0.42
1:4:52:ARG:O	1:4:213:SER:HB2	2.20	0.42
1:1:233:ASP:CG	1:3:233:ASP:OD2	2.62	0.42
1:1:326:GLU:HA	1:1:326:GLU:OE1	2.20	0.42
1:1:463:GLY:HA2	5:1:4629:HOH:O	2.19	0.42
1:3:261:TRP:CZ3	1:3:266:GLN:HB2	2.55	0.42
1:1:899:GLY:HA2	1:1:915:PHE:CE1	2.55	0.42
1:3:751:LEU:HD23	1:3:862:GLY:HA2	2.01	0.42
1:4:377:LEU:O	1:4:381:GLN:HB2	2.20	0.42
1:2:1017:GLN:HB2	5:2:7130:HOH:O	2.19	0.42
1:3:375:ASP:O	1:3:379:MET:HG3	2.20	0.42
1:4:43:ARG:HD2	1:4:261:TRP:CD2	2.55	0.42
1:4:742:THR:CG2	1:4:743:SER:N	2.82	0.42
1:1:502:MET:CB	1:1:537:GLU:HB2	2.50	0.42
1:2:340:GLY:O	1:2:532:PRO:HB3	2.19	0.42
1:4:240:LEU:C	1:4:240:LEU:HD23	2.45	0.42
1:4:782:ASP:OD1	1:4:842:TRP:HH2	2.03	0.42
1:1:126:THR:H	2:1:8609:DMS:H22	1.83	0.42
1:1:225:PHE:HA	1:1:243:GLU:O	2.20	0.42
1:1:524:LEU:HD11	1:1:562:LEU:HG	2.02	0.42
1:1:561:ARG:HD3	1:2:525:SER:O	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:824:GLN:HG2	1:1:825:CYS:N	2.35	0.42
1:2:784:PHE:HA	1:2:881:ARG:O	2.19	0.42
1:3:355:ASN:OD1	1:3:388:ARG:HD3	2.20	0.42
1:3:361:PRO:HB3	1:3:609:ALA:HB1	2.01	0.42
1:3:730:LEU:HA	1:3:731:PRO:HD3	1.92	0.42
1:1:86:VAL:HG13	1:1:87:PRO:HA	2.01	0.42
1:3:133:TRP:CE3	1:3:216:HIS:HB2	2.54	0.42
1:4:35:SER:HB2	1:4:217:LYS:HD2	2.02	0.42
1:4:988:GLY:C	1:4:989:PHE:CD1	2.98	0.42
1:1:127:PHE:CD2	1:1:127:PHE:N	2.88	0.41
1:1:797:GLU:C	1:1:799:THR:N	2.78	0.41
1:1:991:MET:CE	1:1:1003:VAL:HG21	2.50	0.41
1:2:478:VAL:CG1	2:3:8420:DMS:H13	2.50	0.41
1:2:571:VAL:CG2	1:2:609:ALA:HA	2.49	0.41
1:2:583:ASN:HA	1:2:584:PRO:HD3	1.81	0.41
1:2:930:VAL:HA	1:2:973:ARG:HD3	2.01	0.41
1:4:844:HIS:ND1	1:4:845:GLN:HG2	2.34	0.41
1:1:16:TRP:CG	1:1:189:LEU:HD13	2.55	0.41
1:1:78:LEU:HA	1:1:79:PRO:HD3	1.92	0.41
1:1:416:GLU:HG3	1:1:460:ASN:O	2.19	0.41
1:2:842:TRP:HZ3	1:2:852:SER:HB3	1.85	0.41
1:3:907:PRO:HG2	1:3:990:HIS:O	2.21	0.41
1:4:178:ARG:HG2	1:4:182:ASN:OD1	2.20	0.41
1:1:753:ASN:HB2	1:1:771:GLY:HA2	2.00	0.41
1:1:937:LEU:HD12	1:1:957:PHE:C	2.45	0.41
1:2:372:MET:CE	1:2:395:HIS:HB3	2.48	0.41
1:2:995:GLY:O	1:2:996:ASP:C	2.64	0.41
1:3:127:PHE:N	1:3:127:PHE:CD2	2.88	0.41
1:3:254:LEU:C	1:3:255:ARG:HG2	2.45	0.41
1:3:482:ARG:HA	1:3:483:PRO:HD3	1.94	0.41
1:3:963:SER:HB3	1:3:983:TRP:CE2	2.56	0.41
1:4:663:LEU:O	1:4:663:LEU:HD23	2.21	0.41
1:4:726:LEU:HA	5:4:7519:HOH:O	2.20	0.41
1:2:246:MET:SD	1:2:246:MET:C	3.03	0.41
1:3:105:TYR:CE2	1:3:199:ASP:HB2	2.54	0.41
1:3:881:ARG:HE	1:3:987:ASP:CG	2.28	0.41
1:4:559:TYR:HB2	1:4:562:LEU:HD12	2.02	0.41
1:4:702:GLN:HE22	1:4:708:TRP:HH2	1.67	0.41
1:4:894:ARG:HH12	1:4:921:PRO:HG3	1.85	0.41
1:1:237:ARG:HH11	1:1:237:ARG:CB	2.34	0.41
1:2:578:TYR:CE2	1:2:584:PRO:HG3	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:2:646:HIS:CD2	1:2:673:ALA:HA	2.54	0.41
1:2:649:ASN:O	1:2:702:GLN:HA	2.21	0.41
1:4:78:LEU:HD23	5:4:4504:HOH:O	2.20	0.41
1:2:693:GLN:OE1	1:2:724:GLU:HB2	2.21	0.41
1:2:806:TRP:CE2	1:2:809:ARG:NH2	2.89	0.41
1:2:851:ILE:HB	1:2:871:GLU:HB2	2.01	0.41
1:4:105:TYR:CE2	1:4:199:ASP:HB2	2.55	0.41
1:4:663:LEU:HD23	1:4:663:LEU:C	2.46	0.41
1:1:755:ARG:HH21	1:1:772:ASP:HA	1.86	0.41
1:3:995:GLY:O	1:3:996:ASP:C	2.63	0.41
1:4:482:ARG:HA	1:4:483:PRO:HD3	1.96	0.41
1:4:610:ASP:O	1:4:611:ARG:HB2	2.20	0.41
1:2:361:PRO:HB3	1:2:609:ALA:HB1	2.03	0.41
1:3:542:MET:CE	1:3:600:GLN:HG2	2.51	0.41
1:1:513:PRO:O	1:1:514:ALA:HB3	2.21	0.41
1:1:595:THR:HA	1:1:596:PRO:C	2.45	0.41
1:1:601:PHE:CE2	1:1:796:SER:HA	2.56	0.41
1:1:801:ILE:N	1:1:801:ILE:CD1	2.84	0.41
1:1:833:ALA:HB2	1:1:859:ASP:HA	2.02	0.41
1:2:650:GLU:HB3	1:2:670:LEU:HD12	2.02	0.41
1:2:807:VAL:HG13	1:2:808:GLU:N	2.35	0.41
1:2:937:LEU:HA	1:2:957:PHE:O	2.21	0.41
1:3:487:GLU:C	1:3:487:GLU:CD	2.89	0.41
1:3:578:TYR:CE2	1:3:584:PRO:HB3	2.56	0.41
1:3:651:LEU:HD11	1:3:701:VAL:HB	2.03	0.41
1:3:767:GLN:HE21	1:3:774:LYS:HD3	1.86	0.41
1:3:883:GLY:HA3	1:3:987:ASP:HA	2.02	0.41
1:4:789:LEU:HD11	1:4:993:ILE:HG22	2.03	0.41
1:4:836:ILE:O	1:4:855:THR:HA	2.21	0.41
1:1:131:GLU:O	1:1:135:GLN:HG2	2.21	0.41
1:3:816:TYR:CD2	1:3:968:MET:HE1	2.56	0.41
1:1:630:ARG:HH11	1:1:630:ARG:HG2	1.86	0.40
1:1:820:ALA:HB2	1:1:842:TRP:CE2	2.56	0.40
1:2:658:LEU:O	1:2:659:ASP:C	2.64	0.40
1:3:522:LYS:HE3	1:4:560:PRO:HD3	2.03	0.40
1:3:873:ALA:O	1:3:876:THR:HG22	2.20	0.40
1:4:147:ASN:HA	1:4:148:SER:HA	1.64	0.40
1:1:148:SER:HA	1:1:165:SER:OG	2.20	0.40
1:1:577:LYS:O	1:1:584:PRO:HA	2.21	0.40
1:2:200:GLN:HG2	1:2:391:HIS:HB2	2.04	0.40
1:3:146:VAL:HG11	1:3:150:PHE:CD2	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:4:106:PRO:HD3	1:4:204:ARG:NH1	2.36	0.40
1:1:360:HIS:CE1	1:1:362:LEU:HB2	2.56	0.40
1:2:96:ASP:OD1	1:2:96:ASP:C	2.64	0.40
1:2:568:TRP:CD2	1:2:569:ASP:HB3	2.55	0.40
1:3:561:ARG:HD3	1:4:525:SER:O	2.22	0.40
1:3:945:ASN:CG	1:3:1023:LYS:HE2	2.46	0.40
1:4:577:LYS:O	1:4:584:PRO:HA	2.21	0.40
1:4:754:LYS:HE2	1:4:1022:GLN:OE1	2.21	0.40
1:1:478:VAL:HG12	2:4:8611:DMS:H23	2.02	0.40
1:2:359:HIS:CD2	1:2:573:GLN:HA	2.56	0.40
1:2:854:LYS:HA	1:2:867:THR:O	2.21	0.40
1:3:50:GLN:CD	1:3:50:GLN:N	2.79	0.40
1:3:568:TRP:CD1	1:3:568:TRP:C	2.99	0.40
1:3:725:ASN:HB2	5:3:6007:HOH:O	2.21	0.40
1:3:749:ILE:N	1:3:749:ILE:CD1	2.83	0.40
1:4:36:TRP:CH2	2:4:8404:DMS:H23	2.51	0.40
1:4:226:HIS:O	1:4:242:ALA:HA	2.21	0.40
1:1:670:LEU:HD23	1:1:670:LEU:HA	1.90	0.40
1:2:333:ARG:HA	1:2:345:ASN:OD1	2.21	0.40
1:3:630:ARG:HB2	1:3:637:GLU:HB3	2.04	0.40
1:3:778:THR:HG23	1:3:887:GLN:OE1	2.22	0.40
1:3:782:ASP:HB2	1:3:842:TRP:CH2	2.56	0.40
1:3:825:CYS:HA	1:3:837:THR:O	2.22	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%)	948 (94%)	56 (6%)	5 (0%)	24 43
1	2	1009/1023 (99%)	959 (95%)	49 (5%)	1 (0%)	48 68

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	3	1009/1023 (99%)	953 (94%)	55 (6%)	1 (0%)	48	68
1	4	1009/1023 (99%)	946 (94%)	60 (6%)	3 (0%)	36	55
All	All	4036/4092 (99%)	3806 (94%)	220 (6%)	10 (0%)	43	63

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	1	795	VAL
1	1	794	GLY
1	4	794	GLY
1	1	796	SER
1	1	798	ALA
1	3	164	ASP
1	4	164	ASP
1	2	164	ASP
1	1	511	PRO
1	4	511	PRO

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	1	863/874 (99%)	858 (99%)	5 (1%)	78	91
1	2	863/874 (99%)	853 (99%)	10 (1%)	63	83
1	3	863/874 (99%)	854 (99%)	9 (1%)	68	86
1	4	863/874 (99%)	857 (99%)	6 (1%)	76	89
All	All	3452/3496 (99%)	3422 (99%)	30 (1%)	70	87

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	1	519	SER
1	1	546	LEU

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Mol	Chain	Res	Type
1	1	672	VAL
1	1	795	VAL
1	1	885	ASN
1	2	80	GLU
1	2	519	SER
1	2	546	LEU
1	2	604	ASN
1	2	651	LEU
1	2	730	LEU
1	2	768	MET
1	2	800	ARG
1	2	850	PHE
1	2	885	ASN
1	3	80	GLU
1	3	178	ARG
1	3	262	GLN
1	3	519	SER
1	3	546	LEU
1	3	687	GLN
1	3	761	GLN
1	3	956	GLN
1	3	1017	GLN
1	4	112	PRO
1	4	519	SER
1	4	546	LEU
1	4	795	VAL
1	4	801	ILE
1	4	850	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (58) such sidechains are listed below:

Mol	Chain	Res	Type
1	1	266	GLN
1	1	294	ASN
1	1	297	ASN
1	1	363	HIS
1	1	460	ASN
1	1	485	GLN
1	1	510	GLN
1	1	558	GLN
1	1	623	GLN
1	1	624	GLN

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Mol	Chain	Res	Type
1	1	817	GLN
1	1	843	GLN
1	1	845	GLN
1	1	977	HIS
1	2	128	ASN
1	2	226	HIS
1	2	262	GLN
1	2	294	ASN
1	2	381	GLN
1	2	394	ASN
1	2	445	GLN
1	2	604	ASN
1	2	624	GLN
1	2	725	ASN
1	2	804	ASN
1	2	817	GLN
1	2	843	GLN
1	2	890	GLN
1	2	950	GLN
1	2	966	GLN
1	2	1022	GLN
1	3	50	GLN
1	3	245	GLN
1	3	294	ASN
1	3	297	ASN
1	3	394	ASN
1	3	485	GLN
1	3	510	GLN
1	3	623	GLN
1	3	634	GLN
1	3	702	GLN
1	3	739	HIS
1	3	817	GLN
1	3	824	GLN
1	3	885	ASN
1	3	977	HIS
1	4	266	GLN
1	4	297	ASN
1	4	445	GLN
1	4	675	GLN
1	4	702	GLN
1	4	757	GLN

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Mol	Chain	Res	Type
1	4	804	ASN
1	4	817	GLN
1	4	863	GLN
1	4	950	GLN
1	4	1017	GLN
1	4	1022	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 156 ligands modelled in this entry, 26 are monoatomic - leaving 130 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	$\# Z > 2$
2	DMS	2	8603	-	3,3,3	0.27	0	3,3,3	0.64	0
2	DMS	3	8410	-	3,3,3	0.25	0	3,3,3	0.63	0
2	DMS	4	8403	-	3,3,3	0.20	0	3,3,3	0.60	0
2	DMS	4	8405	-	3,3,3	0.29	0	3,3,3	0.68	0
2	DMS	3	8404	-	3,3,3	0.20	0	3,3,3	0.66	0
2	DMS	1	8408	-	3,3,3	0.25	0	3,3,3	0.67	0
2	DMS	4	1024	4	3,3,3	0.22	0	3,3,3	0.62	0
2	DMS	4	8404	-	3,3,3	0.27	0	3,3,3	0.67	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	DMS	3	8401	-	3,3,3	0.22	0	3,3,3	0.63	0
2	DMS	2	8423	-	3,3,3	0.25	0	3,3,3	0.62	0
2	DMS	1	8402	-	3,3,3	0.20	0	3,3,3	0.63	0
2	DMS	4	8417	-	3,3,3	0.24	0	3,3,3	0.63	0
2	DMS	2	8609	-	3,3,3	0.23	0	3,3,3	0.61	0
2	DMS	2	8403	-	3,3,3	0.32	0	3,3,3	0.63	0
2	DMS	2	8613	-	3,3,3	0.26	0	3,3,3	0.65	0
2	DMS	1	8602	-	3,3,3	0.25	0	3,3,3	0.63	0
2	DMS	1	8606	-	3,3,3	0.25	0	3,3,3	0.62	0
2	DMS	2	8404	-	3,3,3	0.26	0	3,3,3	0.63	0
2	DMS	3	8425	4	3,3,3	0.26	0	3,3,3	0.63	0
2	DMS	2	8402	-	3,3,3	0.30	0	3,3,3	0.60	0
2	DMS	3	8607	-	3,3,3	0.22	0	3,3,3	0.60	0
2	DMS	1	8610	-	3,3,3	0.32	0	3,3,3	0.68	0
2	DMS	1	8416	-	3,3,3	0.26	0	3,3,3	0.64	0
2	DMS	2	8604	-	3,3,3	0.25	0	3,3,3	0.63	0
2	DMS	3	8407	-	3,3,3	0.26	0	3,3,3	0.62	0
2	DMS	4	8608	-	3,3,3	0.26	0	3,3,3	0.64	0
2	DMS	4	8410	-	3,3,3	0.26	0	3,3,3	0.66	0
2	DMS	3	8501	-	3,3,3	0.24	0	3,3,3	0.66	0
2	DMS	3	8606	-	3,3,3	0.20	0	3,3,3	0.65	0
2	DMS	3	8409	-	3,3,3	0.28	0	3,3,3	0.62	0
2	DMS	3	8611	-	3,3,3	0.26	0	3,3,3	0.62	0
2	DMS	4	8406	-	3,3,3	0.32	0	3,3,3	0.59	0
2	DMS	2	8405	-	3,3,3	0.18	0	3,3,3	0.59	0
2	DMS	3	8408	-	3,3,3	0.23	0	3,3,3	0.66	0
2	DMS	2	8421	-	3,3,3	0.25	0	3,3,3	0.65	0
2	DMS	3	8416	-	3,3,3	0.26	0	3,3,3	0.64	0
2	DMS	4	8705	-	3,3,3	0.24	0	3,3,3	0.66	0
2	DMS	2	8420	-	3,3,3	0.26	0	3,3,3	0.62	0
2	DMS	1	8414	-	3,3,3	0.27	0	3,3,3	0.62	0
2	DMS	2	8409	-	3,3,3	0.27	0	3,3,3	0.68	0
2	DMS	2	8412	-	3,3,3	0.31	0	3,3,3	0.65	0
2	DMS	2	8611	-	3,3,3	0.27	0	3,3,3	0.66	0
2	DMS	1	8603	-	3,3,3	0.19	0	3,3,3	0.62	0
2	DMS	2	8508	-	3,3,3	0.25	0	3,3,3	0.63	0
2	DMS	3	8428	-	3,3,3	0.23	0	3,3,3	0.64	0
2	DMS	2	8417	-	3,3,3	0.26	0	3,3,3	0.64	0
2	DMS	2	8610	-	3,3,3	0.31	0	3,3,3	0.66	0
2	DMS	2	8408	-	3,3,3	0.25	0	3,3,3	0.65	0
2	DMS	1	8401	-	3,3,3	0.21	0	3,3,3	0.62	0
2	DMS	4	8411	-	3,3,3	0.31	0	3,3,3	0.63	0
2	DMS	4	8614	-	3,3,3	0.25	0	3,3,3	0.61	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	DMS	3	1024	-	3,3,3	0.24	0	3,3,3	0.63	0
2	DMS	4	8421	-	3,3,3	0.24	0	3,3,3	0.66	0
2	DMS	3	8610	-	3,3,3	0.30	0	3,3,3	0.66	0
2	DMS	1	8605	-	3,3,3	0.24	0	3,3,3	0.64	0
2	DMS	3	8603	-	3,3,3	0.23	0	3,3,3	0.61	0
2	DMS	1	8411	-	3,3,3	0.30	0	3,3,3	0.65	0
2	DMS	2	8601	-	3,3,3	0.27	0	3,3,3	0.63	0
2	DMS	4	8613	-	3,3,3	0.25	0	3,3,3	0.64	0
2	DMS	4	8610	-	3,3,3	0.25	0	3,3,3	0.61	0
2	DMS	3	8415	-	3,3,3	0.27	0	3,3,3	0.64	0
2	DMS	3	8403	-	3,3,3	0.31	0	3,3,3	0.67	0
2	DMS	2	8401	-	3,3,3	0.29	0	3,3,3	0.66	0
2	DMS	2	8612	-	3,3,3	0.22	0	3,3,3	0.64	0
2	DMS	4	8503	-	3,3,3	0.28	0	3,3,3	0.63	0
2	DMS	1	8425	4	3,3,3	0.21	0	3,3,3	0.61	0
2	DMS	2	8411	-	3,3,3	0.27	0	3,3,3	0.62	0
2	DMS	2	8414	-	3,3,3	0.25	0	3,3,3	0.63	0
2	DMS	1	8404	-	3,3,3	0.30	0	3,3,3	0.65	0
2	DMS	3	8405	-	3,3,3	0.22	0	3,3,3	0.63	0
2	DMS	4	8701	-	3,3,3	0.30	0	3,3,3	0.61	0
2	DMS	3	8421	-	3,3,3	0.26	0	3,3,3	0.62	0
2	DMS	4	8412	-	3,3,3	0.24	0	3,3,3	0.62	0
2	DMS	1	8604	-	3,3,3	0.26	0	3,3,3	0.64	0
2	DMS	3	8613	-	3,3,3	0.19	0	3,3,3	0.61	0
2	DMS	1	8412	-	3,3,3	0.26	0	3,3,3	0.59	0
2	DMS	2	8602	-	3,3,3	0.27	0	3,3,3	0.61	0
2	DMS	1	8405	-	3,3,3	0.28	0	3,3,3	0.65	0
2	DMS	2	8425	4	3,3,3	0.25	0	3,3,3	0.62	0
2	DMS	3	8604	-	3,3,3	0.21	0	3,3,3	0.60	0
2	DMS	4	8703	-	3,3,3	0.20	0	3,3,3	0.59	0
2	DMS	2	8605	-	3,3,3	0.23	0	3,3,3	0.64	0
2	DMS	1	8503	-	3,3,3	0.25	0	3,3,3	0.62	0
2	DMS	3	8417	-	3,3,3	0.25	0	3,3,3	0.65	0
2	DMS	4	8618	-	3,3,3	0.20	0	3,3,3	0.61	0
2	DMS	3	8609	-	3,3,3	0.21	0	3,3,3	0.64	0
2	DMS	4	8508	-	3,3,3	0.21	0	3,3,3	0.62	0
2	DMS	3	8608	-	3,3,3	0.22	0	3,3,3	0.64	0
2	DMS	1	8609	-	3,3,3	0.24	0	3,3,3	0.66	0
2	DMS	4	8402	-	3,3,3	0.18	0	3,3,3	0.63	0
2	DMS	1	8607	-	3,3,3	0.23	0	3,3,3	0.63	0
2	DMS	4	8615	-	3,3,3	0.25	0	3,3,3	0.65	0
2	DMS	3	8423	-	3,3,3	0.26	0	3,3,3	0.63	0
2	DMS	3	8612	-	3,3,3	0.28	0	3,3,3	0.67	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	DMS	1	8501	-	3,3,3	0.23	0	3,3,3	0.66	0
2	DMS	2	8410	-	3,3,3	0.25	0	3,3,3	0.65	0
2	DMS	1	8419	-	3,3,3	0.24	0	3,3,3	0.64	0
2	DMS	3	8411	-	3,3,3	0.27	0	3,3,3	0.65	0
2	DMS	2	8504	-	3,3,3	0.24	0	3,3,3	0.63	0
2	DMS	1	8403	-	3,3,3	0.24	0	3,3,3	0.63	0
2	DMS	1	8406	-	3,3,3	0.26	0	3,3,3	0.64	0
2	DMS	3	8402	-	3,3,3	0.33	0	3,3,3	0.69	0
2	DMS	4	8414	-	3,3,3	0.27	0	3,3,3	0.63	0
2	DMS	4	8611	-	3,3,3	0.20	0	3,3,3	0.61	0
2	DMS	4	8419	-	3,3,3	0.23	0	3,3,3	0.64	0
2	DMS	2	8608	-	3,3,3	0.25	0	3,3,3	0.63	0
2	DMS	3	8602	-	3,3,3	0.21	0	3,3,3	0.61	0
2	DMS	3	8420	-	3,3,3	0.26	0	3,3,3	0.62	0
2	DMS	4	8501	-	3,3,3	0.25	0	3,3,3	0.63	0
2	DMS	4	8607	-	3,3,3	0.24	0	3,3,3	0.63	0
2	DMS	4	8617	-	3,3,3	0.25	0	3,3,3	0.62	0
2	DMS	4	8409	-	3,3,3	0.26	0	3,3,3	0.66	0
2	DMS	1	8421	-	3,3,3	0.26	0	3,3,3	0.63	0
2	DMS	3	8601	-	3,3,3	0.26	0	3,3,3	0.61	0
2	DMS	2	8606	-	3,3,3	0.23	0	3,3,3	0.62	0
2	DMS	2	8406	-	3,3,3	0.20	0	3,3,3	0.61	0
2	DMS	4	8401	-	3,3,3	0.23	0	3,3,3	0.65	0
2	DMS	1	8504	-	3,3,3	0.29	0	3,3,3	0.63	0
2	DMS	4	8423	-	3,3,3	0.26	0	3,3,3	0.62	0
2	DMS	3	8414	-	3,3,3	0.27	0	3,3,3	0.64	0
2	DMS	3	8605	-	3,3,3	0.23	0	3,3,3	0.64	0
2	DMS	4	8408	-	3,3,3	0.27	0	3,3,3	0.64	0
2	DMS	4	8416	-	3,3,3	0.23	0	3,3,3	0.63	0
2	DMS	2	8607	-	3,3,3	0.21	0	3,3,3	0.62	0
2	DMS	4	8612	-	3,3,3	0.22	0	3,3,3	0.64	0
2	DMS	2	8416	-	3,3,3	0.22	0	3,3,3	0.63	0
2	DMS	2	8502	-	3,3,3	0.25	0	3,3,3	0.65	0
2	DMS	3	8412	-	3,3,3	0.26	0	3,3,3	0.62	0
2	DMS	1	8409	-	3,3,3	0.27	0	3,3,3	0.62	0
2	DMS	1	8407	-	3,3,3	0.23	0	3,3,3	0.64	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

32 monomers are involved in 37 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	4	8403	DMS	1	0
2	4	1024	DMS	1	0
2	4	8404	DMS	2	0
2	4	8417	DMS	1	0
2	2	8403	DMS	1	0
2	3	8425	DMS	1	0
2	3	8606	DMS	1	0
2	3	8409	DMS	1	0
2	4	8406	DMS	1	0
2	2	8405	DMS	1	0
2	3	8408	DMS	1	0
2	3	8416	DMS	1	0
2	4	8705	DMS	1	0
2	2	8420	DMS	1	0
2	2	8409	DMS	2	0
2	2	8611	DMS	1	0
2	2	8610	DMS	1	0
2	2	8408	DMS	1	0
2	3	8603	DMS	1	0
2	2	8612	DMS	1	0
2	2	8414	DMS	1	0
2	3	8405	DMS	1	0
2	1	8609	DMS	1	0
2	1	8607	DMS	1	0
2	4	8615	DMS	1	0
2	3	8612	DMS	1	0
2	4	8611	DMS	3	0
2	3	8602	DMS	1	0
2	3	8420	DMS	2	0
2	2	8606	DMS	1	0
2	1	8409	DMS	1	0
2	1	8407	DMS	1	0

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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%)	-0.52	17 (1%) 69 65	5, 17, 37, 78	0
1	2	1011/1023 (98%)	-0.52	6 (0%) 85 83	4, 17, 37, 68	0
1	3	1011/1023 (98%)	-0.57	6 (0%) 85 83	4, 17, 39, 70	0
1	4	1011/1023 (98%)	-0.55	14 (1%) 73 70	4, 17, 38, 81	0
All	All	4044/4092 (98%)	-0.54	43 (1%) 78 75	4, 17, 38, 81	0

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	1	795	VAL	8.6
1	1	796	SER	8.3
1	4	795	VAL	7.4
1	4	796	SER	6.4
1	1	801	ILE	6.4
1	1	798	ALA	5.9
1	4	798	ALA	4.6
1	4	799	THR	4.4
1	1	797	GLU	4.4
1	4	801	ILE	4.0
1	1	799	THR	4.0
1	4	800	ARG	3.5
1	4	797	GLU	3.3
1	1	686	PRO	3.2
1	1	800	ARG	3.1
1	1	735	HIS	3.1
1	2	733	ALA	2.8
1	1	688	PRO	2.8
1	2	688	PRO	2.6
1	4	686	PRO	2.6
1	3	731	PRO	2.6

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Mol	Chain	Res	Type	RSRZ
1	1	689	GLU	2.6
1	2	745	MET	2.5
1	1	771	GLY	2.5
1	3	730	LEU	2.5
1	3	734	SER	2.5
1	1	634	GLN	2.4
1	4	634	GLN	2.4
1	2	734	SER	2.3
1	3	735	HIS	2.3
1	4	685	LEU	2.3
1	1	730	LEU	2.2
1	3	732	ALA	2.2
1	4	733	ALA	2.2
1	4	732	ALA	2.2
1	2	686	PRO	2.2
1	1	736	ALA	2.2
1	3	745	MET	2.2
1	4	688	PRO	2.2
1	4	735	HIS	2.1
1	2	798	ALA	2.1
1	1	687	GLN	2.1
1	1	772	ASP	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 [i](#)

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
2	DMS	4	8417	4/4	0.70	0.30	91,91,91,92	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	DMS	1	8610	4/4	0.74	0.25	66,66,67,68	0
2	DMS	4	8503	4/4	0.76	0.29	83,83,84,84	0
2	DMS	4	8614	4/4	0.77	0.24	61,61,62,62	0
2	DMS	2	8417	4/4	0.80	0.23	63,65,65,65	0
2	DMS	1	8605	4/4	0.80	0.21	71,71,71,72	0
2	DMS	1	8602	4/4	0.81	0.22	80,80,80,80	0
2	DMS	3	8410	4/4	0.82	0.26	77,78,78,78	0
2	DMS	4	8613	4/4	0.82	0.20	56,57,58,58	0
2	DMS	2	8603	4/4	0.82	0.27	71,72,72,73	0
2	DMS	4	1024	4/4	0.83	0.20	70,70,70,70	0
2	DMS	2	8410	4/4	0.84	0.23	87,87,88,88	0
2	DMS	4	8615	4/4	0.84	0.21	49,50,51,52	0
2	DMS	2	8601	4/4	0.85	0.23	69,69,70,70	0
2	DMS	1	8604	4/4	0.85	0.20	66,67,67,68	0
2	DMS	2	8609	4/4	0.85	0.21	67,68,68,68	0
2	DMS	2	8508	4/4	0.85	0.20	72,72,73,73	0
2	DMS	3	8420	4/4	0.85	0.24	73,74,74,75	0
2	DMS	3	8609	4/4	0.85	0.19	52,52,54,54	0
2	DMS	1	8407	4/4	0.86	0.22	72,72,73,73	0
2	DMS	3	1024	4/4	0.86	0.19	77,77,77,77	0
2	DMS	1	8416	4/4	0.86	0.20	57,57,57,58	0
2	DMS	1	8419	4/4	0.86	0.18	68,68,68,68	0
2	DMS	2	8610	4/4	0.87	0.16	51,51,52,52	0
2	DMS	4	8410	4/4	0.87	0.20	65,65,65,65	0
2	DMS	3	8407	4/4	0.87	0.23	69,70,70,71	0
2	DMS	2	8423	4/4	0.87	0.23	76,76,76,76	0
2	DMS	3	8415	4/4	0.87	0.17	60,61,61,62	0
2	DMS	3	8417	4/4	0.87	0.20	66,66,67,67	0
2	DMS	2	8608	4/4	0.87	0.17	53,54,55,55	0
2	DMS	2	8602	4/4	0.87	0.19	63,64,64,65	0
2	DMS	4	8618	4/4	0.87	0.20	59,59,60,61	0
4	NA	3	3105	1/1	0.87	0.13	54,54,54,54	0
2	DMS	2	8420	4/4	0.88	0.17	57,57,58,58	0
2	DMS	3	8610	4/4	0.88	0.16	49,49,49,50	0
2	DMS	1	8609	4/4	0.88	0.17	43,43,45,45	0
2	DMS	2	8613	4/4	0.89	0.16	63,64,65,65	0
2	DMS	2	8604	4/4	0.89	0.17	62,63,63,63	0
2	DMS	3	8428	4/4	0.89	0.22	79,79,79,79	0
2	DMS	3	8607	4/4	0.89	0.17	65,65,65,66	0
2	DMS	4	8611	4/4	0.89	0.19	67,67,67,68	0
2	DMS	1	8503	4/4	0.89	0.18	63,63,64,64	0
2	DMS	2	8611	4/4	0.89	0.17	60,60,61,62	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	DMS	3	8611	4/4	0.89	0.20	60,61,62,62	0
2	DMS	3	8416	4/4	0.89	0.14	70,70,70,71	0
2	DMS	4	8608	4/4	0.89	0.14	52,52,52,53	0
2	DMS	4	8617	4/4	0.90	0.17	70,71,71,71	0
2	DMS	3	8608	4/4	0.90	0.16	58,59,59,60	0
4	NA	3	3103	1/1	0.90	0.09	34,34,34,34	0
2	DMS	3	8423	4/4	0.90	0.17	63,63,64,64	0
2	DMS	3	8601	4/4	0.91	0.18	59,59,59,60	0
2	DMS	4	8612	4/4	0.91	0.14	60,60,61,62	0
2	DMS	3	8602	4/4	0.91	0.15	55,55,56,56	0
2	DMS	3	8606	4/4	0.91	0.15	54,54,54,55	0
2	DMS	3	8421	4/4	0.91	0.16	58,58,58,58	0
2	DMS	1	8414	4/4	0.91	0.16	50,50,50,52	0
2	DMS	4	8421	4/4	0.91	0.15	60,61,61,62	0
4	NA	1	3103	1/1	0.91	0.06	30,30,30,30	0
4	NA	2	3104	1/1	0.91	0.09	27,27,27,27	0
2	DMS	3	8425	4/4	0.91	0.16	58,58,59,59	0
2	DMS	2	8409	4/4	0.91	0.14	42,43,44,45	0
2	DMS	2	8612	4/4	0.92	0.15	48,49,50,52	0
2	DMS	2	8406	4/4	0.92	0.16	58,58,59,59	0
2	DMS	3	8605	4/4	0.92	0.15	63,64,64,65	0
2	DMS	1	8408	4/4	0.92	0.15	45,46,47,47	0
2	DMS	4	8423	4/4	0.92	0.14	63,63,63,64	0
2	DMS	3	8612	4/4	0.92	0.14	47,47,48,49	0
2	DMS	4	8703	4/4	0.92	0.15	59,59,60,60	0
2	DMS	3	8613	4/4	0.92	0.13	61,61,61,63	0
2	DMS	4	8610	4/4	0.92	0.15	55,55,55,56	0
2	DMS	1	8421	4/4	0.92	0.16	69,69,69,69	0
4	NA	4	3104	1/1	0.92	0.09	54,54,54,54	0
2	DMS	4	8508	4/4	0.93	0.13	58,58,58,58	0
2	DMS	2	8606	4/4	0.93	0.12	46,47,47,48	0
2	DMS	4	8414	4/4	0.93	0.14	52,52,53,54	0
2	DMS	1	8607	4/4	0.93	0.15	53,53,53,54	0
2	DMS	4	8419	4/4	0.93	0.12	52,52,52,53	0
2	DMS	1	8409	4/4	0.93	0.14	51,52,52,52	0
4	NA	3	3104	1/1	0.93	0.06	39,39,39,39	0
2	DMS	3	8603	4/4	0.93	0.15	64,64,64,65	0
4	NA	4	3101	1/1	0.93	0.05	18,18,18,18	0
2	DMS	2	8605	4/4	0.93	0.14	62,62,63,63	0
2	DMS	4	8705	4/4	0.94	0.13	57,57,58,58	0
2	DMS	4	8607	4/4	0.94	0.12	53,54,54,55	0
2	DMS	1	8412	4/4	0.94	0.14	44,45,46,47	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	DMS	1	8425	4/4	0.94	0.13	49,49,49,49	0
2	DMS	1	8603	4/4	0.94	0.12	57,57,58,58	0
2	DMS	4	8409	4/4	0.94	0.12	45,45,45,46	0
2	DMS	1	8501	4/4	0.94	0.14	47,47,48,48	0
2	DMS	3	8408	4/4	0.94	0.12	38,39,40,41	0
2	DMS	4	8416	4/4	0.94	0.12	51,51,52,53	0
2	DMS	4	8406	4/4	0.95	0.12	34,34,34,35	0
2	DMS	4	8408	4/4	0.95	0.10	34,36,37,37	0
2	DMS	2	8412	4/4	0.95	0.11	37,37,38,38	0
2	DMS	3	8412	4/4	0.95	0.11	43,44,44,45	0
2	DMS	4	8411	4/4	0.95	0.12	38,40,41,42	0
2	DMS	1	8406	4/4	0.95	0.15	58,59,60,60	0
2	DMS	1	8402	4/4	0.95	0.12	28,29,31,31	0
2	DMS	2	8421	4/4	0.95	0.12	52,53,53,53	0
2	DMS	1	8504	4/4	0.95	0.12	43,43,43,44	0
3	MG	1	3003	1/1	0.95	0.15	41,41,41,41	0
2	DMS	2	8425	4/4	0.95	0.11	52,53,53,53	0
2	DMS	2	8502	4/4	0.95	0.12	48,48,48,49	0
2	DMS	2	8408	4/4	0.95	0.14	55,55,55,56	0
2	DMS	1	8606	4/4	0.95	0.13	39,40,40,41	0
2	DMS	1	8411	4/4	0.95	0.10	40,41,41,42	0
2	DMS	2	8411	4/4	0.95	0.12	38,40,40,40	0
4	NA	4	3103	1/1	0.95	0.08	41,41,41,41	0
2	DMS	4	8405	4/4	0.95	0.12	42,43,43,43	0
2	DMS	3	8404	4/4	0.96	0.09	30,31,32,33	0
2	DMS	2	8404	4/4	0.96	0.10	38,39,40,41	0
2	DMS	2	8416	4/4	0.96	0.10	52,52,53,54	0
2	DMS	3	8409	4/4	0.96	0.11	39,39,40,40	0
2	DMS	2	8504	4/4	0.96	0.14	63,63,63,63	0
2	DMS	2	8402	4/4	0.96	0.10	28,28,29,32	0
4	NA	4	3102	1/1	0.96	0.04	18,18,18,18	0
2	DMS	3	8414	4/4	0.96	0.11	41,41,43,43	0
2	DMS	3	8402	4/4	0.96	0.10	27,27,27,29	0
2	DMS	3	8501	4/4	0.97	0.09	34,36,36,36	0
2	DMS	2	8607	4/4	0.97	0.09	44,45,46,46	0
3	MG	4	3003	1/1	0.97	0.11	29,29,29,29	0
4	NA	1	3101	1/1	0.97	0.05	17,17,17,17	0
4	NA	1	3102	1/1	0.97	0.07	16,16,16,16	0
2	DMS	4	8402	4/4	0.97	0.08	25,28,28,29	0
4	NA	2	3102	1/1	0.97	0.04	29,29,29,29	0
2	DMS	4	8404	4/4	0.97	0.09	45,46,47,48	0
2	DMS	1	8404	4/4	0.97	0.09	38,39,39,39	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	DMS	2	8405	4/4	0.97	0.10	40,40,40,40	0
2	DMS	3	8604	4/4	0.97	0.10	44,44,44,44	0
2	DMS	4	8501	4/4	0.97	0.10	43,43,44,45	0
2	DMS	2	8414	4/4	0.97	0.10	50,51,52,52	0
2	DMS	3	8403	4/4	0.97	0.11	34,34,35,35	0
2	DMS	4	8701	4/4	0.97	0.08	21,22,23,25	0
4	NA	2	3103	1/1	0.98	0.04	28,28,28,28	0
2	DMS	2	8403	4/4	0.98	0.07	29,29,29,30	0
4	NA	3	3102	1/1	0.98	0.04	17,17,17,17	0
3	MG	4	3001	1/1	0.98	0.03	7,7,7,7	0
3	MG	4	3002	1/1	0.98	0.06	11,11,11,11	0
2	DMS	4	8412	4/4	0.98	0.07	33,33,34,35	0
2	DMS	3	8405	4/4	0.98	0.06	32,32,33,34	0
2	DMS	1	8405	4/4	0.98	0.08	32,32,33,33	0
2	DMS	4	8403	4/4	0.98	0.08	26,26,27,29	0
2	DMS	3	8411	4/4	0.98	0.08	38,38,38,38	0
3	MG	2	3002	1/1	0.99	0.07	12,12,12,12	0
3	MG	3	3001	1/1	0.99	0.02	15,15,15,15	0
4	NA	3	3101	1/1	0.99	0.01	10,10,10,10	0
3	MG	3	3002	1/1	0.99	0.07	17,17,17,17	0
2	DMS	2	8401	4/4	0.99	0.04	14,17,17,19	0
2	DMS	4	8401	4/4	0.99	0.05	20,20,20,22	0
2	DMS	3	8401	4/4	0.99	0.07	16,17,17,18	0
2	DMS	1	8401	4/4	0.99	0.04	12,12,13,15	0
3	MG	1	3002	1/1	0.99	0.05	23,23,23,23	0
2	DMS	1	8403	4/4	0.99	0.05	29,29,30,31	0
3	MG	2	3001	1/1	0.99	0.02	11,11,11,11	0
3	MG	1	3001	1/1	1.00	0.02	14,14,14,14	0
4	NA	2	3101	1/1	1.00	0.01	12,12,12,12	0

6.5 Other polymers ⓘ

There are no such residues in this entry.