



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 09:56 PM GMT

PDB ID : 1R8Y
Title : Crystal Structure of Mouse Glycine N-Methyltransferase (Monoclinic Form)
Authors : Pakhomova, S.; Luka, Z.; Wagner, C.; Newcomer, M.E.
Deposited on : 2003-10-28
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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

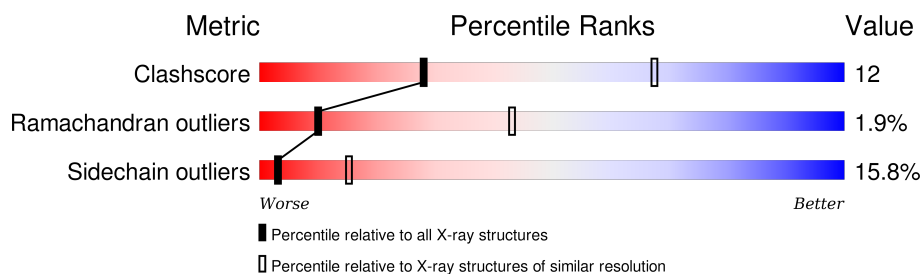
1 Overall quality at a glance ⓘ

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))
Clashscore	102246	1912 (3.00-3.00)
Ramachandran outliers	100387	1853 (3.00-3.00)
Sidechain outliers	100360	1856 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	292	
1	B	292	
1	C	292	
1	D	292	
1	E	292	
1	F	292	
1	G	292	

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Mol	Chain	Length	Quality of chain
1	H	292	 <div>66% 25% 6%</div>

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
2	BME	B	2057	-	-	X	-

2 Entry composition

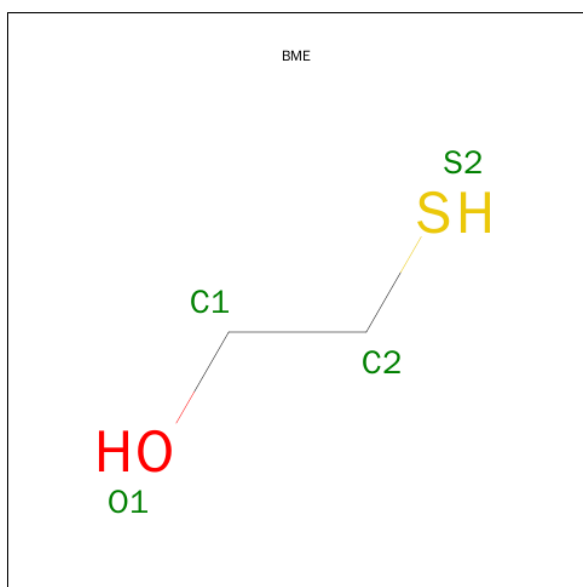
There are 3 unique types of molecules in this entry. The entry contains 17606 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 glycine N-methyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	281	Total	C	N	O	S	0	0	0
			2192	1391	380	408	13			
1	B	282	Total	C	N	O	S	0	0	0
			2201	1399	382	407	13			
1	C	283	Total	C	N	O	S	0	0	0
			2175	1382	378	402	13			
1	D	286	Total	C	N	O	S	0	0	0
			2226	1417	382	414	13			
1	E	286	Total	C	N	O	S	0	0	0
			2216	1408	385	410	13			
1	F	282	Total	C	N	O	S	0	0	0
			2137	1357	370	397	13			
1	G	283	Total	C	N	O	S	0	0	0
			2217	1409	385	410	13			
1	H	282	Total	C	N	O	S	0	0	0
			2188	1391	379	405	13			

- Molecule 2 is BETA-MERCAPTOETHANOL (three-letter code: BME) (formula: C₂H₆OS).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	O	S	0	0
			4	2	1	1		
2	A	1	Total	C	O	S	0	0
			4	2	1	1		
2	B	1	Total	C	O	S	0	0
			4	2	1	1		
2	D	1	Total	C	O	S	0	0
			4	2	1	1		
2	D	1	Total	C	O	S	0	0
			4	2	1	1		
2	D	1	Total	C	O	S	0	0
			4	2	1	1		
2	E	1	Total	C	O	S	0	0
			4	2	1	1		
2	F	1	Total	C	O	S	0	0
			4	2	1	1		
2	G	1	Total	C	O	S	0	0
			4	2	1	1		
2	G	1	Total	C	O	S	0	0
			4	2	1	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	C	4	Total	O	0	0
			4	4		
3	D	3	Total	O	0	0
			3	3		

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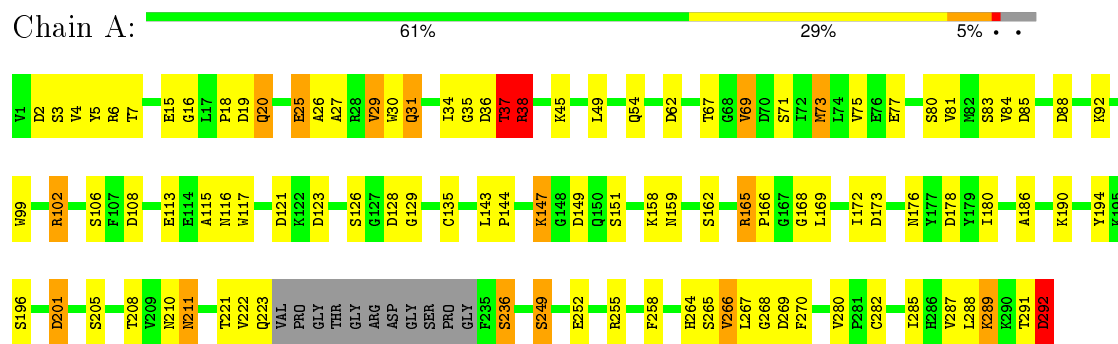
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	E	2	Total	O	0	0
			2	2		
3	F	3	Total	O	0	0
			3	3		
3	G	2	Total	O	0	0
			2	2		

3 Residue-property plots

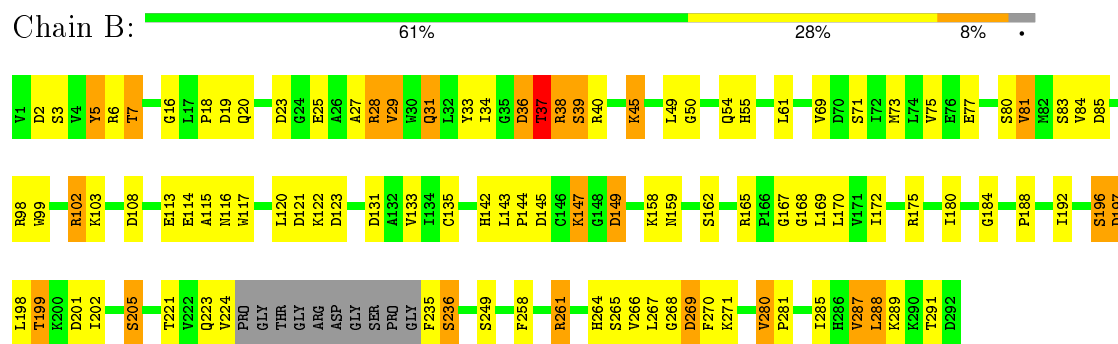
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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.

Note EDS was not executed.

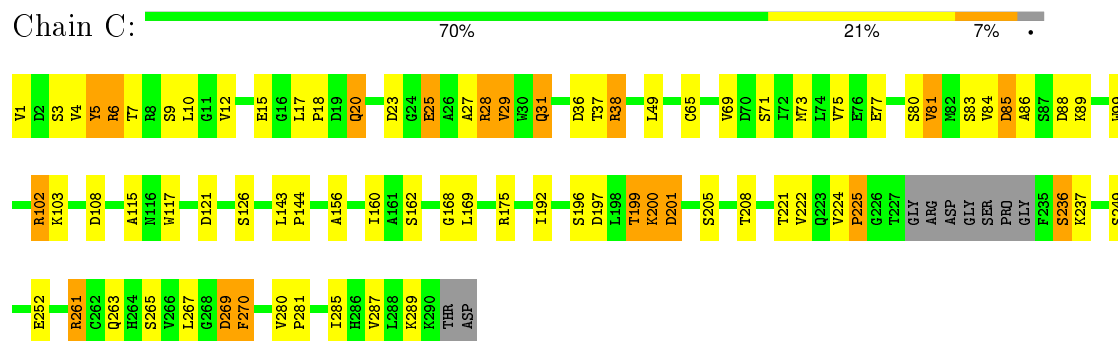
- Molecule 1: glycine N-methyltransferase



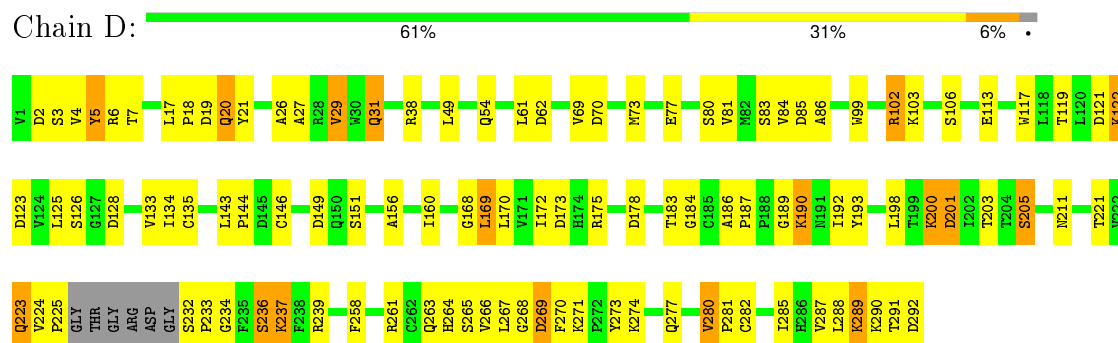
- Molecule 1: glycine N-methyltransferase



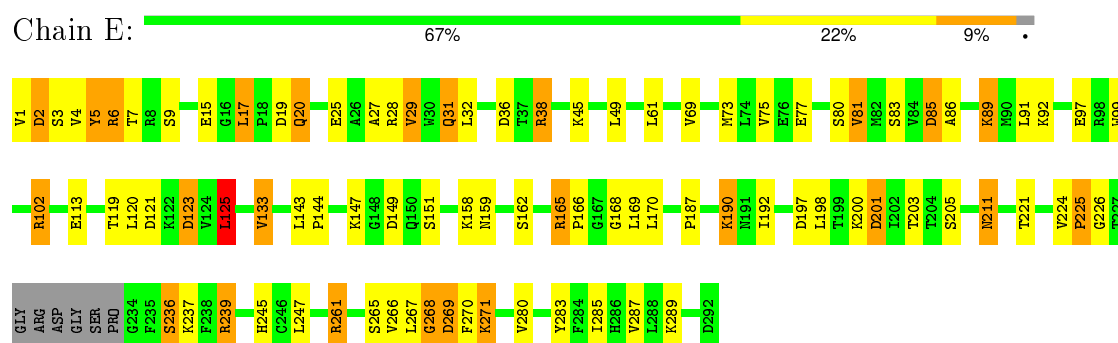
- Molecule 1: glycine N-methyltransferase



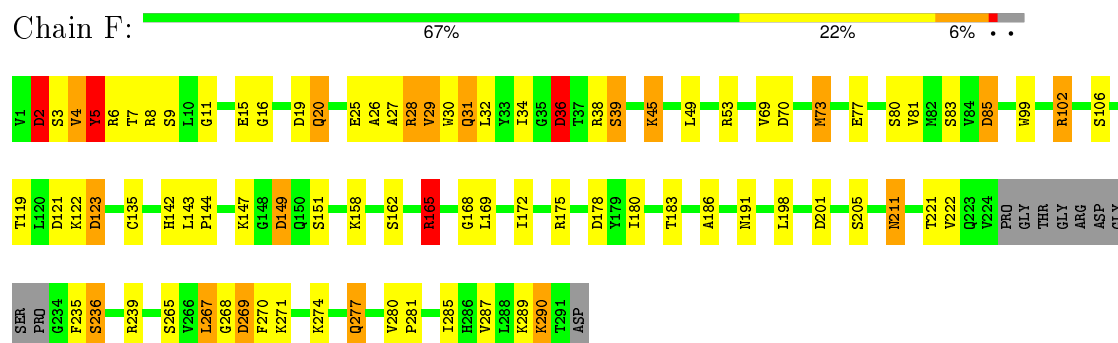
- Molecule 1: glycine N-methyltransferase



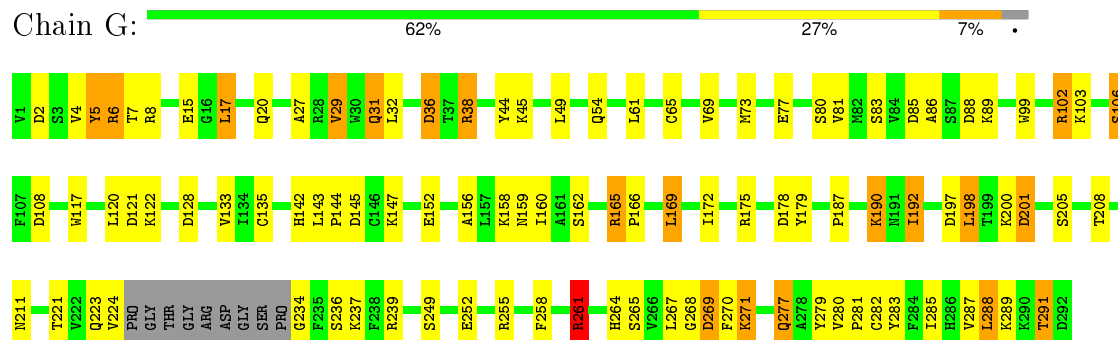
- Molecule 1: glycine N-methyltransferase



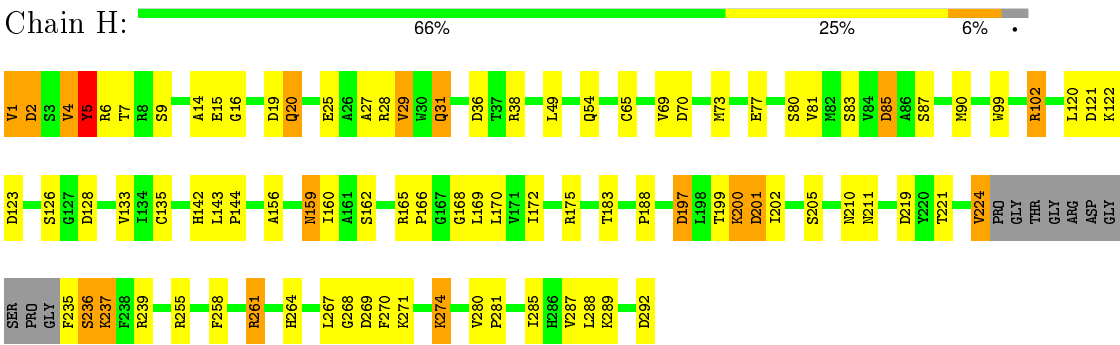
- Molecule 1: glycine N-methyltransferase



- Molecule 1: glycine N-methyltransferase



● Molecule 1: glycine N-methyltransferase



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	107.98Å 108.30Å 119.01Å 90.00° 93.71° 90.00°	Depositor
Resolution (Å)	13.00 – 3.00	Depositor
% Data completeness (in resolution range)	74.2 (13.00-3.00)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.222 , 0.285	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	17606	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

5 Model quality ⓘ

5.1 Standard geometry ⓘ

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

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	A	0.83	1/2244 (0.0%)	0.98	12/3047 (0.4%)
1	B	0.83	3/2253 (0.1%)	0.93	8/3058 (0.3%)
1	C	0.74	0/2228	0.90	7/3029 (0.2%)
1	D	0.83	0/2280	0.96	9/3096 (0.3%)
1	E	0.79	0/2269	0.94	9/3081 (0.3%)
1	F	0.77	0/2189	0.92	9/2978 (0.3%)
1	G	0.85	1/2269 (0.0%)	0.99	13/3076 (0.4%)
1	H	0.79	1/2240 (0.0%)	0.95	6/3043 (0.2%)
All	All	0.81	6/17972 (0.0%)	0.95	73/24408 (0.3%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	C	0	2
1	E	0	1
1	G	0	1
1	H	0	1
All	All	0	6

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	196	SER	CA-CB	5.91	1.61	1.52
1	H	1	VAL	CB-CG1	5.46	1.64	1.52
1	G	152	GLU	CD-OE1	5.41	1.31	1.25
1	B	37	THR	CA-CB	5.29	1.67	1.53
1	A	37	THR	CA-CB	5.18	1.66	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	33	TYR	CE1-CZ	5.08	1.45	1.38

All (73) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	123	ASP	CB-CG-OD2	8.89	126.30	118.30
1	G	198	LEU	CA-CB-CG	7.98	133.65	115.30
1	D	128	ASP	CB-CG-OD2	7.56	125.11	118.30
1	C	269	ASP	CB-CG-OD2	7.45	125.00	118.30
1	H	128	ASP	CB-CG-OD2	7.43	124.99	118.30
1	E	2	ASP	CB-CG-OD2	7.25	124.83	118.30
1	G	36	ASP	CB-CG-OD2	7.01	124.61	118.30
1	C	85	ASP	CB-CG-OD2	6.94	124.54	118.30
1	D	178	ASP	CB-CG-OD2	6.76	124.38	118.30
1	D	269	ASP	CB-CG-OD2	6.68	124.32	118.30
1	F	269	ASP	CB-CG-OD2	6.64	124.28	118.30
1	A	269	ASP	CB-CG-OD2	6.53	124.18	118.30
1	H	123	ASP	CB-CG-OD2	6.53	124.18	118.30
1	F	123	ASP	CB-CG-OD2	6.49	124.14	118.30
1	F	121	ASP	CB-CG-OD2	6.43	124.09	118.30
1	A	128	ASP	CB-CG-OD2	6.42	124.08	118.30
1	B	198	LEU	N-CA-C	-6.40	93.71	111.00
1	D	292	ASP	CB-CG-OD2	6.38	124.04	118.30
1	F	36	ASP	CB-CG-OD2	6.38	124.04	118.30
1	C	121	ASP	CB-CG-OD2	6.25	123.93	118.30
1	B	269	ASP	CB-CG-OD2	6.23	123.91	118.30
1	H	121	ASP	CB-CG-OD2	6.21	123.89	118.30
1	H	85	ASP	CB-CG-OD2	6.21	123.89	118.30
1	E	269	ASP	CB-CG-OD2	6.17	123.85	118.30
1	H	201	ASP	CB-CG-OD2	6.16	123.85	118.30
1	A	178	ASP	CB-CG-OD2	6.12	123.80	118.30
1	F	2	ASP	CB-CG-OD2	6.07	123.77	118.30
1	G	128	ASP	CB-CG-OD2	6.03	123.73	118.30
1	A	123	ASP	CB-CG-OD2	5.99	123.69	118.30
1	F	178	ASP	CB-CG-OD2	5.94	123.64	118.30
1	A	292	ASP	CB-CG-OD2	5.91	123.62	118.30
1	E	197	ASP	CB-CG-OD2	5.90	123.61	118.30
1	A	201	ASP	CB-CG-OD2	5.89	123.60	118.30
1	C	108	ASP	CB-CG-OD2	5.88	123.59	118.30
1	C	197	ASP	CB-CG-OD2	5.86	123.57	118.30
1	F	149	ASP	CB-CG-OD2	5.85	123.57	118.30
1	A	149	ASP	CB-CG-OD2	5.85	123.56	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	85	ASP	CB-CG-OD2	5.81	123.53	118.30
1	A	173	ASP	CB-CG-OD2	5.81	123.53	118.30
1	C	201	ASP	CB-CG-OD2	5.79	123.51	118.30
1	G	178	ASP	CB-CG-OD2	5.79	123.51	118.30
1	E	201	ASP	CB-CG-OD2	5.71	123.44	118.30
1	G	88	ASP	CB-CG-OD1	5.68	123.41	118.30
1	G	108	ASP	CB-CG-OD2	5.63	123.37	118.30
1	B	85	ASP	CB-CG-OD2	5.62	123.36	118.30
1	G	208	THR	OG1-CB-CG2	-5.60	97.11	110.00
1	B	197	ASP	CB-CG-OD2	5.60	123.34	118.30
1	C	88	ASP	CB-CG-OD1	5.60	123.34	118.30
1	E	85	ASP	CB-CG-OD2	5.58	123.32	118.30
1	F	85	ASP	CB-CG-OD2	5.58	123.32	118.30
1	B	149	ASP	CB-CG-OD2	5.56	123.31	118.30
1	H	70	ASP	CB-CG-OD2	5.53	123.28	118.30
1	G	269	ASP	CB-CG-OD2	5.53	123.28	118.30
1	D	123	ASP	CB-CG-OD2	5.51	123.26	118.30
1	E	121	ASP	CB-CG-OD2	5.50	123.25	118.30
1	B	19	ASP	CB-CG-OD2	5.43	123.19	118.30
1	F	70	ASP	CB-CG-OD2	5.42	123.18	118.30
1	G	197	ASP	CB-CG-OD2	5.41	123.16	118.30
1	B	123	ASP	CB-CG-OD2	5.40	123.16	118.30
1	E	239	ARG	NE-CZ-NH2	-5.35	117.62	120.30
1	B	121	ASP	CB-CG-OD2	5.35	123.12	118.30
1	D	62	ASP	CB-CG-OD2	5.33	123.10	118.30
1	A	62	ASP	CB-CG-OD2	5.31	123.08	118.30
1	E	125	LEU	CA-CB-CG	5.30	127.49	115.30
1	D	121	ASP	CB-CG-OD2	5.20	122.98	118.30
1	G	201	ASP	CB-CG-OD2	5.19	122.97	118.30
1	G	261	ARG	CB-CA-C	5.18	120.77	110.40
1	G	145	ASP	CB-CG-OD2	5.11	122.90	118.30
1	D	201	ASP	CB-CG-OD2	5.10	122.89	118.30
1	A	121	ASP	CB-CG-OD2	5.05	122.85	118.30
1	A	88	ASP	CB-CG-OD1	5.02	122.82	118.30
1	D	173	ASP	CB-CG-OD2	5.02	122.82	118.30
1	G	121	ASP	CB-CG-OD2	5.00	122.80	118.30

There are no chirality outliers.

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	38	ARG	Peptide

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Mol	Chain	Res	Type	Group
1	C	36	ASP	Peptide
1	C	38	ARG	Peptide
1	E	1	VAL	Peptide
1	G	38	ARG	Peptide
1	H	2	ASP	Peptide

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2192	0	2106	72	0
1	B	2201	0	2121	78	0
1	C	2175	0	2063	41	0
1	D	2226	0	2156	64	0
1	E	2216	0	2134	58	0
1	F	2137	0	1985	46	0
1	G	2217	0	2158	47	0
1	H	2188	0	2103	53	0
2	A	8	0	10	3	0
2	B	4	0	5	5	0
2	D	12	0	15	1	0
2	E	4	0	5	0	0
2	F	4	0	5	0	0
2	G	8	0	10	2	0
3	C	4	0	0	1	0
3	D	3	0	0	1	0
3	E	2	0	0	3	0
3	F	3	0	0	1	0
3	G	2	0	0	0	0
All	All	17606	0	16876	402	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 12.

All (402) 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:H:4:VAL:HG12	1:H:5:TYR:H	1.15	1.09
1:A:264:HIS:ND1	1:A:288:LEU:HD21	1.71	1.05
1:A:37:THR:O	1:A:38:ARG:HG3	1.62	0.99
1:C:20:GLN:HG2	3:C:293:HOH:O	1.66	0.96
1:E:20:GLN:HG2	3:E:5247:HOH:O	1.67	0.94
1:E:49:LEU:HD11	1:E:77:GLU:HG3	1.48	0.93
1:D:223:GLN:H	1:D:223:GLN:HE21	1.16	0.92
1:F:183:THR:O	1:G:211:ASN:ND2	2.05	0.90
1:D:224:VAL:O	1:D:233:PRO:HB3	1.72	0.89
1:D:190:LYS:HG2	1:D:282:CYS:SG	2.13	0.88
1:E:36:ASP:HA	1:E:38:ARG:HH22	1.36	0.88
1:A:29:VAL:HG21	1:A:236:SER:HB3	1.56	0.86
1:A:102:ARG:HD2	1:B:108:ASP:OD1	1.74	0.86
1:D:119:THR:HB	1:D:122:LYS:HZ3	1.39	0.86
1:E:211:ASN:OD1	1:H:183:THR:O	1.94	0.86
1:E:133:VAL:HG12	1:E:170:LEU:CD1	2.06	0.84
1:H:4:VAL:HG12	1:H:5:TYR:N	1.93	0.83
1:G:32:LEU:O	1:G:36:ASP:HB2	1.79	0.83
1:B:223:GLN:NE2	1:B:235:PHE:HE2	1.77	0.81
1:C:200:LYS:HB2	1:C:222:VAL:HG22	1.62	0.81
1:D:232:SER:N	1:D:233:PRO:HD3	1.96	0.81
1:A:49:LEU:HD11	1:A:77:GLU:HG3	1.62	0.81
1:C:168:GLY:O	1:C:289:LYS:HA	1.82	0.79
1:F:29:VAL:HG21	1:F:236:SER:HB3	1.65	0.79
1:C:75:VAL:HG22	1:C:81:VAL:HG13	1.65	0.79
1:D:119:THR:HB	1:D:122:LYS:NZ	1.97	0.79
1:F:45:LYS:HG2	1:F:73:MET:SD	2.23	0.79
1:A:108:ASP:OD1	1:B:102:ARG:HD2	1.82	0.78
1:B:29:VAL:HG21	1:B:236:SER:HB3	1.66	0.77
1:E:36:ASP:HA	1:E:38:ARG:NH2	1.98	0.77
1:H:49:LEU:HD11	1:H:77:GLU:HG3	1.67	0.77
1:F:25:GLU:O	1:F:28:ARG:HB3	1.86	0.76
1:F:49:LEU:HD11	1:F:77:GLU:HG3	1.68	0.76
1:B:5:TYR:CD1	1:B:5:TYR:O	2.38	0.76
1:D:198:LEU:HB3	1:D:223:GLN:HE22	1.49	0.75
1:B:49:LEU:HD11	1:B:77:GLU:HG3	1.66	0.75
1:C:224:VAL:CG2	1:C:225:PRO:HD2	2.16	0.74
1:A:37:THR:HG23	1:A:194:TYR:HB2	1.67	0.74
1:B:264:HIS:ND1	1:B:288:LEU:HD21	2.02	0.74
1:B:40:ARG:HD2	1:B:45:LYS:CG	2.18	0.74
1:A:38:ARG:HG2	1:A:69:VAL:HG11	1.70	0.74
1:B:38:ARG:HG2	1:B:39:SER:H	1.53	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:5:TYR:O	1:B:5:TYR:HD1	1.71	0.73
1:C:29:VAL:HG21	1:C:236:SER:HB3	1.71	0.73
1:D:29:VAL:HG21	1:D:236:SER:HB3	1.69	0.73
1:H:258:PHE:O	1:H:261:ARG:HG2	1.88	0.73
1:G:29:VAL:HG21	1:G:236:SER:HB3	1.70	0.73
1:B:40:ARG:HG2	1:B:45:LYS:HD2	1.71	0.73
1:G:143:LEU:HD12	1:G:144:PRO:HD2	1.70	0.73
1:E:133:VAL:HG12	1:E:170:LEU:HD12	1.70	0.72
1:D:189:GLY:H	1:D:200:LYS:NZ	1.87	0.72
1:E:61:LEU:HB3	1:E:133:VAL:HG22	1.71	0.71
1:C:49:LEU:HD11	1:C:77:GLU:HG3	1.73	0.70
1:E:125:LEU:HD13	1:E:125:LEU:O	1.90	0.70
1:G:135:CYS:HB3	1:G:172:ILE:HG13	1.71	0.70
1:F:53:ARG:HD2	3:F:6283:HOH:O	1.90	0.70
1:E:85:ASP:OD1	1:F:15:GLU:O	2.10	0.70
1:F:28:ARG:HD3	1:F:29:VAL:HG22	1.74	0.69
1:D:189:GLY:H	1:D:200:LYS:HZ3	1.39	0.69
1:H:29:VAL:HG21	1:H:236:SER:HB3	1.74	0.69
1:G:49:LEU:HD11	1:G:77:GLU:HG3	1.74	0.69
1:E:29:VAL:HG21	1:E:236:SER:HB3	1.74	0.69
1:C:249:SER:O	1:C:252:GLU:HB2	1.93	0.68
1:D:49:LEU:HD11	1:D:77:GLU:HG3	1.75	0.68
1:E:120:LEU:HB3	1:E:159:ASN:HB3	1.76	0.68
1:E:49:LEU:CD1	1:E:77:GLU:HG3	2.24	0.67
1:B:40:ARG:CD	1:B:45:LYS:HG3	2.24	0.67
1:A:190:LYS:HD2	1:A:282:CYS:SG	2.34	0.67
1:E:224:VAL:HG22	1:E:225:PRO:HD2	1.77	0.66
1:H:1:VAL:O	1:H:1:VAL:HG12	1.93	0.66
1:D:223:GLN:HE21	1:D:223:GLN:N	1.91	0.66
1:B:40:ARG:CG	1:B:45:LYS:HD2	2.25	0.66
1:H:4:VAL:CG1	1:H:5:TYR:H	1.93	0.66
1:E:89:LYS:NZ	1:F:11:GLY:O	2.25	0.66
1:E:6:ARG:O	1:F:239:ARG:HD3	1.96	0.66
1:G:179:TYR:HB3	2:G:7282:BME:H12	1.78	0.66
1:H:224:VAL:HB	1:H:235:PHE:N	2.10	0.66
1:A:15:GLU:H	1:B:142:HIS:CE1	2.15	0.65
1:A:264:HIS:ND1	1:A:288:LEU:CD2	2.56	0.65
1:A:210:ASN:OD1	2:D:4185:BME:S2	2.56	0.64
1:A:29:VAL:CG2	1:A:236:SER:HB3	2.26	0.64
1:B:55:HIS:O	2:B:2057:BME:H22	1.97	0.64
1:A:71:SER:O	1:A:75:VAL:HG23	1.98	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:23:ASP:HB2	1:C:28:ARG:NH2	2.13	0.63
1:A:258:PHE:CE2	1:A:288:LEU:HD12	2.34	0.63
1:B:23:ASP:HB2	1:B:28:ARG:CZ	2.29	0.63
1:A:222:VAL:HG12	1:A:223:GLN:H	1.62	0.63
1:C:224:VAL:HG22	1:C:225:PRO:HD2	1.81	0.62
1:A:166:PRO:HB3	1:A:292:ASP:HA	1.81	0.62
1:D:258:PHE:CE2	1:D:288:LEU:HD13	2.34	0.62
1:G:6:ARG:O	1:H:239:ARG:HD3	2.00	0.62
1:A:255:ARG:HD3	2:A:1262:BME:H22	1.81	0.62
1:A:143:LEU:HD12	1:A:144:PRO:HD2	1.81	0.61
1:D:223:GLN:H	1:D:223:GLN:NE2	1.93	0.61
1:C:143:LEU:HD12	1:C:144:PRO:HD2	1.82	0.61
1:A:102:ARG:HH11	1:B:108:ASP:CG	2.03	0.61
1:A:129:GLY:HA3	1:A:165:ARG:HB3	1.82	0.61
1:B:50:GLY:O	1:B:54:GLN:HB2	2.01	0.61
1:B:287:VAL:C	1:B:288:LEU:HD23	2.21	0.61
1:F:168:GLY:O	1:F:289:LYS:HA	2.01	0.61
1:F:45:LYS:CG	1:F:73:MET:SD	2.88	0.60
1:A:210:ASN:ND2	1:D:203:THR:HG23	2.17	0.60
1:E:133:VAL:HG12	1:E:170:LEU:HD13	1.83	0.59
1:B:5:TYR:C	1:B:5:TYR:HD1	2.06	0.59
1:D:168:GLY:O	1:D:289:LYS:HA	2.03	0.59
1:D:221:THR:HG23	1:D:237:LYS:HE2	1.84	0.59
1:B:38:ARG:CG	1:B:39:SER:H	2.14	0.59
1:D:29:VAL:CG2	1:D:236:SER:HB3	2.32	0.59
1:B:40:ARG:HD2	1:B:45:LYS:HG3	1.83	0.59
1:A:117:TRP:CZ3	1:B:16:GLY:HA3	2.37	0.59
1:D:224:VAL:HB	1:D:234:GLY:H	1.68	0.59
1:E:133:VAL:O	1:E:170:LEU:HD12	2.03	0.59
1:E:32:LEU:CD1	1:E:224:VAL:HG23	2.33	0.59
1:E:245:HIS:HE1	3:E:5248:HOH:O	1.86	0.58
1:A:99:TRP:CE2	1:B:98:ARG:CZ	2.85	0.58
1:E:75:VAL:HG22	1:E:81:VAL:HG13	1.85	0.58
1:E:192:ILE:HG12	1:E:283:TYR:CD1	2.37	0.58
1:D:169:LEU:HA	1:D:288:LEU:O	2.04	0.58
1:F:29:VAL:CG2	1:F:236:SER:HB3	2.34	0.57
1:E:239:ARG:N	1:F:9:SER:OG	2.37	0.57
1:H:143:LEU:HD12	1:H:144:PRO:HD2	1.85	0.57
1:F:5:TYR:C	1:F:5:TYR:CD1	2.77	0.57
1:H:27:ALA:O	1:H:31:GLN:HG2	2.04	0.57
1:E:36:ASP:OD1	1:E:38:ARG:NH2	2.37	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:29:VAL:CG2	1:H:236:SER:HB3	2.35	0.57
1:G:29:VAL:CG2	1:G:236:SER:HB3	2.35	0.57
1:B:268:GLY:N	1:B:271:LYS:O	2.33	0.56
1:B:261:ARG:NH2	1:B:291:THR:OG1	2.30	0.56
1:G:17:LEU:HD21	1:H:142:HIS:C	2.26	0.56
1:B:7:THR:HG21	1:C:1:VAL:O	2.05	0.56
1:A:35:GLY:O	1:A:38:ARG:HD3	2.05	0.56
1:G:6:ARG:O	1:H:239:ARG:NH1	2.25	0.56
1:H:219:ASP:HB3	1:H:237:LYS:HE3	1.87	0.56
1:H:1:VAL:O	1:H:2:ASP:C	2.42	0.56
1:D:156:ALA:O	1:D:160:ILE:HD12	2.06	0.56
1:C:27:ALA:O	1:C:31:GLN:HG2	2.06	0.56
1:B:131:ASP:HA	1:B:165:ARG:HE	1.70	0.56
1:B:199:THR:CB	1:B:223:GLN:HG3	2.37	0.55
1:B:29:VAL:CG2	1:B:236:SER:HB3	2.34	0.55
1:B:34:ILE:O	1:B:37:THR:OG1	2.18	0.55
1:E:6:ARG:O	1:F:239:ARG:NH1	2.24	0.55
1:E:29:VAL:CG2	1:E:236:SER:HB3	2.36	0.55
1:C:4:VAL:O	1:C:4:VAL:HG12	2.07	0.55
1:C:224:VAL:HG23	1:C:225:PRO:HD2	1.89	0.55
1:F:5:TYR:HD1	1:F:5:TYR:C	2.10	0.55
1:E:245:HIS:CE1	3:E:5248:HOH:O	2.60	0.55
1:B:120:LEU:HB3	1:B:159:ASN:HB3	1.87	0.54
1:F:4:VAL:HG12	1:F:4:VAL:O	2.07	0.54
1:A:210:ASN:HD21	1:D:203:THR:HA	1.73	0.54
1:G:192:ILE:HG12	1:G:283:TYR:CD1	2.42	0.54
1:D:232:SER:N	1:D:233:PRO:CD	2.70	0.54
1:H:1:VAL:O	1:H:1:VAL:CG1	2.55	0.54
1:H:36:ASP:OD2	1:H:197:ASP:HB2	2.07	0.54
1:B:5:TYR:CD1	1:B:5:TYR:C	2.81	0.54
1:E:203:THR:HA	1:H:210:ASN:HD21	1.73	0.53
1:A:190:LYS:CD	1:A:282:CYS:SG	2.95	0.53
1:G:117:TRP:CZ3	1:H:16:GLY:HA3	2.44	0.53
1:A:27:ALA:O	1:A:31:GLN:HG2	2.07	0.53
1:G:17:LEU:HD21	1:H:142:HIS:O	2.09	0.53
1:B:223:GLN:NE2	1:B:235:PHE:CE2	2.68	0.53
1:A:108:ASP:CG	1:B:102:ARG:HH11	2.11	0.53
1:E:267:LEU:HB2	1:E:285:ILE:HB	1.90	0.53
1:F:8:ARG:NE	1:G:2:ASP:OD2	2.40	0.53
1:C:29:VAL:CG2	1:C:236:SER:HB3	2.39	0.53
1:A:267:LEU:N	1:A:285:ILE:O	2.41	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:267:LEU:HB2	1:C:285:ILE:HB	1.91	0.53
1:B:201:ASP:HB3	1:B:221:THR:HB	1.91	0.53
1:B:267:LEU:HB2	1:B:285:ILE:HB	1.90	0.53
1:D:263:GLN:HB2	1:D:289:LYS:HB2	1.91	0.52
1:F:2:ASP:OD2	1:G:8:ARG:NH2	2.39	0.52
1:H:168:GLY:O	1:H:289:LYS:HA	2.09	0.52
1:A:37:THR:O	1:A:38:ARG:CG	2.47	0.52
1:D:27:ALA:O	1:D:31:GLN:HG2	2.09	0.52
1:H:49:LEU:CD1	1:H:77:GLU:HG3	2.39	0.52
1:C:75:VAL:HG22	1:C:81:VAL:CG1	2.36	0.52
1:E:224:VAL:CG2	1:E:225:PRO:HD2	2.39	0.52
1:B:165:ARG:NH2	2:B:2057:BME:H11	2.25	0.52
1:F:30:TRP:CH2	1:F:34:ILE:HD13	2.44	0.52
1:B:40:ARG:HD2	1:B:45:LYS:CD	2.40	0.51
1:H:268:GLY:N	1:H:271:LYS:O	2.38	0.51
1:G:224:VAL:O	1:G:234:GLY:HA3	2.10	0.51
1:F:268:GLY:N	1:F:271:LYS:O	2.39	0.51
1:B:205:SER:HA	1:C:208:THR:O	2.10	0.51
1:F:30:TRP:CZ2	1:F:34:ILE:HG21	2.46	0.51
1:G:85:ASP:OD1	1:G:86:ALA:N	2.44	0.51
1:B:38:ARG:HG2	1:B:39:SER:N	2.24	0.51
1:B:133:VAL:O	1:B:170:LEU:HD12	2.09	0.51
1:C:224:VAL:HG22	1:C:225:PRO:CD	2.41	0.51
1:F:165:ARG:O	1:F:290:LYS:HD2	2.11	0.51
1:E:36:ASP:CA	1:E:38:ARG:NH2	2.72	0.51
1:D:274:LYS:N	1:D:277:GLN:OE1	2.39	0.51
1:A:116:ASN:ND2	1:B:18:PRO:HG3	2.26	0.51
1:G:187:PRO:HB2	1:G:190:LYS:HD3	1.92	0.51
1:B:40:ARG:HD3	1:B:45:LYS:HG3	1.93	0.51
1:A:201:ASP:HB3	1:A:221:THR:HB	1.91	0.50
1:A:19:ASP:O	1:A:20:GLN:C	2.49	0.50
1:F:119:THR:OG1	1:F:123:ASP:OD2	2.28	0.50
1:D:268:GLY:N	1:D:271:LYS:O	2.40	0.50
1:B:167:GLY:O	2:B:2057:BME:H12	2.10	0.50
1:E:203:THR:HG23	1:H:210:ASN:ND2	2.27	0.50
1:E:165:ARG:HG2	1:E:166:PRO:N	2.26	0.50
1:D:267:LEU:HB2	1:D:285:ILE:HB	1.94	0.50
1:G:268:GLY:N	1:G:271:LYS:O	2.34	0.50
1:B:196:SER:O	1:B:197:ASP:C	2.49	0.50
1:D:273:TYR:HD1	1:D:277:GLN:NE2	2.09	0.50
1:A:18:PRO:HG3	1:B:116:ASN:HD21	1.76	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:61:LEU:HB3	1:D:133:VAL:HG22	1.94	0.50
1:B:143:LEU:HD12	1:B:144:PRO:HD2	1.94	0.50
1:E:99:TRP:O	1:E:102:ARG:HB3	2.11	0.49
1:A:211:ASN:ND2	1:D:183:THR:O	2.45	0.49
1:D:4:VAL:O	1:D:4:VAL:HG23	2.11	0.49
1:A:116:ASN:HD21	1:B:18:PRO:HG3	1.77	0.49
1:F:143:LEU:HD12	1:F:144:PRO:HD2	1.93	0.49
1:F:38:ARG:O	1:F:39:SER:C	2.51	0.49
1:H:267:LEU:HB2	1:H:285:ILE:HB	1.95	0.49
1:G:190:LYS:HG2	1:G:282:CYS:SG	2.52	0.49
1:F:32:LEU:O	1:F:36:ASP:HB2	2.13	0.48
1:A:267:LEU:HB2	1:A:285:ILE:HB	1.95	0.48
1:E:143:LEU:HD12	1:E:144:PRO:HD2	1.95	0.48
1:G:120:LEU:HB3	1:G:159:ASN:HB3	1.95	0.48
1:A:49:LEU:CD1	1:A:77:GLU:HG3	2.38	0.48
1:C:201:ASP:HB3	1:C:221:THR:HB	1.96	0.48
1:A:2:ASP:O	1:A:3:SER:HB3	2.14	0.48
1:H:133:VAL:O	1:H:170:LEU:HD12	2.13	0.48
1:G:15:GLU:H	1:H:142:HIS:CE1	2.31	0.48
1:B:168:GLY:O	1:B:289:LYS:HA	2.13	0.48
1:H:258:PHE:O	1:H:261:ARG:CG	2.59	0.48
1:E:86:ALA:HB3	1:F:16:GLY:O	2.14	0.48
1:H:19:ASP:O	1:H:20:GLN:C	2.51	0.48
1:E:168:GLY:O	1:E:289:LYS:HA	2.13	0.48
1:C:6:ARG:O	1:D:239:ARG:HD3	2.13	0.48
1:F:149:ASP:C	1:F:151:SER:H	2.17	0.48
1:F:201:ASP:HB3	1:F:221:THR:HB	1.95	0.48
1:A:15:GLU:OE2	1:B:142:HIS:HE1	1.97	0.47
1:G:258:PHE:O	1:G:261:ARG:NH2	2.43	0.47
1:B:131:ASP:O	2:B:2057:BME:S2	2.72	0.47
1:C:99:TRP:O	1:C:102:ARG:HB3	2.14	0.47
1:G:267:LEU:HB2	1:G:285:ILE:HB	1.96	0.47
1:H:135:CYS:HB3	1:H:172:ILE:HG13	1.96	0.47
1:G:223:GLN:HA	1:G:234:GLY:O	2.14	0.47
1:H:172:ILE:HG23	1:H:172:ILE:O	2.13	0.47
1:A:92:LYS:HZ1	1:B:114:GLU:CD	2.17	0.47
1:G:249:SER:O	1:G:252:GLU:HB2	2.14	0.47
1:E:32:LEU:HD13	1:E:224:VAL:HG23	1.95	0.47
1:C:156:ALA:O	1:C:160:ILE:HD12	2.15	0.47
1:A:135:CYS:HB3	1:A:172:ILE:HG13	1.96	0.47
1:D:273:TYR:CD1	1:D:277:GLN:CD	2.89	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:176:ASN:HB3	2:A:1282:BME:H12	1.97	0.46
1:A:45:LYS:HG3	1:A:73:MET:SD	2.55	0.46
1:B:27:ALA:O	1:B:31:GLN:HG2	2.14	0.46
1:G:277:GLN:HE21	1:G:279:TYR:HB3	1.81	0.46
1:E:271:LYS:HG2	1:E:271:LYS:H	1.56	0.46
1:H:261:ARG:HG2	1:H:261:ARG:H	1.63	0.46
1:E:187:PRO:HB2	1:E:190:LYS:HG2	1.97	0.46
1:C:84:VAL:HB	1:C:115:ALA:HB3	1.98	0.46
1:H:156:ALA:O	1:H:160:ILE:HD12	2.16	0.46
1:D:223:GLN:HG2	1:D:225:PRO:HD3	1.98	0.46
1:D:187:PRO:HB2	1:D:190:LYS:HD2	1.97	0.46
1:F:28:ARG:HG2	1:F:29:VAL:N	2.31	0.46
1:D:117:TRP:CD1	1:D:160:ILE:HD11	2.51	0.46
1:D:143:LEU:HD12	1:D:144:PRO:HD2	1.98	0.46
1:G:4:VAL:HG23	1:G:4:VAL:O	2.16	0.46
1:G:142:HIS:CG	1:H:14:ALA:HB1	2.51	0.46
1:A:147:LYS:HD3	1:A:151:SER:OG	2.16	0.46
1:C:117:TRP:CD1	1:C:160:ILE:HD11	2.50	0.45
1:D:189:GLY:HA2	1:D:200:LYS:HZ2	1.81	0.45
1:C:23:ASP:HB2	1:C:28:ARG:HH21	1.79	0.45
1:F:267:LEU:HB2	1:F:285:ILE:HB	1.97	0.45
1:E:201:ASP:HB3	1:E:221:THR:HB	1.98	0.45
1:H:201:ASP:HB3	1:H:221:THR:HB	1.99	0.45
1:H:274:LYS:H	1:H:274:LYS:HD2	1.81	0.45
1:E:133:VAL:CG1	1:E:170:LEU:HD13	2.46	0.45
1:G:27:ALA:O	1:G:31:GLN:HG2	2.17	0.45
1:E:17:LEU:HD21	1:F:142:HIS:O	2.17	0.45
1:H:4:VAL:CG1	1:H:5:TYR:N	2.64	0.45
1:A:108:ASP:OD2	1:B:103:LYS:NZ	2.34	0.45
1:F:222:VAL:O	1:F:235:PHE:HA	2.16	0.45
1:A:99:TRP:CD1	1:B:98:ARG:NH2	2.85	0.45
1:C:15:GLU:O	1:D:85:ASP:OD1	2.35	0.45
1:F:274:LYS:O	1:F:277:GLN:HG2	2.16	0.45
1:A:190:LYS:HG3	1:A:282:CYS:SG	2.57	0.45
1:G:8:ARG:HA	1:H:239:ARG:HG2	1.99	0.45
1:G:169:LEU:HD12	1:G:288:LEU:O	2.17	0.45
1:H:65:CYS:HB3	1:H:85:ASP:HB2	1.98	0.45
1:H:99:TRP:O	1:H:102:ARG:HB3	2.16	0.45
1:A:180:ILE:HG12	1:A:186:ALA:HA	1.98	0.45
1:A:255:ARG:CD	2:A:1262:BME:H22	2.45	0.45
1:B:36:ASP:HA	1:B:38:ARG:NH1	2.31	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:38:ARG:O	1:B:39:SER:C	2.55	0.45
1:C:4:VAL:O	1:C:6:ARG:N	2.50	0.45
1:H:255:ARG:HH21	1:H:264:HIS:CD2	2.35	0.45
1:G:201:ASP:HB3	1:G:221:THR:HB	1.99	0.45
1:F:19:ASP:O	1:F:20:GLN:C	2.55	0.45
1:B:75:VAL:HG23	1:B:81:VAL:HG13	1.98	0.45
1:D:280:VAL:HA	1:D:281:PRO:HD3	1.87	0.44
1:B:165:ARG:HH21	2:B:2057:BME:H21	1.81	0.44
1:F:27:ALA:O	1:F:31:GLN:HG2	2.16	0.44
1:D:192:ILE:HG22	1:D:193:TYR:CE2	2.52	0.44
1:E:224:VAL:O	1:E:226:GLY:N	2.50	0.44
1:A:25:GLU:HG3	1:A:26:ALA:N	2.31	0.44
1:C:9:SER:HB2	1:C:12:VAL:HG13	1.98	0.44
1:B:49:LEU:CD1	1:B:77:GLU:HG3	2.42	0.44
1:A:15:GLU:N	1:B:142:HIS:ND1	2.65	0.44
1:B:180:ILE:O	1:B:184:GLY:N	2.38	0.44
1:D:273:TYR:HA	1:D:277:GLN:OE1	2.18	0.44
1:A:211:ASN:HD22	1:D:184:GLY:HA3	1.82	0.44
1:E:5:TYR:C	1:E:5:TYR:CD1	2.91	0.44
1:D:135:CYS:HB3	1:D:172:ILE:HG13	2.00	0.44
1:H:120:LEU:HB3	1:H:159:ASN:HB3	2.00	0.44
1:A:249:SER:O	1:A:252:GLU:HB2	2.18	0.44
1:G:165:ARG:HD2	1:G:166:PRO:O	2.18	0.44
1:E:85:ASP:OD1	1:E:86:ALA:N	2.51	0.44
1:B:135:CYS:HB3	1:B:172:ILE:HG13	2.00	0.44
1:E:27:ALA:O	1:E:31:GLN:HG2	2.18	0.44
1:H:188:PRO:HB3	1:H:202:ILE:HD12	2.00	0.43
1:C:199:THR:O	1:C:199:THR:HG22	2.18	0.43
1:H:166:PRO:HB3	1:H:292:ASP:HA	2.01	0.43
1:G:65:CYS:HB3	1:G:85:ASP:HB2	2.00	0.43
1:A:208:THR:O	1:D:205:SER:HA	2.17	0.43
1:A:38:ARG:HH21	1:A:67:THR:CB	2.30	0.43
1:F:5:TYR:HB3	1:G:5:TYR:HD2	1.84	0.43
1:D:273:TYR:CD1	1:D:277:GLN:NE2	2.86	0.43
1:A:16:GLY:HA3	1:B:117:TRP:CZ3	2.53	0.43
1:D:201:ASP:HB3	1:D:221:THR:HB	2.00	0.43
1:G:85:ASP:OD1	1:H:15:GLU:O	2.37	0.43
1:B:61:LEU:HB3	1:B:133:VAL:HG22	2.00	0.43
1:E:19:ASP:O	1:E:20:GLN:C	2.57	0.43
1:C:280:VAL:HA	1:C:281:PRO:HD3	1.83	0.43
1:D:170:LEU:HB3	1:D:288:LEU:HD12	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:99:TRP:O	1:A:102:ARG:HB3	2.19	0.43
1:D:169:LEU:CD1	1:D:289:LYS:HG2	2.48	0.43
1:G:156:ALA:O	1:G:160:ILE:HD12	2.18	0.43
1:C:192:ILE:HD12	1:C:270:PHE:CZ	2.54	0.42
1:F:280:VAL:HA	1:F:281:PRO:HD3	1.86	0.42
1:D:99:TRP:O	1:D:102:ARG:HB3	2.18	0.42
1:E:261:ARG:HG3	1:E:261:ARG:H	1.61	0.42
1:B:5:TYR:HB3	1:C:5:TYR:HB3	2.01	0.42
1:A:84:VAL:HA	1:A:113:GLU:O	2.18	0.42
1:A:168:GLY:O	1:A:289:LYS:HA	2.18	0.42
1:E:9:SER:HB3	1:F:26:ALA:CB	2.49	0.42
1:C:71:SER:O	1:C:75:VAL:HG23	2.20	0.42
1:G:280:VAL:HA	1:G:281:PRO:HD3	1.82	0.42
1:B:99:TRP:O	1:B:102:ARG:HB3	2.20	0.42
1:F:5:TYR:HD1	1:F:5:TYR:O	2.03	0.42
1:C:261:ARG:HA	1:C:261:ARG:HD3	1.71	0.42
1:F:135:CYS:HB3	1:F:172:ILE:HG13	2.02	0.42
1:D:70:ASP:HB2	3:D:4263:HOH:O	2.18	0.42
1:A:222:VAL:HG12	1:A:223:GLN:N	2.33	0.42
1:A:266:VAL:CG2	1:A:267:LEU:N	2.82	0.42
1:E:119:THR:OG1	1:E:123:ASP:OD2	2.35	0.42
1:G:99:TRP:O	1:G:102:ARG:HB3	2.20	0.42
1:E:15:GLU:O	1:F:85:ASP:OD1	2.37	0.41
1:E:97:GLU:OE1	1:E:97:GLU:HA	2.19	0.41
1:B:84:VAL:HB	1:B:115:ALA:HB3	2.02	0.41
1:C:10:LEU:HB3	1:D:21:TYR:CG	2.54	0.41
1:D:119:THR:CB	1:D:122:LYS:NZ	2.77	0.41
1:A:84:VAL:HB	1:A:115:ALA:HB3	2.01	0.41
1:A:30:TRP:CZ2	1:A:34:ILE:HG21	2.56	0.41
1:A:255:ARG:HH21	1:A:264:HIS:CD2	2.39	0.41
1:E:268:GLY:N	1:E:271:LYS:O	2.45	0.41
1:H:200:LYS:H	1:H:200:LYS:HG2	1.74	0.41
1:A:99:TRP:CE2	1:B:98:ARG:NH2	2.88	0.41
1:G:179:TYR:CB	2:G:7282:BME:H12	2.49	0.41
1:D:149:ASP:C	1:D:151:SER:H	2.23	0.41
1:C:85:ASP:OD1	1:C:86:ALA:N	2.54	0.41
1:A:264:HIS:HB2	1:A:288:LEU:HD22	2.02	0.41
1:D:119:THR:CG2	1:D:122:LYS:HZ2	2.33	0.41
1:G:142:HIS:ND1	1:H:14:ALA:HB1	2.36	0.41
1:E:5:TYR:O	1:E:5:TYR:HD1	2.04	0.41
1:G:239:ARG:N	1:H:9:SER:OG	2.44	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:61:LEU:HB3	1:G:133:VAL:HG22	2.03	0.41
1:B:145:ASP:OD1	1:B:149:ASP:O	2.38	0.41
1:A:117:TRP:CZ3	1:B:16:GLY:CA	3.04	0.41
1:C:9:SER:HB3	1:D:26:ALA:CB	2.50	0.41
1:B:147:LYS:HD3	1:B:149:ASP:OD1	2.21	0.41
1:F:99:TRP:O	1:F:102:ARG:HB3	2.19	0.41
1:H:280:VAL:HA	1:H:281:PRO:HD3	1.83	0.41
1:B:258:PHE:CE2	1:B:288:LEU:HD12	2.55	0.41
1:H:172:ILE:O	1:H:172:ILE:CG2	2.69	0.41
1:D:85:ASP:OD1	1:D:86:ALA:N	2.54	0.41
1:C:65:CYS:HB3	1:C:85:ASP:HB2	2.02	0.41
1:E:91:LEU:O	1:E:92:LYS:C	2.58	0.41
1:D:19:ASP:O	1:D:20:GLN:C	2.58	0.41
1:B:188:PRO:HB3	1:B:202:ILE:HD12	2.03	0.41
1:E:149:ASP:C	1:E:151:SER:H	2.24	0.41
1:E:247:LEU:HD11	1:E:266:VAL:HG21	2.03	0.41
1:D:84:VAL:HA	1:D:113:GLU:O	2.20	0.41
1:D:17:LEU:HA	1:D:18:PRO:HD3	1.98	0.41
1:G:255:ARG:HH21	1:G:264:HIS:CD2	2.39	0.41
1:G:44:TYR:O	1:G:45:LYS:C	2.60	0.40
1:D:186:ALA:HA	1:D:187:PRO:HD3	1.86	0.40
1:A:30:TRP:CH2	1:A:34:ILE:HD13	2.57	0.40
1:C:17:LEU:HA	1:C:18:PRO:HD3	1.95	0.40
1:H:87:SER:OG	1:H:90:MET:HG3	2.20	0.40
1:A:190:LYS:HD3	1:A:190:LYS:HA	1.92	0.40
1:A:190:LYS:CG	1:A:282:CYS:SG	3.10	0.40
1:B:84:VAL:HA	1:B:113:GLU:O	2.21	0.40
1:B:40:ARG:CD	1:B:45:LYS:CG	2.87	0.40
1:D:134:ILE:HG13	1:D:134:ILE:O	2.21	0.40
1:D:264:HIS:ND1	1:D:288:LEU:HD21	2.37	0.40
1:F:180:ILE:HG12	1:F:186:ALA:HA	2.04	0.40
1:B:280:VAL:HA	1:B:281:PRO:HD3	1.86	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

of similar resolution.

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	277/292 (95%)	250 (90%)	22 (8%)	5 (2%)	11	45
1	B	278/292 (95%)	250 (90%)	23 (8%)	5 (2%)	11	45
1	C	279/292 (96%)	253 (91%)	20 (7%)	6 (2%)	8	38
1	D	282/292 (97%)	258 (92%)	21 (7%)	3 (1%)	17	58
1	E	282/292 (97%)	253 (90%)	24 (8%)	5 (2%)	11	45
1	F	278/292 (95%)	245 (88%)	24 (9%)	9 (3%)	5	27
1	G	279/292 (96%)	250 (90%)	24 (9%)	5 (2%)	11	45
1	H	278/292 (95%)	252 (91%)	22 (8%)	4 (1%)	14	51
All	All	2233/2336 (96%)	2011 (90%)	180 (8%)	42 (2%)	10	43

All (42) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	6	ARG
1	A	20	GLN
1	B	20	GLN
1	B	39	SER
1	C	5	TYR
1	C	20	GLN
1	C	25	GLU
1	D	5	TYR
1	D	20	GLN
1	E	20	GLN
1	F	20	GLN
1	F	39	SER
1	G	20	GLN
1	H	4	VAL
1	H	5	TYR
1	H	20	GLN
1	A	126	SER
1	B	37	THR
1	B	199	THR
1	C	37	THR
1	F	4	VAL
1	A	268	GLY
1	B	270	PHE

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Mol	Chain	Res	Type
1	C	225	PRO
1	D	270	PHE
1	E	270	PHE
1	F	191	ASN
1	F	267	LEU
1	F	270	PHE
1	G	5	TYR
1	G	270	PHE
1	A	270	PHE
1	F	165	ARG
1	F	211	ASN
1	G	291	THR
1	H	270	PHE
1	C	270	PHE
1	G	106	SER
1	E	225	PRO
1	F	5	TYR
1	E	4	VAL
1	E	268	GLY

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	230/246 (94%)	195 (85%)	35 (15%)	3	16
1	B	230/246 (94%)	193 (84%)	37 (16%)	3	14
1	C	222/246 (90%)	191 (86%)	31 (14%)	4	19
1	D	236/246 (96%)	197 (84%)	39 (16%)	3	13
1	E	231/246 (94%)	191 (83%)	40 (17%)	2	12
1	F	211/246 (86%)	178 (84%)	33 (16%)	3	16
1	G	235/246 (96%)	197 (84%)	38 (16%)	3	14
1	H	228/246 (93%)	193 (85%)	35 (15%)	3	16
All	All	1823/1968 (93%)	1535 (84%)	288 (16%)	3	15

All (288) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	VAL
1	A	5	TYR
1	A	7	THR
1	A	25	GLU
1	A	29	VAL
1	A	31	GLN
1	A	36	ASP
1	A	37	THR
1	A	38	ARG
1	A	54	GLN
1	A	69	VAL
1	A	73	MET
1	A	80	SER
1	A	81	VAL
1	A	83	SER
1	A	102	ARG
1	A	106	SER
1	A	147	LYS
1	A	158	LYS
1	A	159	ASN
1	A	162	SER
1	A	165	ARG
1	A	169	LEU
1	A	196	SER
1	A	205	SER
1	A	211	ASN
1	A	236	SER
1	A	249	SER
1	A	265	SER
1	A	266	VAL
1	A	280	VAL
1	A	287	VAL
1	A	289	LYS
1	A	291	THR
1	A	292	ASP
1	B	2	ASP
1	B	3	SER
1	B	5	TYR
1	B	6	ARG
1	B	7	THR
1	B	25	GLU
1	B	28	ARG

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Mol	Chain	Res	Type
1	B	29	VAL
1	B	31	GLN
1	B	36	ASP
1	B	38	ARG
1	B	45	LYS
1	B	69	VAL
1	B	71	SER
1	B	73	MET
1	B	80	SER
1	B	81	VAL
1	B	83	SER
1	B	102	ARG
1	B	122	LYS
1	B	147	LYS
1	B	158	LYS
1	B	162	SER
1	B	169	LEU
1	B	175	ARG
1	B	192	ILE
1	B	205	SER
1	B	224	VAL
1	B	236	SER
1	B	249	SER
1	B	261	ARG
1	B	265	SER
1	B	266	VAL
1	B	269	ASP
1	B	280	VAL
1	B	287	VAL
1	B	288	LEU
1	C	3	SER
1	C	6	ARG
1	C	7	THR
1	C	25	GLU
1	C	28	ARG
1	C	29	VAL
1	C	31	GLN
1	C	38	ARG
1	C	69	VAL
1	C	73	MET
1	C	80	SER
1	C	81	VAL

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Mol	Chain	Res	Type
1	C	83	SER
1	C	89	LYS
1	C	102	ARG
1	C	103	LYS
1	C	126	SER
1	C	162	SER
1	C	169	LEU
1	C	175	ARG
1	C	196	SER
1	C	199	THR
1	C	200	LYS
1	C	205	SER
1	C	236	SER
1	C	237	LYS
1	C	261	ARG
1	C	263	GLN
1	C	265	SER
1	C	269	ASP
1	C	287	VAL
1	D	2	ASP
1	D	3	SER
1	D	5	TYR
1	D	6	ARG
1	D	7	THR
1	D	29	VAL
1	D	31	GLN
1	D	38	ARG
1	D	54	GLN
1	D	69	VAL
1	D	73	MET
1	D	80	SER
1	D	81	VAL
1	D	83	SER
1	D	102	ARG
1	D	103	LYS
1	D	106	SER
1	D	122	LYS
1	D	125	LEU
1	D	126	SER
1	D	146	CYS
1	D	169	LEU
1	D	175	ARG

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Mol	Chain	Res	Type
1	D	190	LYS
1	D	200	LYS
1	D	205	SER
1	D	211	ASN
1	D	223	GLN
1	D	236	SER
1	D	237	LYS
1	D	261	ARG
1	D	265	SER
1	D	266	VAL
1	D	269	ASP
1	D	280	VAL
1	D	287	VAL
1	D	289	LYS
1	D	290	LYS
1	D	291	THR
1	E	2	ASP
1	E	3	SER
1	E	5	TYR
1	E	6	ARG
1	E	7	THR
1	E	17	LEU
1	E	25	GLU
1	E	28	ARG
1	E	29	VAL
1	E	31	GLN
1	E	38	ARG
1	E	45	LYS
1	E	69	VAL
1	E	73	MET
1	E	80	SER
1	E	81	VAL
1	E	83	SER
1	E	89	LYS
1	E	102	ARG
1	E	113	GLU
1	E	125	LEU
1	E	133	VAL
1	E	147	LYS
1	E	158	LYS
1	E	162	SER
1	E	165	ARG

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Mol	Chain	Res	Type
1	E	169	LEU
1	E	190	LYS
1	E	198	LEU
1	E	200	LYS
1	E	205	SER
1	E	211	ASN
1	E	236	SER
1	E	237	LYS
1	E	261	ARG
1	E	265	SER
1	E	269	ASP
1	E	271	LYS
1	E	280	VAL
1	E	287	VAL
1	F	2	ASP
1	F	3	SER
1	F	5	TYR
1	F	6	ARG
1	F	7	THR
1	F	28	ARG
1	F	29	VAL
1	F	31	GLN
1	F	36	ASP
1	F	45	LYS
1	F	69	VAL
1	F	73	MET
1	F	80	SER
1	F	81	VAL
1	F	83	SER
1	F	102	ARG
1	F	106	SER
1	F	122	LYS
1	F	147	LYS
1	F	158	LYS
1	F	162	SER
1	F	165	ARG
1	F	169	LEU
1	F	175	ARG
1	F	198	LEU
1	F	205	SER
1	F	211	ASN
1	F	236	SER

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Mol	Chain	Res	Type
1	F	265	SER
1	F	269	ASP
1	F	277	GLN
1	F	287	VAL
1	F	290	LYS
1	G	6	ARG
1	G	7	THR
1	G	17	LEU
1	G	29	VAL
1	G	31	GLN
1	G	38	ARG
1	G	54	GLN
1	G	69	VAL
1	G	73	MET
1	G	80	SER
1	G	81	VAL
1	G	83	SER
1	G	89	LYS
1	G	102	ARG
1	G	103	LYS
1	G	106	SER
1	G	122	LYS
1	G	147	LYS
1	G	158	LYS
1	G	162	SER
1	G	165	ARG
1	G	169	LEU
1	G	175	ARG
1	G	190	LYS
1	G	192	ILE
1	G	198	LEU
1	G	200	LYS
1	G	205	SER
1	G	237	LYS
1	G	261	ARG
1	G	265	SER
1	G	269	ASP
1	G	271	LYS
1	G	277	GLN
1	G	287	VAL
1	G	288	LEU
1	G	289	LYS

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Mol	Chain	Res	Type
1	G	291	THR
1	H	5	TYR
1	H	6	ARG
1	H	7	THR
1	H	25	GLU
1	H	28	ARG
1	H	29	VAL
1	H	31	GLN
1	H	38	ARG
1	H	54	GLN
1	H	69	VAL
1	H	73	MET
1	H	80	SER
1	H	81	VAL
1	H	83	SER
1	H	102	ARG
1	H	122	LYS
1	H	126	SER
1	H	159	ASN
1	H	162	SER
1	H	165	ARG
1	H	169	LEU
1	H	175	ARG
1	H	197	ASP
1	H	199	THR
1	H	200	LYS
1	H	205	SER
1	H	211	ASN
1	H	224	VAL
1	H	236	SER
1	H	237	LYS
1	H	261	ARG
1	H	269	ASP
1	H	274	LYS
1	H	287	VAL
1	H	288	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	116	ASN
1	A	159	ASN

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Mol	Chain	Res	Type
1	A	211	ASN
1	B	116	ASN
1	B	223	GLN
1	C	159	ASN
1	C	264	HIS
1	D	20	GLN
1	D	211	ASN
1	D	223	GLN
1	E	150	GLN
1	E	159	ASN
1	E	245	HIS
1	E	264	HIS
1	F	211	ASN
1	G	159	ASN
1	G	245	HIS
1	H	150	GLN
1	H	159	ASN
1	H	211	ASN
1	H	264	HIS

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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

10 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected

value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	BME	A	1262	1	3,3,3	0.57	0	2,2,2	0.51	0
2	BME	A	1282	1	3,3,3	0.93	0	2,2,2	1.78	0
2	BME	B	2057	1	3,3,3	0.47	0	2,2,2	0.39	0
2	BME	D	4185	1	3,3,3	0.76	0	2,2,2	1.32	0
2	BME	D	4246	1	3,3,3	0.28	0	2,2,2	1.03	0
2	BME	D	4262	1	3,3,3	0.31	0	2,2,2	0.24	0
2	BME	E	5246	1	3,3,3	0.36	0	2,2,2	0.49	0
2	BME	F	6282	1	3,3,3	0.34	0	2,2,2	0.40	0
2	BME	G	7246	1	3,3,3	0.78	0	2,2,2	0.82	0
2	BME	G	7282	1	3,3,3	0.46	0	2,2,2	0.73	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
2	BME	A	1262	1	-	0/1/1/1	0/0/0/0
2	BME	A	1282	1	-	0/1/1/1	0/0/0/0
2	BME	B	2057	1	-	0/1/1/1	0/0/0/0
2	BME	D	4185	1	-	0/1/1/1	0/0/0/0
2	BME	D	4246	1	-	0/1/1/1	0/0/0/0
2	BME	D	4262	1	-	0/1/1/1	0/0/0/0
2	BME	E	5246	1	-	0/1/1/1	0/0/0/0
2	BME	F	6282	1	-	0/1/1/1	0/0/0/0
2	BME	G	7246	1	-	0/1/1/1	0/0/0/0
2	BME	G	7282	1	-	0/1/1/1	0/0/0/0

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.

5 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1262	BME	2	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1282	BME	1	0
2	B	2057	BME	5	0
2	D	4185	BME	1	0
2	G	7282	BME	2	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 ⓘ

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.