



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 04:29 pm GMT

PDB ID : 1S5U
Title : Crystal Structure of Hypothetical Protein EC709 from Escherichia coli
Authors : Kim, Y.; Joachimiak, A.; Skarina, T.; Savchenko, A.; Edwards, A.; Midwest Center for Structural Genomics (MCSG)
Deposited on : 2004-01-21
Resolution : 1.70 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	:	1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix)	:	1.9-1692
EDS	:	trunk28620
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	recalc28949

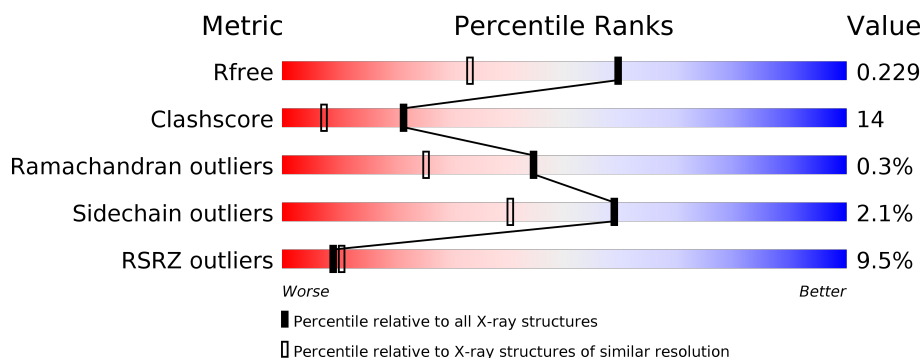
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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}	100719	3453 (1.70-1.70)
Clashscore	112137	3876 (1.70-1.70)
Ramachandran outliers	110173	3815 (1.70-1.70)
Sidechain outliers	110143	3815 (1.70-1.70)
RSRZ outliers	101464	3491 (1.70-1.70)

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.

Mol	Chain	Length	Quality of chain
1	A	138	<div> <div>8%</div> <div> <div></div> <div>73%</div> <div>18%</div> <div>• 7%</div> </div> </div>
1	B	138	<div> <div>7%</div> <div> <div></div> <div>70%</div> <div>22%</div> <div>7%</div> </div> </div>
1	C	138	<div> <div>11%</div> <div> <div></div> <div>69%</div> <div>25%</div> <div>6%</div> </div> </div>
1	D	138	<div> <div>6%</div> <div> <div></div> <div>70%</div> <div>23%</div> <div>• 7%</div> </div> </div>
1	E	138	<div> <div>14%</div> <div> <div></div> <div>62%</div> <div>36%</div> <div>•</div> </div> </div>
1	F	138	<div> <div>6%</div> <div> <div></div> <div>73%</div> <div>20%</div> <div>• 6%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
1	G	138	<div><div></div><div>10%</div><div></div><div>70%</div><div></div><div>20%</div><div></div><div>6%</div></div>
1	H	138	<div><div></div><div>11%</div><div></div><div>70%</div><div></div><div>22%</div><div></div><div>6%</div></div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9316 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 Protein ybgC.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	129	Total	C	N	O	S	0	0	0
			1051	673	182	189	7			
1	B	128	Total	C	N	O	S	0	0	0
			1044	669	181	187	7			
1	C	130	Total	C	N	O	S	0	0	0
			1060	679	184	190	7			
1	D	129	Total	C	N	O	S	0	0	0
			1051	673	182	189	7			
1	E	136	Total	C	N	O	S	0	0	0
			1106	705	194	199	8			
1	F	130	Total	C	N	O	S	0	0	0
			1060	679	184	190	7			
1	G	130	Total	C	N	O	S	0	0	0
			1058	677	183	191	7			
1	H	130	Total	C	N	O	S	0	0	0
			1058	677	183	191	7			

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	CLONING ARTIFACT	UNP P0A8Z3
A	0	HIS	-	CLONING ARTIFACT	UNP P0A8Z3
A	135	GLY	-	CLONING ARTIFACT	UNP P0A8Z3
A	136	SER	-	CLONING ARTIFACT	UNP P0A8Z3
B	-1	GLY	-	CLONING ARTIFACT	UNP P0A8Z3
B	0	HIS	-	CLONING ARTIFACT	UNP P0A8Z3
B	135	GLY	-	CLONING ARTIFACT	UNP P0A8Z3
B	136	SER	-	CLONING ARTIFACT	UNP P0A8Z3
C	-1	GLY	-	CLONING ARTIFACT	UNP P0A8Z3
C	0	HIS	-	CLONING ARTIFACT	UNP P0A8Z3
C	135	GLY	-	CLONING ARTIFACT	UNP P0A8Z3
C	136	SER	-	CLONING ARTIFACT	UNP P0A8Z3
D	-1	GLY	-	CLONING ARTIFACT	UNP P0A8Z3

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Chain	Residue	Modelled	Actual	Comment	Reference
D	0	HIS	-	CLONING ARTIFACT	UNP P0A8Z3
D	135	GLY	-	CLONING ARTIFACT	UNP P0A8Z3
D	136	SER	-	CLONING ARTIFACT	UNP P0A8Z3
E	-1	GLY	-	CLONING ARTIFACT	UNP P0A8Z3
E	0	HIS	-	CLONING ARTIFACT	UNP P0A8Z3
E	135	GLY	-	CLONING ARTIFACT	UNP P0A8Z3
E	136	SER	-	CLONING ARTIFACT	UNP P0A8Z3
F	-1	GLY	-	CLONING ARTIFACT	UNP P0A8Z3
F	0	HIS	-	CLONING ARTIFACT	UNP P0A8Z3
F	135	GLY	-	CLONING ARTIFACT	UNP P0A8Z3
F	136	SER	-	CLONING ARTIFACT	UNP P0A8Z3
G	-1	GLY	-	CLONING ARTIFACT	UNP P0A8Z3
G	0	HIS	-	CLONING ARTIFACT	UNP P0A8Z3
G	135	GLY	-	CLONING ARTIFACT	UNP P0A8Z3
G	136	SER	-	CLONING ARTIFACT	UNP P0A8Z3
H	-1	GLY	-	CLONING ARTIFACT	UNP P0A8Z3
H	0	HIS	-	CLONING ARTIFACT	UNP P0A8Z3
H	135	GLY	-	CLONING ARTIFACT	UNP P0A8Z3
H	136	SER	-	CLONING ARTIFACT	UNP P0A8Z3

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



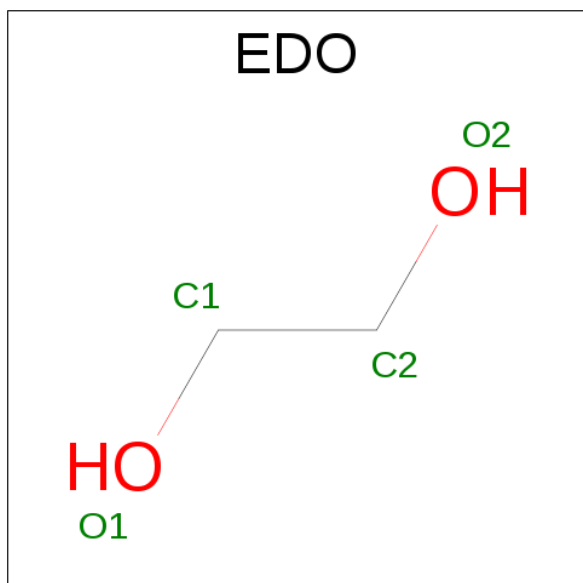
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	D	1	Total	O	S	0	0
			5	4	1		
2	G	1	Total	O	S	0	0
			5	4	1		
2	H	1	Total	O	S	0	0
			5	4	1		
2	E	1	Total	O	S	0	0
			5	4	1		
2	F	1	Total	O	S	0	0
			5	4	1		
2	F	1	Total	O	S	0	0
			5	4	1		
2	H	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).



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

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	E	1	Total C O 4 2 2	0	0
3	F	1	Total C O 4 2 2	0	0
3	A	1	Total C O 4 2 2	0	0
3	C	1	Total C O 4 2 2	0	0

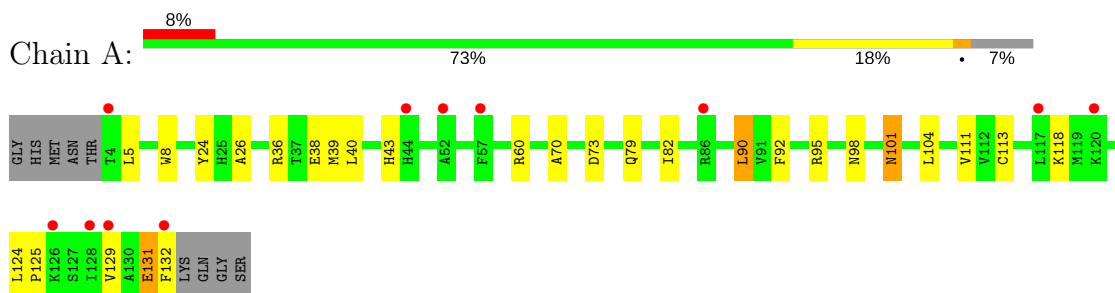
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	89	Total O 89 89	0	0
4	B	87	Total O 87 87	0	0
4	C	93	Total O 93 93	0	0
4	D	86	Total O 86 86	0	0
4	E	81	Total O 81 81	0	0
4	F	102	Total O 102 102	0	0
4	G	105	Total O 105 105	0	0
4	H	105	Total O 105 105	0	0

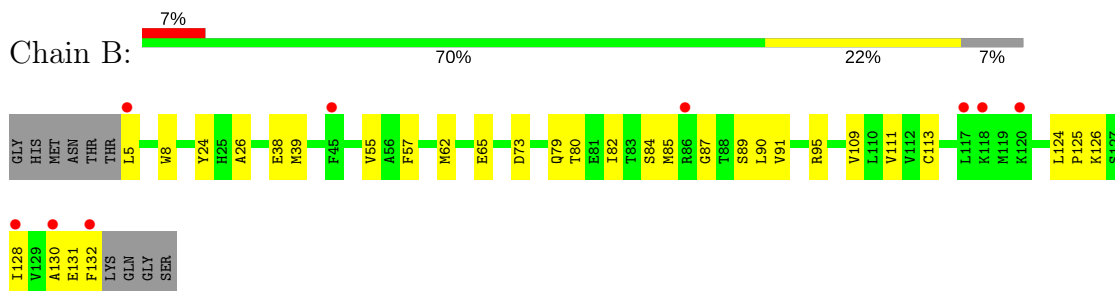
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

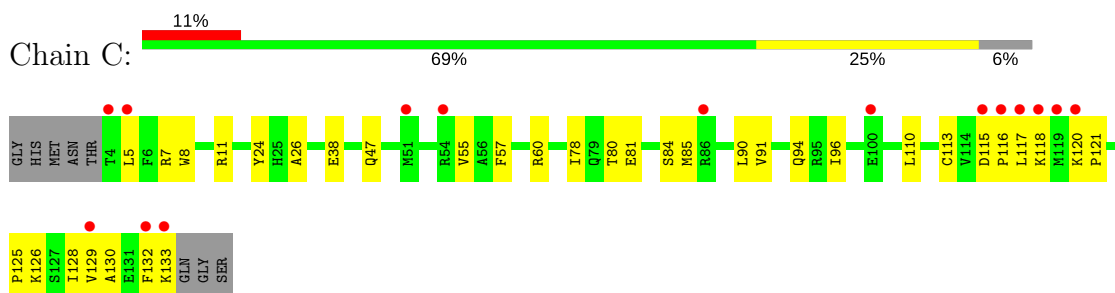
• Molecule 1: Protein ybgC



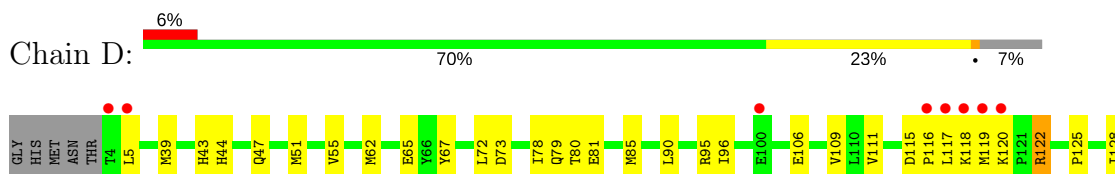
• Molecule 1: Protein ybgC



• Molecule 1: Protein ybgC

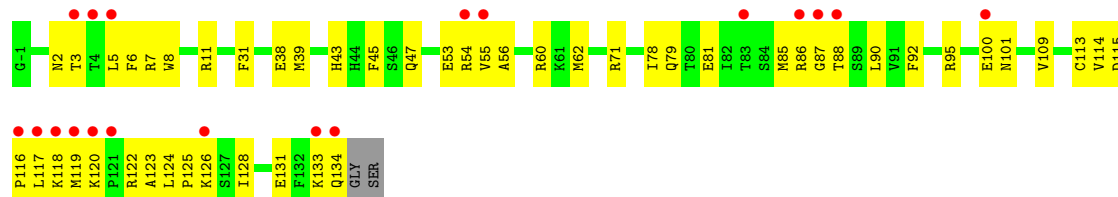


• Molecule 1: Protein ybgC

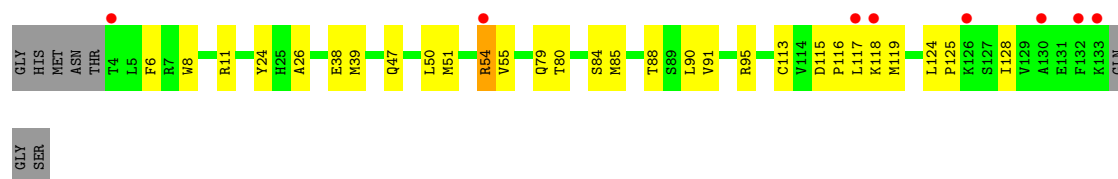




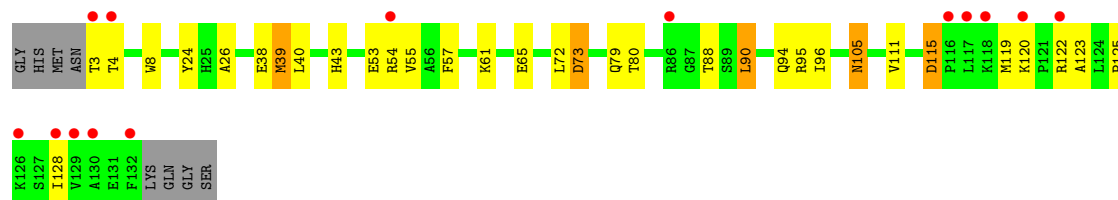
• Molecule 1: Protein ybgC



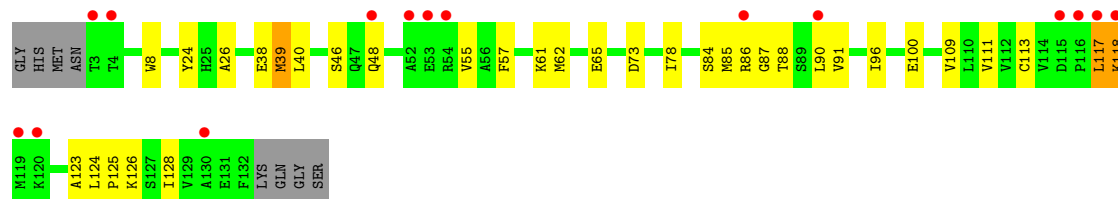
• Molecule 1: Protein ybgC



• Molecule 1: Protein ybgC



• Molecule 1: Protein ybgC



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	57.11Å 70.19Å 71.53Å 103.47° 95.15° 97.58°	Depositor
Resolution (Å)	23.64 – 1.70 24.77 – 1.70	Depositor EDS
% Data completeness (in resolution range)	86.8 (23.64-1.70) 82.3 (24.77-1.70)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.54 (at 1.71Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.200 , 0.230 0.200 , 0.229	Depositor DCC
R_{free} test set	10122 reflections (10.03%)	DCC
Wilson B-factor (Å ²)	17.2	Xtriage
Anisotropy	0.231	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 59.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.013 for -h,-l,-k	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	9316	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.62% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, EDO

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	0.33	0/1074	0.59	0/1457
1	B	0.34	0/1067	0.59	0/1447
1	C	0.33	0/1083	0.61	0/1468
1	D	0.34	0/1074	0.59	0/1457
1	E	0.34	0/1130	0.58	0/1531
1	F	0.35	0/1083	0.62	0/1468
1	G	0.33	0/1081	0.59	0/1467
1	H	0.34	0/1081	0.59	0/1467
All	All	0.34	0/8673	0.60	0/11762

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1051	0	1050	23	0
1	B	1044	0	1043	22	0
1	C	1060	0	1063	28	0
1	D	1051	0	1050	24	0
1	E	1106	0	1106	49	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	F	1060	0	1063	35	0
1	G	1058	0	1057	30	0
1	H	1058	0	1057	32	0
2	A	5	0	0	0	0
2	B	10	0	0	0	0
2	C	10	0	0	1	0
2	D	5	0	0	0	0
2	E	5	0	0	0	0
2	F	10	0	0	0	0
2	G	5	0	0	0	0
2	H	10	0	0	0	0
3	A	4	0	6	0	0
3	C	4	0	6	0	0
3	E	4	0	6	2	0
3	F	4	0	6	1	0
3	H	4	0	6	1	0
4	A	89	0	0	2	0
4	B	87	0	0	0	0
4	C	93	0	0	3	0
4	D	86	0	0	0	0
4	E	81	0	0	3	0
4	F	102	0	0	3	0
4	G	105	0	0	3	0
4	H	105	0	0	1	0
All	All	9316	0	8519	231	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 14.

All (231) 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:117:LEU:HB3	1:H:118:LYS:HE2	1.25	1.17
1:F:51:MET:HA	4:F:995:HOH:O	1.58	1.03
1:H:86:ARG:HG2	1:H:87:GLY:H	1.34	0.93
1:E:85:MET:HE3	1:E:128:ILE:HG22	1.51	0.91
1:F:55:VAL:HG11	1:F:125:PRO:HG3	1.55	0.89
1:D:122:ARG:HH11	1:D:122:ARG:HB2	1.37	0.88
1:E:55:VAL:HG21	1:E:125:PRO:HG2	1.55	0.88
1:C:55:VAL:HG21	1:C:125:PRO:CG	2.11	0.81
1:C:60:ARG:HB3	1:C:110:LEU:HD12	1.61	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:47:GLN:HB2	1:G:73:ASP:OD2	1.81	0.80
1:H:118:LYS:H	1:H:118:LYS:CE	1.96	0.77
1:C:11:ARG:HH12	3:H:901:EDO:H12	1.50	0.76
1:H:118:LYS:H	1:H:118:LYS:HE2	1.50	0.76
1:F:47:GLN:HB3	4:G:891:HOH:O	1.88	0.74
1:D:115:ASP:HB2	1:D:122:ARG:NH1	2.02	0.73
1:E:2:ASN:ND2	1:E:101:ASN:HD21	1.86	0.73
1:E:133:LYS:HB2	1:E:134:GLN:HB2	1.69	0.73
1:A:43:HIS:HB3	1:A:131:GLU:HG2	1.71	0.72
1:E:90:LEU:HD11	1:E:92:PHE:CE2	2.25	0.71
1:E:86:ARG:HG2	1:E:87:GLY:H	1.55	0.71
1:F:54:ARG:HD2	1:F:116:PRO:HD2	1.71	0.71
1:C:55:VAL:HG21	1:C:125:PRO:HG3	1.72	0.71
1:E:126:LYS:HD3	1:E:126:LYS:O	1.90	0.70
4:E:956:HOH:O	1:F:118:LYS:HD2	1.90	0.69
4:A:964:HOH:O	1:D:47:GLN:HG2	1.93	0.67
1:H:117:LEU:HD12	1:H:118:LYS:HG3	1.76	0.67
1:F:11:ARG:HH22	3:F:903:EDO:H22	1.60	0.66
1:C:129:VAL:HB	4:C:978:HOH:O	1.95	0.65
1:H:84:SER:HB2	1:H:91:VAL:HB	1.79	0.64
1:G:90:LEU:N	1:G:90:LEU:HD23	2.11	0.64
1:H:55:VAL:HG23	1:H:113:CYS:SG	2.37	0.64
1:H:85:MET:HE2	1:H:90:LEU:HD21	1.79	0.64
1:E:31:PHE:HB3	1:E:78:ILE:HD11	1.79	0.64
1:E:5:LEU:HD12	1:E:95:ARG:HH22	1.63	0.64
1:F:55:VAL:HG11	1:F:125:PRO:CG	2.27	0.64
1:C:85:MET:HE3	2:C:810:SO4:O1	1.98	0.63
1:F:116:PRO:HG2	4:F:995:HOH:O	1.97	0.63
1:E:85:MET:CE	1:E:124:LEU:HD23	2.28	0.63
1:E:47:GLN:HG2	4:H:940:HOH:O	1.98	0.63
1:H:125:PRO:HD2	1:H:128:ILE:HD12	1.81	0.63
1:H:39:MET:HG3	1:H:40:LEU:N	2.14	0.62
1:D:85:MET:HG3	1:D:90:LEU:CD2	2.29	0.62
1:G:115:ASP:OD2	1:G:120:LYS:HG2	2.00	0.62
1:A:90:LEU:CD1	1:A:111:VAL:HB	2.31	0.61
1:H:117:LEU:HB3	1:H:118:LYS:CE	2.17	0.61
1:H:85:MET:CE	1:H:90:LEU:HD21	2.30	0.61
1:D:79:GLN:OE1	1:D:95:ARG:HD2	2.01	0.60
1:F:117:LEU:O	1:F:118:LYS:HG2	2.02	0.60
1:F:54:ARG:CD	1:F:116:PRO:HD2	2.31	0.60
1:C:118:LYS:HD3	1:C:120:LYS:HE2	1.84	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:85:MET:HE2	1:E:124:LEU:HD23	1.83	0.59
1:E:11:ARG:HH22	3:E:902:EDO:H21	1.68	0.59
1:F:125:PRO:HD2	1:F:128:ILE:HD12	1.85	0.58
1:H:86:ARG:HG2	1:H:87:GLY:N	2.14	0.58
1:D:125:PRO:HD2	1:D:128:ILE:HD12	1.84	0.58
1:E:118:LYS:HG3	1:E:120:LYS:HE2	1.85	0.58
1:B:125:PRO:O	1:B:128:ILE:HG22	2.03	0.58
1:C:130:ALA:HA	1:C:133:LYS:HE2	1.86	0.57
1:A:36:ARG:O	1:A:40:LEU:HD13	2.04	0.57
1:A:90:LEU:HD13	1:A:92:PHE:CE2	2.40	0.57
1:G:96:ILE:HD12	1:G:105:ASN:HD21	1.70	0.56
1:G:54:ARG:HD2	4:G:890:HOH:O	2.05	0.56
1:E:85:MET:CE	1:E:128:ILE:HG22	2.31	0.56
1:B:130:ALA:O	1:F:117:LEU:HD23	2.06	0.56
1:F:85:MET:HG3	1:F:90:LEU:HD23	1.87	0.56
1:E:114:VAL:CG1	1:E:119:MET:HA	2.36	0.56
1:A:113:CYS:HB2	1:A:124:LEU:HD23	1.86	0.56
1:G:72:LEU:O	1:G:73:ASP:HB2	2.05	0.55
1:F:47:GLN:HG2	1:G:72:LEU:CD2	2.36	0.55
1:C:78:ILE:HD12	1:C:96:ILE:HG12	1.87	0.55
1:F:115:ASP:OD2	1:F:117:LEU:HB3	2.06	0.55
1:B:62:MET:HG3	1:B:109:VAL:HG22	1.89	0.55
1:D:55:VAL:HG12	1:D:122:ARG:HD3	1.89	0.55
1:A:39:MET:HE1	1:A:92:PHE:HE1	1.72	0.55
1:E:55:VAL:CG2	1:E:125:PRO:HG2	2.34	0.54
1:G:115:ASP:O	1:G:119:MET:HA	2.06	0.54
1:G:39:MET:HE2	1:G:40:LEU:N	2.22	0.54
1:H:85:MET:HE2	1:H:124:LEU:HD23	1.90	0.54
1:E:115:ASP:OD1	1:E:117:LEU:HD23	2.06	0.54
1:E:118:LYS:HG3	1:E:118:LYS:O	2.07	0.54
1:E:11:ARG:HH12	3:E:902:EDO:H22	1.73	0.54
1:F:88:THR:HA	1:F:124:LEU:HD13	1.88	0.54
1:C:80:THR:HG23	1:C:94:GLN:HG2	1.90	0.54
1:H:85:MET:HE1	1:H:128:ILE:HG22	1.90	0.54
1:A:125:PRO:O	1:A:129:VAL:HG23	2.08	0.54
1:F:84:SER:HB3	1:F:91:VAL:HB	1.88	0.54
1:H:118:LYS:N	1:H:118:LYS:HE2	2.19	0.54
1:F:79:GLN:OE1	1:F:95:ARG:NH2	2.41	0.53
1:C:91:VAL:HG22	1:C:110:LEU:HD23	1.91	0.53
1:G:3:THR:HG22	1:G:4:THR:N	2.25	0.52
1:D:122:ARG:HH11	1:D:122:ARG:CB	2.16	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:2:ASN:HD22	1:E:101:ASN:HD21	1.55	0.52
1:E:95:ARG:HG3	1:E:95:ARG:HH11	1.74	0.52
1:H:118:LYS:H	1:H:118:LYS:HE3	1.74	0.52
1:A:8:TRP:CE2	1:A:38:GLU:HG3	2.44	0.51
1:E:90:LEU:HD12	1:E:90:LEU:C	2.30	0.51
1:E:86:ARG:HG2	1:E:86:ARG:HH11	1.75	0.51
1:G:125:PRO:HD2	1:G:128:ILE:HD12	1.91	0.51
1:G:61:LYS:HG3	1:H:65:GLU:HG2	1.93	0.51
1:E:5:LEU:HD22	1:E:81:GLU:HB3	1.92	0.51
1:A:43:HIS:CD2	1:A:132:PHE:HE2	2.28	0.51
1:F:6:PHE:HB3	1:F:39:MET:HE3	1.93	0.51
1:G:80:THR:HG23	1:G:94:GLN:HG2	1.93	0.51
1:D:43:HIS:HD2	1:D:131:GLU:OE2	1.95	0.50
1:G:65:GLU:HG2	1:H:61:LYS:HG3	1.93	0.50
1:C:113:CYS:O	1:C:121:PRO:HA	2.11	0.50
1:E:56:ALA:HB2	1:E:116:PRO:HG3	1.93	0.50
1:E:114:VAL:HG11	1:E:119:MET:HA	1.92	0.50
1:A:39:MET:HE1	1:A:82:ILE:HD11	1.94	0.49
1:D:67:TYR:HE2	1:D:106:GLU:HG2	1.76	0.49
1:B:57:PHE:HB3	1:B:111:VAL:HG11	1.93	0.49
1:C:8:TRP:CE2	1:C:38:GLU:HG3	2.47	0.49
1:B:55:VAL:HG13	1:B:113:CYS:SG	2.52	0.49
1:C:85:MET:HE1	1:C:132:PHE:HB2	1.93	0.49
1:F:47:GLN:HG2	1:G:72:LEU:HD21	1.93	0.49
1:A:79:GLN:OE1	1:A:95:ARG:HD2	2.13	0.49
1:E:8:TRP:CE2	1:E:38:GLU:HG3	2.47	0.49
1:H:46:SER:OG	1:H:48:GLN:HG2	2.12	0.49
1:D:5:LEU:HD11	1:D:79:GLN:HB3	1.94	0.49
1:B:84:SER:HB3	1:B:91:VAL:HB	1.94	0.49
1:C:24:TYR:CE2	1:C:26:ALA:HB3	2.47	0.49
1:B:85:MET:CE	1:B:132:PHE:HB2	2.42	0.48
1:E:43:HIS:HD2	1:E:131:GLU:OE2	1.96	0.48
1:G:39:MET:HE2	1:G:40:LEU:HA	1.94	0.48
1:C:7:ARG:NE	1:H:126:LYS:NZ	2.62	0.48
1:C:7:ARG:NE	1:H:126:LYS:HZ1	2.11	0.48
1:C:115:ASP:OD2	1:C:117:LEU:HB2	2.14	0.48
1:D:118:LYS:HB3	1:D:120:LYS:HG3	1.96	0.48
1:C:85:MET:HG3	1:C:90:LEU:HD23	1.95	0.48
1:F:50:LEU:O	1:F:55:VAL:HG22	2.14	0.48
1:E:62:MET:HG3	1:E:109:VAL:HG22	1.96	0.47
1:F:39:MET:HE1	1:F:80:THR:O	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:57:PHE:HB3	1:B:111:VAL:CG1	2.44	0.47
1:B:39:MET:HE1	1:B:80:THR:HG22	1.96	0.47
1:A:8:TRP:CD2	1:A:38:GLU:HG3	2.49	0.47
1:E:53:GLU:O	1:E:55:VAL:HG23	2.14	0.47
1:B:82:ILE:HD12	1:B:132:PHE:CD2	2.50	0.47
1:B:39:MET:CE	1:B:80:THR:HG22	2.45	0.47
1:A:5:LEU:HD11	1:A:79:GLN:HB3	1.96	0.47
1:D:78:ILE:HD12	1:D:96:ILE:HG12	1.97	0.47
1:E:45:PHE:HE2	1:E:131:GLU:HG2	1.80	0.47
1:G:39:MET:HE2	1:G:40:LEU:CA	2.45	0.47
1:E:54:ARG:HE	1:E:117:LEU:HD23	1.80	0.46
1:F:117:LEU:O	1:F:118:LYS:HE2	2.15	0.46
1:H:57:PHE:HB3	1:H:111:VAL:CG1	2.45	0.46
1:B:113:CYS:HB2	1:B:124:LEU:HD23	1.96	0.46
1:H:57:PHE:HB3	1:H:111:VAL:HG11	1.97	0.46
1:A:98:ASN:HB2	4:A:919:HOH:O	2.16	0.46
1:E:85:MET:HE1	1:E:124:LEU:HD23	1.97	0.46
1:F:50:LEU:HB3	1:F:55:VAL:HG23	1.96	0.46
1:C:55:VAL:HG21	1:C:125:PRO:HG2	1.93	0.46
1:E:79:GLN:NE2	1:E:95:ARG:NH2	2.64	0.46
1:B:89:SER:O	1:B:90:LEU:HD12	2.16	0.46
1:A:39:MET:HE1	1:A:92:PHE:CE1	2.51	0.45
1:G:43:HIS:HE1	4:G:900:HOH:O	1.99	0.45
1:D:51:MET:HE3	1:D:116:PRO:HG2	1.97	0.45
1:F:113:CYS:HB2	1:F:124:LEU:HD12	1.99	0.45
1:A:60:ARG:HH12	1:B:65:GLU:CG	2.30	0.45
1:D:65:GLU:HG2	1:D:67:TYR:CZ	2.51	0.45
1:F:117:LEU:O	1:F:118:LYS:CG	2.64	0.45
1:C:84:SER:HB3	1:C:91:VAL:HB	1.98	0.44
1:G:24:TYR:CE2	1:G:26:ALA:HB3	2.52	0.44
1:H:87:GLY:O	1:H:124:LEU:HD13	2.16	0.44
1:C:125:PRO:HD2	1:C:128:ILE:HD12	1.99	0.44
1:E:100:GLU:O	1:E:101:ASN:HB2	2.18	0.44
1:E:86:ARG:NH1	1:E:86:ARG:HG2	2.33	0.44
1:E:6:PHE:HB3	1:E:39:MET:HG3	1.99	0.44
1:C:126:LYS:HA	4:C:978:HOH:O	2.17	0.44
1:E:88:THR:HG22	1:E:123:ALA:HA	2.00	0.44
1:B:79:GLN:OE1	1:B:95:ARG:HD2	2.18	0.44
1:G:8:TRP:CE2	1:G:38:GLU:HG3	2.53	0.44
1:D:39:MET:HE3	1:D:80:THR:O	2.18	0.44
1:A:90:LEU:HD12	1:A:111:VAL:HB	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:88:THR:HG22	1:G:123:ALA:HA	2.00	0.44
1:B:5:LEU:HD11	1:B:79:GLN:HB3	1.98	0.43
1:D:72:LEU:O	1:D:73:ASP:HB2	2.17	0.43
1:G:90:LEU:HD23	1:G:90:LEU:H	1.83	0.43
1:G:55:VAL:HG12	1:G:122:ARG:HD3	1.99	0.43
1:F:6:PHE:HB2	1:F:39:MET:HE2	2.00	0.43
1:G:90:LEU:CD2	1:G:90:LEU:N	2.79	0.43
1:D:39:MET:HE1	1:D:81:GLU:HA	2.00	0.43
1:E:86:ARG:HG2	1:E:87:GLY:N	2.28	0.43
1:H:24:TYR:CE2	1:H:26:ALA:HB3	2.52	0.43
1:H:88:THR:HA	1:H:124:LEU:HD13	2.01	0.43
1:C:85:MET:HG3	1:C:90:LEU:CD2	2.49	0.43
1:F:8:TRP:CE2	1:F:38:GLU:HG3	2.54	0.43
1:F:24:TYR:CE2	1:F:26:ALA:HB3	2.54	0.43
1:A:24:TYR:CE2	1:A:26:ALA:HB3	2.54	0.42
1:B:8:TRP:CE2	1:B:38:GLU:HG3	2.54	0.42
1:C:129:VAL:CB	4:C:978:HOH:O	2.58	0.42
1:E:60:ARG:HG3	1:E:60:ARG:HH11	1.84	0.42
1:F:118:LYS:HA	1:F:118:LYS:HD3	1.59	0.42
1:F:39:MET:HE3	1:F:80:THR:HB	2.02	0.42
1:E:3:THR:HG22	1:E:7:ARG:NH1	2.35	0.42
1:G:57:PHE:HB3	1:G:111:VAL:CG1	2.50	0.42
1:B:126:LYS:HD3	1:B:126:LYS:HA	1.87	0.42
1:B:55:VAL:HG21	1:B:125:PRO:HG2	2.02	0.42
1:D:90:LEU:HB2	1:D:111:VAL:HG22	2.01	0.42
1:G:53:GLU:HB2	1:G:55:VAL:HG22	2.01	0.42
1:C:7:ARG:HE	1:H:126:LYS:NZ	2.18	0.42
1:F:47:GLN:HG2	1:G:72:LEU:HD23	2.01	0.42
1:F:54:ARG:HG3	4:F:1005:HOH:O	2.19	0.42
1:E:55:VAL:HG13	1:E:113:CYS:SG	2.59	0.42
1:E:3:THR:HG21	1:E:79:GLN:HG2	2.02	0.42
1:E:71:ARG:NH2	4:E:956:HOH:O	2.52	0.41
1:C:5:LEU:HA	1:C:81:GLU:HG2	2.02	0.41
1:G:39:MET:C	1:G:39:MET:HE2	2.41	0.41
1:H:78:ILE:HD12	1:H:96:ILE:HG12	2.02	0.41
1:G:39:MET:CE	1:G:40:LEU:HD23	2.50	0.41
1:A:101:ASN:HD22	1:A:101:ASN:HA	1.75	0.41
1:E:31:PHE:HB3	1:E:78:ILE:CD1	2.48	0.41
1:B:24:TYR:CE2	1:B:26:ALA:HB3	2.55	0.41
1:D:62:MET:HG3	1:D:109:VAL:HG22	2.01	0.41
1:A:118:LYS:HB3	1:A:118:LYS:HE2	1.88	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:60:ARG:HH12	1:B:65:GLU:HG2	1.85	0.41
1:E:114:VAL:O	1:E:116:PRO:HD3	2.20	0.41
1:H:8:TRP:CE2	1:H:38:GLU:HG3	2.56	0.41
1:B:85:MET:HG3	1:B:90:LEU:CD1	2.51	0.41
1:A:70:ALA:HB2	1:A:104:LEU:HD13	2.02	0.41
1:E:113:CYS:SG	1:E:122:ARG:HG3	2.60	0.41
1:D:55:VAL:HG12	1:D:122:ARG:CD	2.50	0.40
1:E:71:ARG:NE	4:E:956:HOH:O	2.52	0.40
1:A:43:HIS:ND1	1:A:131:GLU:HG3	2.36	0.40
1:D:43:HIS:O	1:D:44:HIS:HB2	2.21	0.40
1:H:88:THR:HG22	1:H:123:ALA:HA	2.03	0.40
1:H:62:MET:HG3	1:H:109:VAL:HG22	2.03	0.40
1:C:57:PHE:HZ	1:C:128:ILE:HD13	1.86	0.40
1:G:79:GLN:OE1	1:G:95:ARG:HD2	2.21	0.40
1:D:117:LEU:C	1:D:117:LEU:HD13	2.42	0.40
1:D:118:LYS:O	1:D:119:MET:HB2	2.22	0.40
1:F:117:LEU:HD12	1:F:117:LEU:HA	1.94	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	127/138 (92%)	125 (98%)	2 (2%)	0	100	100
1	B	126/138 (91%)	122 (97%)	3 (2%)	1 (1%)	22	7
1	C	128/138 (93%)	124 (97%)	3 (2%)	1 (1%)	22	7
1	D	127/138 (92%)	124 (98%)	3 (2%)	0	100	100
1	E	134/138 (97%)	129 (96%)	5 (4%)	0	100	100
1	F	128/138 (93%)	125 (98%)	2 (2%)	1 (1%)	22	7
1	G	128/138 (93%)	124 (97%)	4 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	H	128/138 (93%)	123 (96%)	5 (4%)	0	100	100
All	All	1026/1104 (93%)	996 (97%)	27 (3%)	3 (0%)	44	25

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	F	119	MET
1	B	87	GLY
1	C	116	PRO

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	113/120 (94%)	109 (96%)	4 (4%)	41	19
1	B	112/120 (93%)	110 (98%)	2 (2%)	64	47
1	C	114/120 (95%)	113 (99%)	1 (1%)	82	74
1	D	113/120 (94%)	112 (99%)	1 (1%)	82	74
1	E	119/120 (99%)	119 (100%)	0	100	100
1	F	114/120 (95%)	113 (99%)	1 (1%)	82	74
1	G	114/120 (95%)	109 (96%)	5 (4%)	33	13
1	H	114/120 (95%)	109 (96%)	5 (4%)	33	13
All	All	913/960 (95%)	894 (98%)	19 (2%)	59	40

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	73	ASP
1	A	90	LEU
1	A	101	ASN
1	A	131	GLU
1	B	73	ASP

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Mol	Chain	Res	Type
1	B	131	GLU
1	C	47	GLN
1	D	122	ARG
1	F	54	ARG
1	G	39	MET
1	G	73	ASP
1	G	90	LEU
1	G	105	ASN
1	G	115	ASP
1	H	39	MET
1	H	73	ASP
1	H	100	GLU
1	H	117	LEU
1	H	118	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	101	ASN
1	B	47	GLN
1	C	47	GLN
1	D	43	HIS
1	D	47	GLN
1	E	2	ASN
1	E	43	HIS
1	E	47	GLN
1	E	79	GLN
1	F	48	GLN
1	G	47	GLN
1	G	105	ASN
1	H	47	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 [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

17 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	SO4	A	801	-	4,4,4	0.33	0	6,6,6	0.06	0
3	EDO	A	904	-	3,3,3	0.54	0	2,2,2	0.44	0
2	SO4	B	802	-	4,4,4	0.30	0	6,6,6	0.07	0
2	SO4	B	811	-	4,4,4	0.31	0	6,6,6	0.07	0
2	SO4	C	810	-	4,4,4	0.30	0	6,6,6	0.09	0
2	SO4	C	812	-	4,4,4	0.31	0	6,6,6	0.08	0
3	EDO	C	905	-	3,3,3	0.61	0	2,2,2	0.43	0
2	SO4	D	803	-	4,4,4	0.31	0	6,6,6	0.05	0
2	SO4	E	806	-	4,4,4	0.30	0	6,6,6	0.06	0
3	EDO	E	902	-	3,3,3	0.64	0	2,2,2	0.48	0
2	SO4	F	807	-	4,4,4	0.33	0	6,6,6	0.12	0
2	SO4	F	808	-	4,4,4	0.32	0	6,6,6	0.07	0
3	EDO	F	903	-	3,3,3	0.62	0	2,2,2	0.38	0
2	SO4	G	804	-	4,4,4	0.30	0	6,6,6	0.08	0
2	SO4	H	805	-	4,4,4	0.33	0	6,6,6	0.09	0
2	SO4	H	809	-	4,4,4	0.32	0	6,6,6	0.07	0
3	EDO	H	901	-	3,3,3	0.58	0	2,2,2	0.51	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	SO4	A	801	-	-	0/0/0/0	0/0/0/0
3	EDO	A	904	-	-	0/1/1/1	0/0/0/0
2	SO4	B	802	-	-	0/0/0/0	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	SO4	B	811	-	-	0/0/0/0	0/0/0/0
2	SO4	C	810	-	-	0/0/0/0	0/0/0/0
2	SO4	C	812	-	-	0/0/0/0	0/0/0/0
3	EDO	C	905	-	-	0/1/1/1	0/0/0/0
2	SO4	D	803	-	-	0/0/0/0	0/0/0/0
2	SO4	E	806	-	-	0/0/0/0	0/0/0/0
3	EDO	E	902	-	-	0/1/1/1	0/0/0/0
2	SO4	F	807	-	-	0/0/0/0	0/0/0/0
2	SO4	F	808	-	-	0/0/0/0	0/0/0/0
3	EDO	F	903	-	-	0/1/1/1	0/0/0/0
2	SO4	G	804	-	-	0/0/0/0	0/0/0/0
2	SO4	H	805	-	-	0/0/0/0	0/0/0/0
2	SO4	H	809	-	-	0/0/0/0	0/0/0/0
3	EDO	H	901	-	-	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.

4 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	810	SO4	1	0
3	E	902	EDO	2	0
3	F	903	EDO	1	0
3	H	901	EDO	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	129/138 (93%)	0.44	11 (8%) 11 13	13, 22, 39, 45	0
1	B	128/138 (92%)	0.44	9 (7%) 17 20	13, 23, 40, 44	0
1	C	130/138 (94%)	0.77	15 (11%) 5 7	13, 24, 41, 48	0
1	D	129/138 (93%)	0.33	8 (6%) 21 24	11, 19, 36, 44	0
1	E	136/138 (98%)	0.77	19 (13%) 3 4	12, 24, 42, 48	0
1	F	130/138 (94%)	0.41	8 (6%) 21 24	10, 19, 37, 49	0
1	G	130/138 (94%)	0.47	14 (10%) 6 7	10, 19, 39, 42	0
1	H	130/138 (94%)	0.60	15 (11%) 5 7	11, 20, 42, 46	0
All	All	1042/1104 (94%)	0.53	99 (9%) 9 10	10, 21, 41, 49	0

All (99) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	117	LEU	11.3
1	E	117	LEU	10.3
1	C	117	LEU	9.0
1	H	3	THR	8.7
1	E	118	LYS	8.7
1	F	117	LEU	8.1
1	D	4	THR	8.0
1	H	117	LEU	7.7
1	G	117	LEU	7.1
1	C	116	PRO	6.4
1	C	54	ARG	6.0
1	H	54	ARG	5.8
1	C	118	LYS	5.7
1	F	133	LYS	5.6
1	H	86	ARG	5.5
1	H	52	ALA	5.4

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Mol	Chain	Res	Type	RSRZ
1	B	128	ILE	5.3
1	A	4	THR	5.1
1	B	130	ALA	5.0
1	G	120	LYS	5.0
1	E	120	LYS	4.9
1	D	118	LYS	4.9
1	C	119	MET	4.8
1	E	119	MET	4.7
1	C	120	LYS	4.6
1	E	134	GLN	4.5
1	H	120	LYS	4.5
1	G	118	LYS	4.4
1	B	117	LEU	4.4
1	A	86	ARG	4.4
1	E	4	THR	4.4
1	A	117	LEU	4.2
1	C	86	ARG	4.2
1	F	118	LYS	4.2
1	C	4	THR	4.1
1	F	54	ARG	4.1
1	F	132	PHE	4.0
1	G	86	ARG	3.8
1	H	118	LYS	3.8
1	A	120	LYS	3.7
1	H	4	THR	3.6
1	E	121	PRO	3.5
1	E	86	ARG	3.5
1	A	132	PHE	3.4
1	B	86	ARG	3.4
1	H	48	GLN	3.3
1	C	133	LYS	3.3
1	G	128	ILE	3.3
1	E	100	GLU	3.3
1	E	133	LYS	3.3
1	G	126	LYS	3.2
1	E	116	PRO	3.2
1	C	129	VAL	3.2
1	G	54	ARG	3.2
1	G	3	THR	3.1
1	D	119	MET	3.1
1	E	83	THR	3.1
1	D	120	LYS	3.0

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Mol	Chain	Res	Type	RSRZ
1	E	5	LEU	3.0
1	E	87	GLY	3.0
1	H	130	ALA	3.0
1	F	130	ALA	2.9
1	C	100	GLU	2.9
1	B	132	PHE	2.9
1	E	88	THR	2.9
1	A	126	LYS	2.8
1	G	130	ALA	2.8
1	A	129	VAL	2.8
1	E	54	ARG	2.7
1	E	3	THR	2.7
1	C	5	LEU	2.7
1	G	4	THR	2.6
1	D	100	GLU	2.6
1	G	129	VAL	2.6
1	C	132	PHE	2.5
1	A	128	ILE	2.5
1	A	44	HIS	2.5
1	C	51	MET	2.4
1	C	115	ASP	2.4
1	G	116	PRO	2.4
1	A	52	ALA	2.4
1	H	119	MET	2.4
1	H	90	LEU	2.3
1	G	122	ARG	2.3
1	F	126	LYS	2.3
1	B	120	LYS	2.3
1	B	118	LYS	2.3
1	H	116	PRO	2.2
1	F	4	THR	2.2
1	B	45	PHE	2.2
1	H	53	GLU	2.2
1	D	116	PRO	2.2
1	G	132	PHE	2.2
1	E	126	LYS	2.1
1	A	57	PHE	2.1
1	H	115	ASP	2.1
1	E	55	VAL	2.1
1	B	5	LEU	2.1
1	D	5	LEU	2.0

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
2	SO4	E	806	5/5	0.94	0.22	0.74	47,47,48,48	4
2	SO4	G	804	5/5	0.96	0.10	0.70	32,36,37,38	3
2	SO4	H	805	5/5	0.96	0.09	0.24	31,33,35,36	1
3	EDO	H	901	4/4	0.87	0.14	-0.06	28,34,34,37	0
2	SO4	C	810	5/5	0.95	0.17	-0.20	18,29,32,32	3
2	SO4	B	802	5/5	0.96	0.09	-0.37	38,39,40,40	3
2	SO4	F	807	5/5	0.95	0.10	-0.60	36,37,39,40	2
2	SO4	H	809	5/5	0.96	0.11	-1.14	48,49,50,51	5
2	SO4	C	812	5/5	0.93	0.20	-	45,46,47,48	3
3	EDO	E	902	4/4	0.85	0.14	-	23,29,30,33	1
3	EDO	C	905	4/4	0.87	0.13	-	33,33,33,34	2
2	SO4	F	808	5/5	0.93	0.18	-	48,48,49,49	5
2	SO4	B	811	5/5	0.97	0.18	-	37,40,41,42	3
2	SO4	A	801	5/5	0.94	0.10	-	37,39,40,42	3
2	SO4	D	803	5/5	0.84	0.26	-	48,48,49,50	3
3	EDO	A	904	4/4	0.91	0.14	-	26,31,32,34	1
3	EDO	F	903	4/4	0.89	0.11	-	25,29,29,30	1

6.5 Other polymers [i](#)

There are no such residues in this entry.