



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 23, 2017 – 04:47 PM EDT

PDB ID : 3H5R
Title : Crystal structure of E. coli MccB + Succinimide
Authors : Regni, C.A.; Roush, R.F.; Miller, D.; Nourse, A.; Walsh, C.T.; Schulman, B.A.
Deposited on : unknown
Resolution : 2.10 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 : rb-20030345
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) : rb-20030345

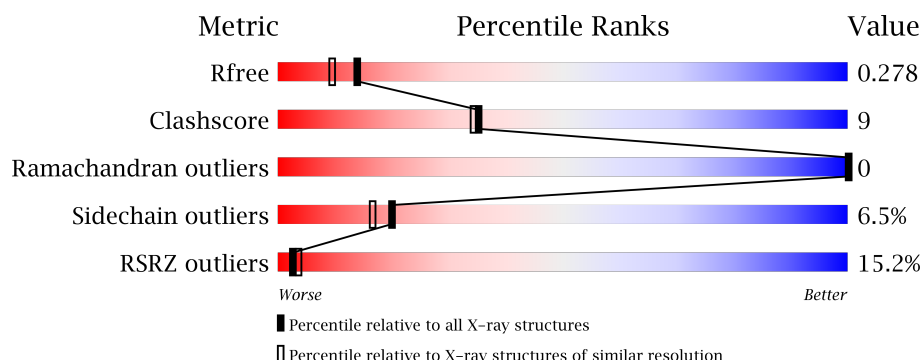
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	4243 (2.10-2.10)
Clashscore	112137	4788 (2.10-2.10)
Ramachandran outliers	110173	4740 (2.10-2.10)
Sidechain outliers	110143	4741 (2.10-2.10)
RSRZ outliers	101464	4275 (2.10-2.10)

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	353	<div> <div>14%</div> <div> <div>81%</div> <div>14%</div> <div>• •</div> </div> </div>
1	B	353	<div> <div>14%</div> <div> <div>81%</div> <div>12%</div> <div>• 5%</div> </div> </div>
1	C	353	<div> <div>11%</div> <div> <div>77%</div> <div>18%</div> <div>• •</div> </div> </div>
1	D	353	<div> <div>17%</div> <div> <div>70%</div> <div>22%</div> <div>• 6%</div> </div> </div>
2	E	7	<div> <div>29%</div> <div> <div>57%</div> <div>43%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
2	F	7	<div> <div>57%</div> <div> <div></div> <div>57%</div> <div>43%</div> </div> </div>
2	G	7	<div> <div>14%</div> <div> <div></div> <div>14%</div> <div>86%</div> </div> </div>
2	H	7	<div> <div>14%</div> <div> <div></div> <div>14%</div> <div>86%</div> </div> </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	SNN	F	77	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 10712 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 MccB protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	340	Total	C	N	O	S	0	0	0
			2599	1659	449	480	11			
1	B	334	Total	C	N	O	S	0	0	0
			2551	1636	427	477	11			
1	C	343	Total	C	N	O	S	0	0	0
			2621	1675	446	489	11			
1	D	333	Total	C	N	O	S	0	0	0
			2546	1627	434	474	11			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	EXPRESSION TAG	UNP Q47506
A	-1	SER	-	EXPRESSION TAG	UNP Q47506
A	0	HIS	-	EXPRESSION TAG	UNP Q47506
B	-2	GLY	-	EXPRESSION TAG	UNP Q47506
B	-1	SER	-	EXPRESSION TAG	UNP Q47506
B	0	HIS	-	EXPRESSION TAG	UNP Q47506
C	-2	GLY	-	EXPRESSION TAG	UNP Q47506
C	-1	SER	-	EXPRESSION TAG	UNP Q47506
C	0	HIS	-	EXPRESSION TAG	UNP Q47506
D	-2	GLY	-	EXPRESSION TAG	UNP Q47506
D	-1	SER	-	EXPRESSION TAG	UNP Q47506
D	0	HIS	-	EXPRESSION TAG	UNP Q47506

- Molecule 2 is a protein called Microcin C7 analog.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	7	Total	C	N	O	S	0	0	0
			51	28	12	10	1			
2	F	7	Total	C	N	O	S	0	0	0
			51	28	12	10	1			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	G	1	Total	C	N	O	S	0	0	0
			8	5	1	1	1			
2	H	1	Total	C	N	O	S	0	0	0
			8	5	1	1	1			

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Zn	0	0
			1	1		
3	A	1	Total	Zn	0	0
			1	1		
3	D	1	Total	Zn	0	0
			1	1		
3	C	1	Total	Zn	0	0
			1	1		

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

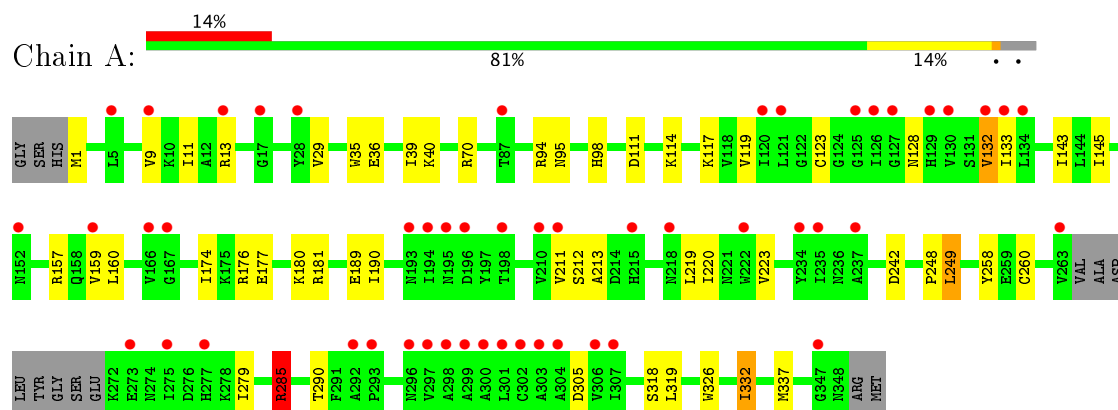
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	82	Total 82	O 82	0	0
5	B	70	Total 70	O 70	0	0
5	C	69	Total 69	O 69	0	0
5	D	38	Total 38	O 38	0	0
5	E	1	Total 1	O 1	0	0
5	F	3	Total 3	O 3	0	0

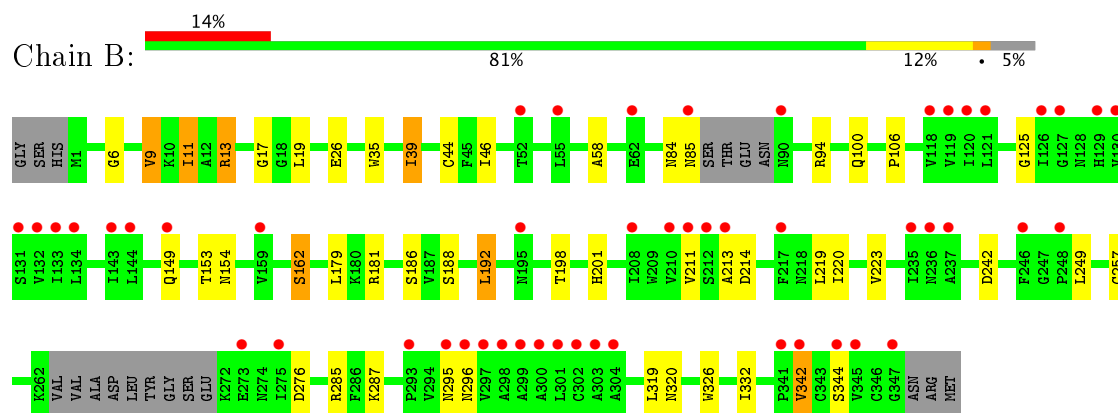
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

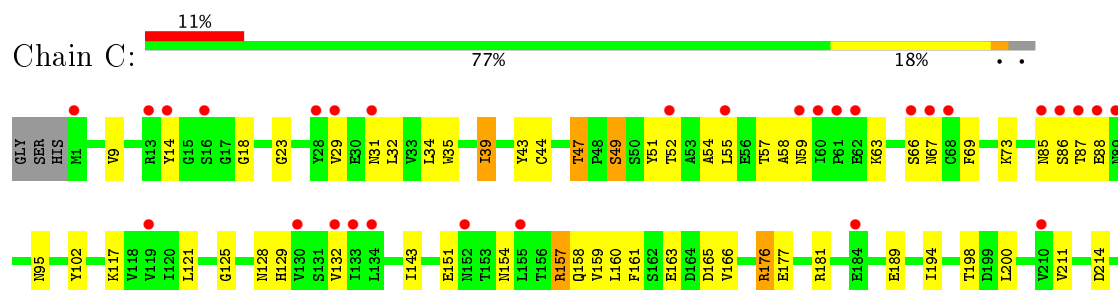
• Molecule 1: MccB protein

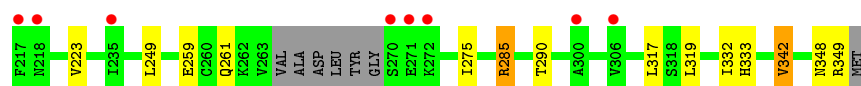


• Molecule 1: MccB protein

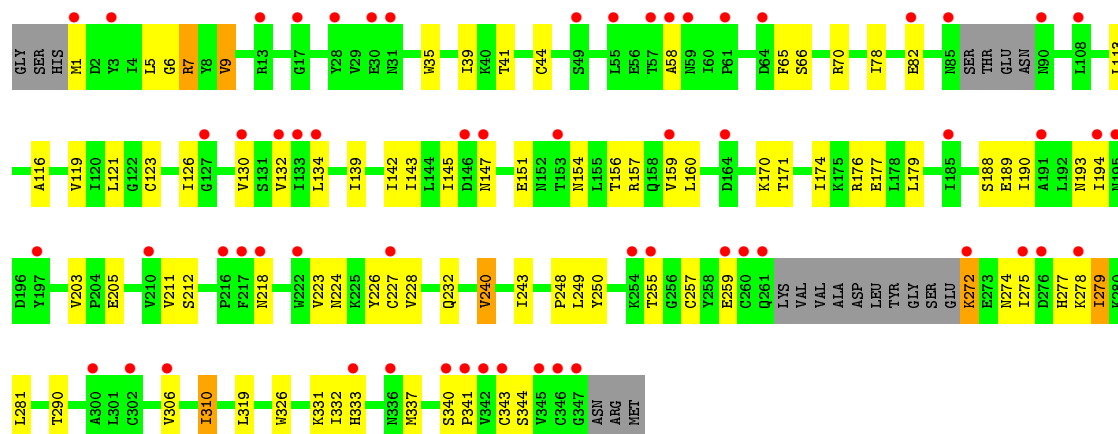


• Molecule 1: MccB protein

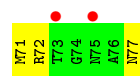




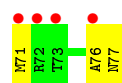
• Molecule 1: MccB protein



• Molecule 2: Microcin C7 analog



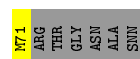
• Molecule 2: Microcin C7 analog



• Molecule 2: Microcin C7 analog



• Molecule 2: Microcin C7 analog



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	55.93Å 137.97Å 80.13Å 90.00° 92.10° 90.00°	Depositor
Resolution (Å)	27.95 – 2.10 27.39 – 2.09	Depositor EDS
% Data completeness (in resolution range)	99.9 (27.95-2.10) 99.1 (27.39-2.09)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.35 (at 2.08Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.194 , 0.250 0.236 , 0.278	Depositor DCC
R_{free} test set	3547 reflections (5.03%)	DCC
Wilson B-factor (Å ²)	37.6	Xtriage
Anisotropy	0.258	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 46.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.034 for h,-k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	10712	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.34% 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: SNN, ZN, SO4

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	0.69	0/2653	0.77	3/3607 (0.1%)
1	B	0.69	0/2603	0.74	1/3540 (0.0%)
1	C	0.64	0/2675	0.69	0/3637
1	D	0.73	5/2598 (0.2%)	0.69	1/3534 (0.0%)
2	E	0.45	0/42	0.70	0/54
2	F	0.45	0/42	0.89	0/54
2	G	0.71	0/7	1.07	0/7
2	H	0.73	0/7	0.44	0/7
All	All	0.69	5/10627 (0.0%)	0.72	5/14440 (0.0%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	272	LYS	N-CA	11.13	1.68	1.46
1	D	205	GLU	CD-OE1	7.53	1.33	1.25
1	D	205	GLU	CD-OE2	5.63	1.31	1.25
1	D	250	TYR	C-O	5.22	1.33	1.23
1	D	250	TYR	C-N	5.04	1.45	1.34

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	285	ARG	NE-CZ-NH2	-7.45	116.58	120.30
1	D	226	TYR	CB-CG-CD2	-7.04	116.77	121.00
1	A	332	ILE	CG1-CB-CG2	-6.91	96.21	111.40
1	A	249	LEU	CA-CB-CG	6.51	130.28	115.30
1	B	192	LEU	CA-CB-CG	5.88	128.81	115.30

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	A	2599	0	2526	46	0
1	B	2551	0	2480	41	0
1	C	2621	0	2551	56	0
1	D	2546	0	2471	69	0
2	E	51	0	46	4	0
2	F	51	0	46	14	0
2	G	8	0	8	0	0
2	H	8	0	8	5	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	B	5	0	0	0	0
4	C	5	0	0	0	0
5	A	82	0	0	11	0
5	B	70	0	0	7	0
5	C	69	0	0	10	0
5	D	38	0	0	2	0
5	E	1	0	0	0	0
5	F	3	0	0	0	0
All	All	10712	0	10136	192	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 (192) 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:D:272:LYS:N	1:D:272:LYS:CA	1.68	1.52
1:C:157:ARG:HD3	1:C:290:THR:HG23	1.31	1.09
1:C:157:ARG:HD3	1:C:290:THR:CG2	1.85	1.04
1:A:181:ARG:CZ	5:A:402:HOH:O	2.11	0.97
1:D:171:THR:CG2	1:D:189:GLU:HB3	2.00	0.92
1:A:332:ILE:HG13	1:B:332:ILE:HG13	1.52	0.90
1:B:186:SER:HB2	5:D:364:HOH:O	1.76	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:157:ARG:NH2	5:D:385:HOH:O	2.11	0.84
1:D:171:THR:HG21	1:D:189:GLU:HB3	1.59	0.82
1:C:157:ARG:NH1	5:C:408:HOH:O	2.12	0.82
1:C:55:LEU:HA	5:C:392:HOH:O	1.79	0.80
1:A:285:ARG:HD3	1:B:46:ILE:HG23	1.61	0.79
1:B:214:ASP:HA	2:F:77:SNN:C3	2.12	0.78
1:B:214:ASP:N	2:F:77:SNN:H42	1.98	0.78
1:C:177:GLU:OE1	1:C:181:ARG:NH2	2.18	0.76
1:A:290:THR:OG1	1:B:94:ARG:CB	2.35	0.75
1:A:159:VAL:O	1:B:181:ARG:NH1	2.20	0.74
1:D:333:HIS:HD2	2:H:71:MET:HE1	1.52	0.74
1:A:117:LYS:HE3	1:A:143:ILE:HD11	1.69	0.72
1:D:171:THR:HG22	1:D:189:GLU:OE1	1.90	0.72
1:C:47:THR:HB	5:C:364:HOH:O	1.89	0.72
1:C:157:ARG:HD3	1:C:290:THR:HG21	1.72	0.72
1:B:213:ALA:O	2:F:77:SNN:H3	1.89	0.71
1:D:248:PRO:HG3	1:D:337:MET:CE	2.21	0.71
1:D:333:HIS:HD2	2:H:71:MET:CE	2.04	0.70
1:D:333:HIS:CD2	2:H:71:MET:HE1	2.27	0.70
1:A:176:ARG:NH1	1:A:177:GLU:OE2	2.25	0.70
1:A:248:PRO:HG3	1:A:337:MET:HE1	1.74	0.69
1:B:211:VAL:HG21	1:B:223:VAL:HG11	1.76	0.68
1:B:214:ASP:HA	2:F:77:SNN:N3	2.08	0.68
1:C:157:ARG:NH2	5:C:362:HOH:O	2.27	0.67
1:C:157:ARG:CD	1:C:290:THR:HG23	2.18	0.67
1:C:58:ALA:HB3	5:C:392:HOH:O	1.94	0.67
1:D:132:VAL:HG11	1:D:160:LEU:HD21	1.76	0.67
1:C:285:ARG:HG2	1:D:7:ARG:HD2	1.77	0.66
1:D:113:LEU:HA	1:D:310:ILE:HD11	1.77	0.66
1:A:258:TYR:HA	1:A:337:MET:HE1	1.77	0.66
1:A:117:LYS:HE3	1:A:143:ILE:CD1	2.25	0.65
1:A:94:ARG:O	5:A:384:HOH:O	2.13	0.65
1:C:39:ILE:HD11	1:D:278:LYS:HB3	1.78	0.65
1:C:128:ASN:O	1:C:132:VAL:HG23	1.98	0.64
1:A:213:ALA:HB3	1:A:219:LEU:HD21	1.80	0.64
1:A:248:PRO:HG3	1:A:337:MET:CE	2.26	0.64
1:A:119:VAL:HG13	1:A:145:ILE:HD12	1.80	0.63
1:A:145:ILE:HG12	1:A:190:ILE:HB	1.81	0.63
1:A:211:VAL:HG21	1:A:223:VAL:HG11	1.81	0.62
1:C:333:HIS:NE2	5:C:400:HOH:O	2.19	0.62
1:D:248:PRO:HG3	1:D:337:MET:HE2	1.82	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:132:VAL:HG22	1:C:160:LEU:HD11	1.83	0.61
1:D:227:CYS:HB3	1:D:232:GLN:O	2.01	0.60
1:D:147:ASN:HB3	1:D:193:ASN:ND2	2.16	0.60
1:D:333:HIS:CD2	2:H:71:MET:CE	2.85	0.59
1:B:213:ALA:HB3	1:B:219:LEU:HD22	1.85	0.58
1:C:332:ILE:HD11	1:D:332:ILE:HG13	1.85	0.58
1:D:272:LYS:C	1:D:272:LYS:N	2.56	0.57
1:D:6:GLY:O	1:D:9:VAL:HG13	2.04	0.57
1:A:181:ARG:NE	5:A:402:HOH:O	2.29	0.57
1:D:116:ALA:HB2	1:D:310:ILE:HD13	1.87	0.56
1:C:158:GLN:HE21	1:C:161:PHE:HE2	1.54	0.56
1:B:213:ALA:C	2:F:77:SNN:H3	2.26	0.56
1:A:114:LYS:HE3	5:A:406:HOH:O	2.06	0.56
5:B:389:HOH:O	2:F:76:ALA:HB2	2.06	0.55
1:A:258:TYR:HA	1:A:337:MET:CE	2.36	0.55
1:A:70:ARG:NH1	5:A:422:HOH:O	2.39	0.55
5:B:389:HOH:O	2:F:76:ALA:CB	2.54	0.55
1:D:151:GLU:H	1:D:154:ASN:ND2	2.04	0.55
1:C:165:ASP:OD1	1:C:176:ARG:NH2	2.39	0.55
1:A:181:ARG:HD3	5:B:395:HOH:O	2.07	0.55
1:C:342:VAL:O	1:C:342:VAL:HG13	2.07	0.54
1:A:305:ASP:OD2	1:A:318:SER:OG	2.20	0.54
1:C:177:GLU:HB3	1:C:181:ARG:HH21	1.72	0.54
1:D:275:ILE:O	1:D:279:ILE:HG23	2.07	0.54
1:C:58:ALA:CB	5:C:392:HOH:O	2.54	0.54
1:B:26:GLU:OE1	2:E:72:ARG:HB2	2.07	0.54
1:A:35:TRP:CZ2	1:A:39:ILE:HD11	2.43	0.54
1:B:214:ASP:N	2:F:77:SNN:C4	2.71	0.54
1:B:6:GLY:O	1:B:9:VAL:HG13	2.08	0.54
1:C:342:VAL:O	1:C:342:VAL:CG1	2.55	0.53
1:A:181:ARG:HH11	1:A:181:ARG:HG2	1.74	0.53
1:D:326:TRP:CZ2	1:D:331:LYS:HE2	2.44	0.53
1:C:151:GLU:H	1:C:154:ASN:ND2	2.07	0.52
1:A:181:ARG:NH2	5:A:402:HOH:O	2.35	0.52
1:D:113:LEU:HD23	1:D:310:ILE:CG1	2.40	0.52
1:D:211:VAL:HG21	1:D:223:VAL:HG11	1.90	0.52
1:C:39:ILE:HD11	1:D:278:LYS:CB	2.39	0.52
1:D:248:PRO:HG3	1:D:337:MET:HE1	1.90	0.52
1:B:214:ASP:HA	2:F:77:SNN:C4	2.39	0.52
1:D:171:THR:CG2	1:D:189:GLU:CB	2.83	0.52
1:A:111:ASP:CB	5:A:396:HOH:O	2.58	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:163:GLU:O	1:C:166:VAL:HG12	2.10	0.52
1:B:84:ASN:O	1:B:85:ASN:C	2.49	0.51
1:D:113:LEU:HD23	1:D:310:ILE:HG13	1.91	0.51
1:D:310:ILE:O	1:D:310:ILE:HD12	2.10	0.51
1:C:211:VAL:HG21	1:C:223:VAL:HG11	1.92	0.50
1:C:158:GLN:NE2	1:C:161:PHE:HE2	2.10	0.50
1:A:180:LYS:HB3	1:B:162:SER:OG	2.12	0.50
1:C:194:ILE:N	1:C:194:ILE:HD12	2.26	0.50
1:D:134:LEU:HB3	1:D:139:ILE:HG13	1.94	0.50
1:A:279:ILE:HG13	1:B:39:ILE:CD1	2.43	0.49
1:B:214:ASP:HA	2:F:77:SNN:H3	1.91	0.49
1:C:44:CYS:SG	1:C:58:ALA:HB2	2.51	0.49
5:B:378:HOH:O	2:E:71:MET:N	2.46	0.49
1:B:213:ALA:C	2:F:77:SNN:H42	2.33	0.49
1:D:171:THR:HG22	1:D:189:GLU:HB3	1.91	0.49
1:A:279:ILE:HG23	1:B:11:ILE:HG22	1.95	0.48
1:D:145:ILE:CD1	1:D:203:VAL:HG13	2.44	0.48
1:D:145:ILE:HG12	1:D:190:ILE:HB	1.94	0.48
1:C:121:LEU:HD11	1:C:194:ILE:HD13	1.95	0.48
1:A:95:ASN:O	1:A:98:HIS:HB3	2.13	0.48
1:D:132:VAL:CG1	1:D:160:LEU:HD21	2.42	0.48
1:C:117:LYS:HE3	1:C:143:ILE:HD11	1.94	0.48
1:A:143:ILE:HD12	1:A:143:ILE:N	2.29	0.47
5:A:431:HOH:O	1:B:154:ASN:HA	2.13	0.47
1:A:94:ARG:NH2	5:A:431:HOH:O	2.32	0.47
1:B:201:HIS:HB3	1:C:261:GLN:OE1	2.15	0.47
1:C:285:ARG:O	1:D:7:ARG:HG3	2.14	0.47
1:A:177:GLU:O	1:A:181:ARG:HG3	2.14	0.47
1:B:295:ASN:ND2	5:B:390:HOH:O	2.47	0.47
1:D:119:VAL:HG13	1:D:145:ILE:HD12	1.96	0.46
1:D:126:ILE:O	1:D:130:VAL:HG23	2.14	0.46
1:D:274:ASN:HA	1:D:277:HIS:HD2	1.81	0.46
5:C:357:HOH:O	1:D:156:THR:HG22	2.16	0.46
1:A:133:ILE:HD11	1:B:296:ASN:HD22	1.80	0.46
1:D:194:ILE:HD12	1:D:194:ILE:N	2.31	0.46
1:A:123:CYS:HB3	1:A:174:ILE:HD12	1.98	0.46
1:D:176:ARG:NH1	1:D:177:GLU:OE2	2.48	0.46
1:D:44:CYS:SG	1:D:58:ALA:HB2	2.55	0.46
1:A:212:SER:O	2:E:77:SNN:H42	2.16	0.46
1:C:129:HIS:CE1	1:C:159:VAL:HB	2.51	0.46
1:A:133:ILE:HD11	1:B:296:ASN:ND2	2.31	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:143:ILE:N	1:D:143:ILE:HD12	2.31	0.45
1:C:31:ASN:HD22	1:C:34:LEU:HB2	1.81	0.45
1:C:51:TYR:HH	1:C:66:SER:HG	1.64	0.45
1:D:123:CYS:HB3	1:D:174:ILE:HD12	1.98	0.44
1:D:41:THR:HG23	1:D:65:PHE:CE1	2.52	0.44
1:B:125:GLY:HA3	1:B:295:ASN:HD21	1.81	0.44
1:D:340:SER:HA	1:D:341:PRO:HD2	1.86	0.44
1:B:13:ARG:HD3	1:B:17:GLY:O	2.18	0.44
1:A:36:GLU:O	1:A:40:LYS:HG3	2.18	0.44
5:A:421:HOH:O	1:B:285:ARG:HD3	2.17	0.44
1:D:66:SER:O	1:D:70:ARG:HG3	2.18	0.44
1:D:255:THR:HG21	1:D:319:LEU:HD22	2.00	0.44
1:D:143:ILE:HG13	1:D:188:SER:HB2	2.00	0.44
1:C:49:SER:HB2	1:C:54:ALA:HB2	1.98	0.43
1:C:121:LEU:HD11	1:C:194:ILE:CD1	2.48	0.43
1:C:102:TYR:OH	5:C:388:HOH:O	2.18	0.43
1:C:95:ASN:ND2	1:D:156:THR:HG21	2.34	0.43
1:C:43:TYR:CE1	1:D:281:LEU:HD23	2.53	0.43
1:A:132:VAL:HG11	1:A:160:LEU:HD21	2.01	0.43
1:B:44:CYS:SG	1:B:58:ALA:HB2	2.57	0.43
1:A:94:ARG:HG2	1:B:153:THR:OG1	2.18	0.43
1:C:125:GLY:HA2	1:C:158:GLN:HG2	2.01	0.43
1:D:154:ASN:ND2	1:D:170:LYS:HE2	2.33	0.43
1:D:240:VAL:HG21	2:H:71:MET:HA	2.01	0.43
1:C:29:VAL:HG13	1:C:35:TRP:HB2	2.01	0.43
1:C:348:ASN:O	1:C:349:ARG:CB	2.67	0.43
1:D:121:LEU:HD11	1:D:194:ILE:HD11	2.00	0.43
1:D:224:ASN:O	1:D:228:VAL:HG23	2.19	0.43
1:B:19:LEU:HD11	2:E:72:ARG:HB3	2.01	0.43
1:C:332:ILE:CD1	1:D:332:ILE:HG13	2.48	0.42
1:C:35:TRP:O	1:C:39:ILE:HG23	2.19	0.42
1:C:69:PHE:CE2	1:C:73:LYS:HD2	2.54	0.42
1:C:23:GLY:HA2	1:D:243:ILE:HD12	2.00	0.42
1:B:342:VAL:O	1:B:342:VAL:CG2	2.67	0.42
1:A:285:ARG:HD2	5:B:352:HOH:O	2.17	0.42
1:B:326:TRP:CH2	2:F:71:MET:HG3	2.54	0.42
1:A:285:ARG:CD	1:B:46:ILE:HG23	2.42	0.42
1:B:214:ASP:CA	2:F:77:SNN:C4	2.97	0.42
1:B:332:ILE:HG21	1:B:332:ILE:HD13	1.68	0.42
1:D:5:LEU:HA	1:D:78:ILE:HG22	2.00	0.42
1:A:94:ARG:NH1	5:A:370:HOH:O	2.51	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:14:TYR:H	1:C:18:GLY:HA2	1.84	0.42
1:A:128:ASN:O	1:A:132:VAL:HG22	2.20	0.42
1:C:52:THR:HA	1:C:55:LEU:HB3	2.02	0.41
1:A:123:CYS:HB3	1:A:174:ILE:CD1	2.50	0.41
1:A:160:LEU:O	1:A:177:GLU:HG3	2.20	0.41
1:D:151:GLU:O	1:D:154:ASN:HB2	2.20	0.41
1:C:43:TYR:HE1	1:D:281:LEU:HD23	1.86	0.41
5:B:389:HOH:O	2:F:76:ALA:HB3	2.20	0.41
1:C:86:SER:O	1:C:88:GLU:N	2.54	0.41
1:D:274:ASN:HA	1:D:277:HIS:CD2	2.56	0.41
1:C:39:ILE:CD1	1:D:278:LYS:HB3	2.48	0.41
1:D:306:VAL:O	1:D:310:ILE:HG23	2.20	0.41
1:B:35:TRP:O	1:B:39:ILE:HG23	2.21	0.41
1:D:78:ILE:HG13	1:D:78:ILE:O	2.20	0.41
1:C:63:LYS:CB	5:C:386:HOH:O	2.68	0.41
1:D:35:TRP:CZ2	1:D:39:ILE:HD11	2.56	0.41
1:D:340:SER:HB3	1:D:343:CYS:HB2	2.01	0.41
1:B:100:GLN:HG3	1:B:106:PRO:HG3	2.03	0.40
1:B:319:LEU:O	1:B:320:ASN:HB2	2.20	0.40
1:C:181:ARG:NH1	1:D:159:VAL:O	2.55	0.40
1:C:194:ILE:HG12	1:C:200:LEU:HD23	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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	336/353 (95%)	325 (97%)	11 (3%)	0	100	100
1	B	328/353 (93%)	318 (97%)	10 (3%)	0	100	100
1	C	339/353 (96%)	329 (97%)	10 (3%)	0	100	100
1	D	327/353 (93%)	317 (97%)	10 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	E	4/7 (57%)	4 (100%)	0	0	100	100
2	F	4/7 (57%)	4 (100%)	0	0	100	100
All	All	1338/1426 (94%)	1297 (97%)	41 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	272/304 (90%)	257 (94%)	15 (6%)	25	22
1	B	269/304 (88%)	251 (93%)	18 (7%)	19	15
1	C	277/304 (91%)	255 (92%)	22 (8%)	14	10
1	D	268/304 (88%)	252 (94%)	16 (6%)	22	19
2	E	4/4 (100%)	4 (100%)	0	100	100
2	F	4/4 (100%)	4 (100%)	0	100	100
2	G	1/4 (25%)	1 (100%)	0	100	100
2	H	1/4 (25%)	1 (100%)	0	100	100
All	All	1096/1232 (89%)	1025 (94%)	71 (6%)	20	16

All (71) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	9	VAL
1	A	11	ILE
1	A	13	ARG
1	A	29	VAL
1	A	132	VAL
1	A	157	ARG
1	A	189	GLU
1	A	220	ILE

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Mol	Chain	Res	Type
1	A	242	ASP
1	A	249	LEU
1	A	260	CYS
1	A	285	ARG
1	A	319	LEU
1	A	326	TRP
1	B	9	VAL
1	B	11	ILE
1	B	13	ARG
1	B	39	ILE
1	B	149	GLN
1	B	162	SER
1	B	179	LEU
1	B	188	SER
1	B	192	LEU
1	B	198	THR
1	B	220	ILE
1	B	242	ASP
1	B	249	LEU
1	B	257	CYS
1	B	276	ASP
1	B	287	LYS
1	B	342	VAL
1	B	344	SER
1	C	9	VAL
1	C	32	LEU
1	C	39	ILE
1	C	47	THR
1	C	49	SER
1	C	57	THR
1	C	59	ASN
1	C	67	ASN
1	C	85	ASN
1	C	87	THR
1	C	157	ARG
1	C	176	ARG
1	C	189	GLU
1	C	198	THR
1	C	214	ASP
1	C	249	LEU
1	C	259	GLU
1	C	275	ILE

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Mol	Chain	Res	Type
1	C	285	ARG
1	C	317	LEU
1	C	319	LEU
1	C	342	VAL
1	D	1	MET
1	D	7	ARG
1	D	9	VAL
1	D	82	GLU
1	D	142	ILE
1	D	179	LEU
1	D	212	SER
1	D	218	ASN
1	D	240	VAL
1	D	249	LEU
1	D	257	CYS
1	D	259	GLU
1	D	279	ILE
1	D	290	THR
1	D	310	ILE
1	D	344	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	295	ASN
1	A	296	ASN
1	B	295	ASN
1	C	31	ASN
1	C	59	ASN
1	C	129	HIS
1	C	154	ASN
1	D	95	ASN
1	D	154	ASN
1	D	236	ASN
1	D	277	HIS
1	D	333	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues 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	SNN	E	77	2	7,8,8	1.68	2 (28%)	5,11,11	2.04	2 (40%)
2	SNN	F	77	2	7,8,8	1.69	2 (28%)	5,11,11	2.03	2 (40%)

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	SNN	E	77	2	-	0/0/12/12	0/1/1/1
2	SNN	F	77	2	-	0/0/12/12	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	77	SNN	C2-N1	-3.55	1.33	1.37
2	E	77	SNN	C2-N1	-3.53	1.33	1.37
2	F	77	SNN	C5-N1	-2.58	1.33	1.37
2	E	77	SNN	C5-N1	-2.54	1.33	1.37

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	77	SNN	O2-C2-C3	-3.37	123.82	126.23
2	F	77	SNN	O2-C2-C3	-3.34	123.84	126.23
2	F	77	SNN	O5-C5-C4	-2.88	122.51	126.45
2	E	77	SNN	O5-C5-C4	-2.86	122.53	126.45

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	77	SNN	1	0
2	F	77	SNN	10	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	SO4	B	351	-	4,4,4	0.21	0	6,6,6	0.45	0
4	SO4	C	351	-	4,4,4	0.20	0	6,6,6	0.18	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
4	SO4	B	351	-	-	0/0/0/0	0/0/0/0
4	SO4	C	351	-	-	0/0/0/0	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.

No monomer is involved in short contacts.

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	A	340/353 (96%)	0.99	51 (15%) 3 4	31, 43, 60, 69	0
1	B	334/353 (94%)	0.95	51 (15%) 2 3	33, 44, 57, 71	0
1	C	343/353 (97%)	0.80	38 (11%) 6 7	28, 43, 58, 69	0
1	D	333/353 (94%)	1.04	60 (18%) 2 2	24, 44, 56, 65	0
2	E	6/7 (85%)	1.85	2 (33%) 0 0	62, 63, 65, 67	0
2	F	6/7 (85%)	2.38	4 (66%) 0 0	70, 74, 77, 80	0
2	G	1/7 (14%)	2.39	1 (100%) 0 0	82, 82, 82, 82	0
2	H	1/7 (14%)	1.62	0 100 100	83, 83, 83, 83	0
All	All	1364/1440 (94%)	0.96	207 (15%) 2 3	24, 44, 60, 83	0

All (207) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	89	ASN	8.2
1	D	342	VAL	7.1
1	D	341	PRO	7.0
1	D	347	GLY	6.2
1	B	130	VAL	6.1
1	D	336	ASN	5.8
1	D	345	VAL	5.7
1	B	210	VAL	5.3
1	B	299	ALA	4.8
1	B	302	CYS	4.7
1	A	235	ILE	4.6
1	C	28	TYR	4.5
1	D	340	SER	4.4
1	A	210	VAL	4.4
1	B	217	PHE	4.4
1	A	126	ILE	4.4

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Mol	Chain	Res	Type	RSRZ
1	A	133	ILE	4.4
1	C	60	ILE	4.4
1	C	217	PHE	4.4
1	B	235	ILE	4.4
1	A	13	ARG	4.4
1	A	195	ASN	4.3
1	A	196	ASP	4.3
1	B	211	VAL	4.3
1	B	300	ALA	4.2
1	D	343	CYS	4.2
1	C	87	THR	4.1
1	A	130	VAL	4.1
1	C	62	GLU	4.1
1	B	303	ALA	4.1
1	B	127	GLY	4.1
2	E	73	THR	4.1
1	B	132	VAL	4.0
1	C	306	VAL	3.9
1	D	346	CYS	3.9
1	B	133	ILE	3.9
1	D	133	ILE	3.8
1	B	298	ALA	3.7
1	D	28	TYR	3.7
1	D	55	LEU	3.7
1	D	130	VAL	3.7
1	B	344	SER	3.7
1	D	195	ASN	3.7
1	A	299	ALA	3.7
1	C	16	SER	3.6
1	B	301	LEU	3.6
1	B	237	ALA	3.6
1	B	297	VAL	3.6
1	A	222	TRP	3.6
1	D	194	ILE	3.6
1	B	120	ILE	3.5
1	A	292	ALA	3.5
1	A	215	HIS	3.5
1	D	1	MET	3.5
1	D	222	TRP	3.5
1	B	342	VAL	3.4
1	A	211	VAL	3.4
1	C	14	TYR	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	306	VAL	3.4
1	D	216	PRO	3.4
1	D	300	ALA	3.4
1	C	61	PRO	3.4
1	B	134	LEU	3.4
1	A	127	GLY	3.4
1	A	218	ASN	3.3
1	A	300	ALA	3.3
1	B	131	SER	3.3
1	A	193	ASN	3.3
1	B	347	GLY	3.3
2	F	72	ARG	3.3
1	D	132	VAL	3.3
1	B	62	GLU	3.2
1	B	119	VAL	3.2
1	A	296	ASN	3.2
1	B	121	LEU	3.2
1	B	159	VAL	3.2
1	A	303	ALA	3.2
1	C	130	VAL	3.1
1	A	132	VAL	3.1
1	A	304	ALA	3.1
1	B	275	ILE	3.1
1	C	133	ILE	3.1
1	A	347	GLY	3.0
1	C	132	VAL	3.0
1	A	263	VAL	3.0
1	A	134	LEU	3.0
1	A	277	HIS	3.0
1	D	82	GLU	3.0
1	B	341	PRO	3.0
1	C	52	THR	2.9
2	E	75	ASN	2.9
1	A	297	VAL	2.9
1	B	126	ILE	2.9
1	C	1	MET	2.9
1	A	5	LEU	2.9
1	A	129	HIS	2.9
1	D	259	GLU	2.9
1	B	345	VAL	2.9
1	C	85	ASN	2.9
1	D	90	ASN	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	120	ILE	2.9
1	D	17	GLY	2.9
1	B	296	ASN	2.9
2	F	71	MET	2.9
1	D	227	CYS	2.9
1	C	210	VAL	2.8
1	C	13	ARG	2.8
1	C	300	ALA	2.8
1	D	153	THR	2.8
1	C	218	ASN	2.8
1	B	273	GLU	2.8
1	C	134	LEU	2.7
1	D	127	GLY	2.7
1	D	147	ASN	2.7
1	D	260	CYS	2.7
2	F	73	THR	2.7
1	A	198	THR	2.7
1	C	59	ASN	2.7
1	A	298	ALA	2.7
1	D	254	LYS	2.6
1	D	261	GLN	2.6
1	B	55	LEU	2.6
1	D	197	TYR	2.6
1	A	302	CYS	2.6
1	B	212	SER	2.6
1	A	9	VAL	2.6
1	A	28	TYR	2.6
1	D	218	ASN	2.6
1	C	272	LYS	2.6
1	B	129	HIS	2.6
1	D	210	VAL	2.5
1	D	108	LEU	2.5
1	D	49	SER	2.5
1	D	57	THR	2.5
1	B	304	ALA	2.5
1	A	234	TYR	2.5
1	D	30	GLU	2.5
1	D	185	ILE	2.5
1	D	3	TYR	2.5
1	C	152	ASN	2.5
1	A	301	LEU	2.5
1	A	166	VAL	2.4

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Mol	Chain	Res	Type	RSRZ
1	D	58	ALA	2.4
1	B	246	PHE	2.4
1	D	31	ASN	2.4
1	D	164	ASP	2.4
1	D	306	VAL	2.4
1	A	307	ILE	2.4
1	B	295	ASN	2.4
1	D	278	LYS	2.4
1	B	85	ASN	2.4
1	B	90	ASN	2.4
2	G	71	MET	2.4
1	A	121	LEU	2.4
1	C	155	LEU	2.4
1	B	149	GLN	2.4
1	B	236	ASN	2.4
1	C	270	SER	2.4
1	D	61	PRO	2.4
1	C	235	ILE	2.4
1	D	134	LEU	2.3
1	C	271	GLU	2.3
1	B	213	ALA	2.3
1	B	118	VAL	2.3
1	A	17	GLY	2.3
1	B	144	LEU	2.3
2	F	76	ALA	2.3
1	B	293	PRO	2.3
1	D	59	ASN	2.3
1	D	85	ASN	2.3
1	B	208	ILE	2.3
1	D	275	ILE	2.3
1	A	293	PRO	2.2
1	C	66	SER	2.2
1	C	86	SER	2.2
1	D	302	CYS	2.2
1	C	67	ASN	2.2
1	D	272	LYS	2.2
1	C	119	VAL	2.2
1	A	87	THR	2.2
1	A	125	GLY	2.2
1	D	13	ARG	2.2
1	A	273	GLU	2.2
1	A	275	ILE	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	146	ASP	2.1
1	B	143	ILE	2.1
1	C	29	VAL	2.1
1	A	237	ALA	2.1
1	C	88	GLU	2.1
1	B	248	PRO	2.1
1	C	31	ASN	2.1
1	D	64	ASP	2.1
1	D	276	ASP	2.1
1	D	159	VAL	2.1
1	D	217	PHE	2.1
1	C	68	CYS	2.1
1	A	194	ILE	2.1
1	A	167	GLY	2.1
1	A	159	VAL	2.1
1	D	255	THR	2.1
1	C	55	LEU	2.0
1	B	52	THR	2.0
1	D	191	ALA	2.0
1	D	333	HIS	2.0
1	A	152	ASN	2.0
1	B	195	ASN	2.0
1	C	184	GLU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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	SNN	E	77	8/8	0.83	0.26	-	63,65,65,67	0
2	SNN	F	77	8/8	0.57	0.38	-	63,65,67,68	0

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
3	ZN	A	500	1/1	0.97	0.07	-1.59	50,50,50,50	0
3	ZN	B	500	1/1	0.98	0.04	-1.83	45,45,45,45	0
3	ZN	C	500	1/1	0.99	0.06	-2.23	37,37,37,37	0
3	ZN	D	500	1/1	0.94	0.04	-2.51	71,71,71,71	0
4	SO4	B	351	5/5	0.92	0.12	-	66,67,68,69	0
4	SO4	C	351	5/5	0.91	0.11	-	85,85,85,85	0

6.5 Other polymers [i](#)

There are no such residues in this entry.