



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2014 – 01:13 AM GMT

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 : 2009-04-22
Resolution : 2.10 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

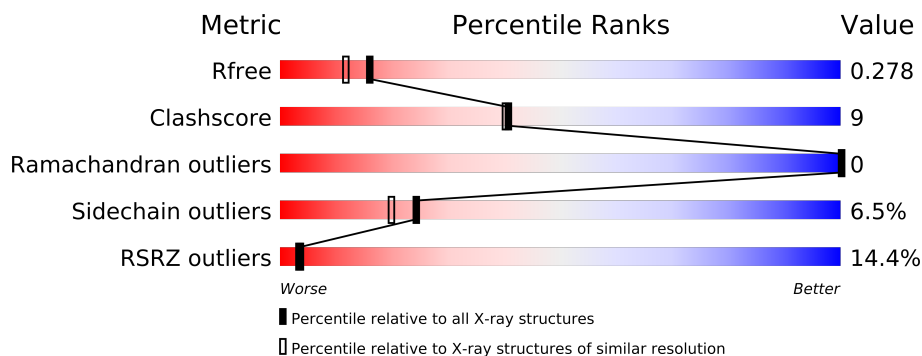
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

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}	66092	3012 (2.10-2.10)
Clashscore	79885	3649 (2.10-2.10)
Ramachandran outliers	78287	3610 (2.10-2.10)
Sidechain outliers	78261	3611 (2.10-2.10)
RSRZ outliers	66119	3013 (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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	353	
1	B	353	
1	C	353	
1	D	353	
2	E	7	
2	F	7	
2	G	7	
2	H	7	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 10712 atoms, of which 0 are hydrogen and 0 are deuterium.

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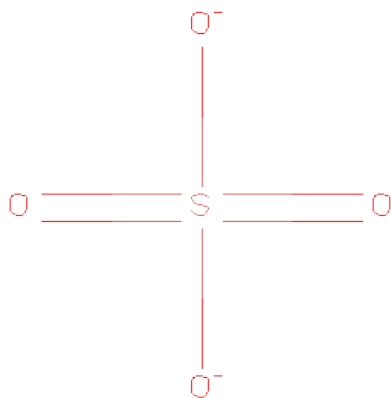
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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	G	1	Total	C	N	O	S	0	0	0
			8	5	1	1	1			
2	H	1	Total	C	N	O	S	0	0	0
			8	5	1	1	1			

- 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.

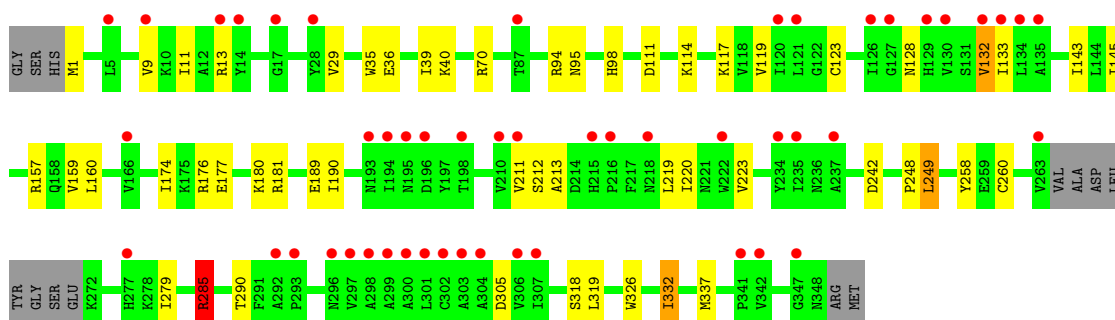
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	82	Total 82	O 82	0	0
5	B	70	Total 70	O 70	0	0
5	C	69	Total 69	O 69	0	0
5	D	38	Total 38	O 38	0	0
5	E	1	Total 1	O 1	0	0
5	F	3	Total 3	O 3	0	0

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 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.

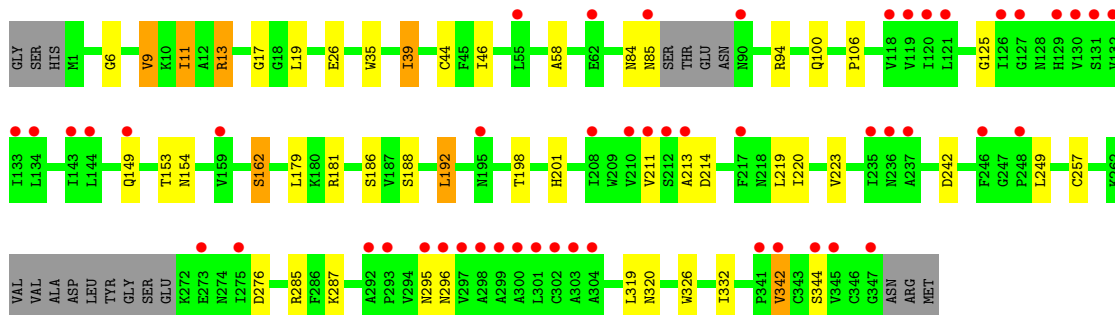
• Molecule 1: MccB protein

Chain A: 



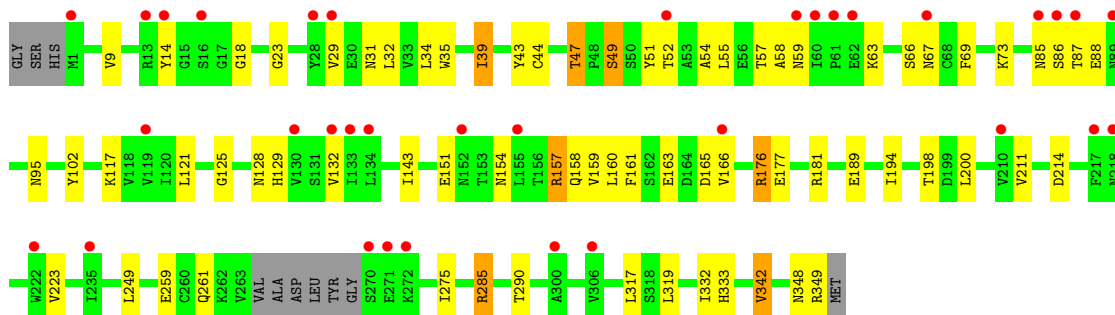
• Molecule 1: MccB protein

Chain B: 



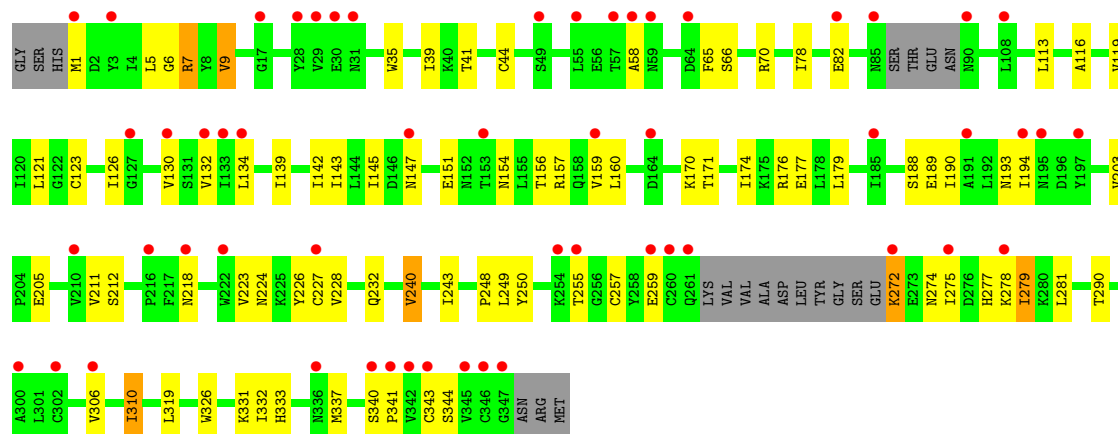
• Molecule 1: MccB protein

Chain C: 



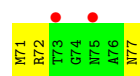
- Molecule 1: MccB protein

Chain D:



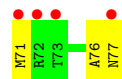
- Molecule 2: Microcin C7 analog

Chain E:



- Molecule 2: Microcin C7 analog

Chain F:



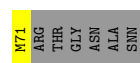
- Molecule 2: Microcin C7 analog

Chain G:



- Molecule 2: Microcin C7 analog

Chain H:



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 5.2.0019	Depositor
R, R_{free}	0.194 , 0.250 0.237 , 0.278	Depositor DCC
R_{free} test set	3597 reflections (5.04%)	DCC
Wilson B-factor (Å ²)	37.6	Xtriage
Anisotropy	0.258	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 45.6	EDS
Estimated twinning fraction	0.034 for h,-k,-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	2 of 71396 reflections (0.003%)	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.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

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 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

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

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 9.

All (192) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:272:LYS:CA	1:D:272:LYS:N	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	Distance(Å)	Clash(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:248:PRO:HG3	1:A:337:MET:CE	2.26	0.64
1:A:213:ALA:HB3	1:A:219:LEU:HD21	1.80	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
1:C:132:VAL:HG22	1:C:160:LEU:HD11	1.83	0.61

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
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:CG1	1:C:342:VAL:O	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:248:PRO:HG3	1:D:337:MET:HE1	1.90	0.52
1:C:39:ILE:HD11	1:D:278:LYS:CB	2.39	0.52
1:D:211:VAL:HG21	1:D:223:VAL:HG11	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:C:163:GLU:O	1:C:166:VAL:HG12	2.10	0.52
1:A:111:ASP:CB	5:A:396:HOH:O	2.58	0.52
1:B:84:ASN:O	1:B:85:ASN:C	2.49	0.51

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
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:C:194:ILE:N	1:C:194:ILE:HD12	2.26	0.50
1:A:180:LYS:HB3	1:B:162:SER:OG	2.12	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:C:44:CYS:SG	1:C:58:ALA:HB2	2.51	0.49
1:B:214:ASP:HA	2:F:77:SNN:H3	1.91	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:C:285:ARG:O	1:D:7:ARG:HG3	2.14	0.47
1:B:201:HIS:HB3	1:C:261:GLN:OE1	2.15	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:274:ASN:HA	1:D:277:HIS:HD2	1.81	0.46
1:D:126:ILE:O	1:D:130:VAL:HG23	2.14	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:C:129:HIS:CE1	1:C:159:VAL:HB	2.51	0.46
1:A:212:SER:O	2:E:77:SNN:H42	2.16	0.46
1:A:133:ILE:HD11	1:B:296:ASN:ND2	2.31	0.45
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

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
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
1:D:66:SER:O	1:D:70:ARG:HG3	2.18	0.44
5:A:421:HOH:O	1:B:285:ARG:HD3	2.17	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:95:ASN:ND2	1:D:156:THR:HG21	2.34	0.43
1:C:102:TYR:OH	5:C:388:HOH:O	2.18	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: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:C:348:ASN:O	1:C:349:ARG:CB	2.67	0.43
1:B:19:LEU:HD11	2:E:72:ARG:HB3	2.01	0.43
1:C:35:TRP:O	1:C:39:ILE:HG23	2.19	0.42
1:C:332:ILE:CD1	1:D:332:ILE:HG13	2.48	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:A:285:ARG:CD	1:B:46:ILE:HG23	2.42	0.42
1:B:326:TRP:CH2	2:F:71:MET:HG3	2.54	0.42
1:B:332:ILE:HG21	1:B:332:ILE:HD13	1.68	0.42
1:B:214:ASP:CA	2:F:77:SNN:C4	2.97	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
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:160:LEU:O	1:A:177:GLU:HG3	2.20	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:123:CYS:HB3	1:A:174:ILE:CD1	2.50	0.41
1:D:151:GLU:O	1:D:154:ASN:HB2	2.20	0.41
5:B:389:HOH:O	2:F:76:ALA:HB3	2.20	0.41
1:C:43:TYR:HE1	1:D:281:LEU:HD23	1.86	0.41
1:D:306:VAL:O	1:D:310:ILE:HG23	2.20	0.41
1:C:39:ILE:CD1	1:D:278:LYS:HB3	2.48	0.41
1:D:274:ASN:HA	1:D:277:HIS:CD2	2.56	0.41
1:C:86:SER:O	1:C:88:GLU:N	2.54	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:D:35:TRP:CZ2	1:D:39:ILE:HD11	2.56	0.41
1:C:63:LYS:CB	5:C:386:HOH:O	2.68	0.41
1:D:340:SER:HB3	1:D:343:CYS:HB2	2.01	0.41
1:C:181:ARG:NH1	1:D:159:VAL:O	2.55	0.40
1:B:319:LEU:O	1:B:320:ASN:HB2	2.20	0.40
1:B:100:GLN:HG3	1:B:106:PRO:HG3	2.03	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

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	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
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%)	30	26
1	B	269/304 (88%)	251 (93%)	18 (7%)	23	18
1	C	277/304 (91%)	255 (92%)	22 (8%)	18	13
1	D	268/304 (88%)	252 (94%)	16 (6%)	27	22
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%)	24	20

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
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

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Mol	Chain	Res	Type
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
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

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Mol	Chain	Res	Type
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 chains 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	8,8,8	1.50	2 (25%)	10,11,11	2.15	5 (50%)
2	SNN	F	77	2	8,8,8	1.51	2 (25%)	10,11,11	2.15	5 (50%)

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	-3.51	1.33	1.37
2	E	77	SNN	C2-N1	-3.49	1.33	1.37
2	F	77	SNN	C5-N1	-2.24	1.33	1.37
2	E	77	SNN	C5-N1	-2.20	1.33	1.37

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	77	SNN	C3-C2-N1	4.29	110.75	107.11
2	F	77	SNN	C3-C2-N1	4.28	110.74	107.11
2	E	77	SNN	O2-C2-C3	-2.75	123.82	126.77
2	F	77	SNN	O2-C2-C3	-2.73	123.84	126.77
2	F	77	SNN	O5-C5-C4	-2.64	122.51	126.41
2	E	77	SNN	O5-C5-C4	-2.63	122.53	126.41
2	E	77	SNN	C4-C5-N1	2.29	111.51	108.06
2	F	77	SNN	C4-C5-N1	2.28	111.50	108.06
2	E	77	SNN	C5-N1-C2	-2.06	111.65	113.93
2	F	77	SNN	C5-N1-C2	-2.03	111.68	113.93

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.5 Carbohydrates

There are no carbohydrates in this entry.

5.6 Ligand geometry

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	0.04	0	6,6,6	0.38	0
4	SO4	C	351	-	4,4,4	0.10	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.

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%)	0.98	50 (14%) 3 3	31, 43, 60, 69	0
1	B	334/353 (94%)	0.95	51 (15%) 3 3	33, 44, 57, 71	0
1	C	343/353 (97%)	0.81	34 (9%) 8 8	28, 43, 58, 69	0
1	D	333/353 (94%)	1.03	55 (16%) 2 2	24, 44, 56, 65	0
2	E	7/7 (100%)	1.61	2 (28%) 1 1	62, 63, 66, 67	0
2	F	7/7 (100%)	2.30	4 (57%) 0 1	66, 73, 77, 80	0
2	G	1/7 (14%)	2.71	1 (100%) 0 0	82, 82, 82, 82	0
2	H	1/7 (14%)	1.48	0 100 100	83, 83, 83, 83	0
All	All	1366/1440 (94%)	0.95	197 (14%) 3 3	24, 44, 60, 83	0

All (197) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	89	ASN	7.7
1	D	342	VAL	7.3
1	D	341	PRO	6.8
1	D	347	GLY	6.4
1	B	130	VAL	5.7
1	D	345	VAL	5.6
1	B	210	VAL	5.5
1	D	336	ASN	5.1
1	C	28	TYR	5.0
1	B	299	ALA	4.8
1	D	340	SER	4.6
1	A	195	ASN	4.3
1	B	217	PHE	4.3
1	C	60	ILE	4.3
1	B	302	CYS	4.2
1	B	235	ILE	4.2

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Mol	Chain	Res	Type	RSRZ
1	B	211	VAL	4.2
1	A	235	ILE	4.2
1	A	210	VAL	4.1
1	B	300	ALA	4.1
1	B	303	ALA	4.0
1	A	133	ILE	4.0
1	B	127	GLY	4.0
1	C	62	GLU	3.9
1	A	130	VAL	3.9
1	A	126	ILE	3.9
1	C	16	SER	3.9
1	B	132	VAL	3.9
1	C	217	PHE	3.9
1	B	298	ALA	3.8
1	D	195	ASN	3.8
1	B	344	SER	3.8
1	D	130	VAL	3.7
1	A	196	ASP	3.7
1	C	14	TYR	3.7
1	D	133	ILE	3.7
1	B	342	VAL	3.7
1	B	133	ILE	3.7
1	C	306	VAL	3.7
1	A	13	ARG	3.6
1	C	87	THR	3.6
1	A	299	ALA	3.5
1	D	28	TYR	3.5
1	D	343	CYS	3.5
2	E	73	THR	3.5
1	A	222	TRP	3.5
1	A	215	HIS	3.5
1	D	1	MET	3.5
1	B	237	ALA	3.5
1	A	292	ALA	3.5
1	B	297	VAL	3.5
1	B	120	ILE	3.4
1	A	306	VAL	3.4
1	D	132	VAL	3.4
1	D	194	ILE	3.4
1	B	347	GLY	3.4
2	F	72	ARG	3.4
1	B	119	VAL	3.3

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Mol	Chain	Res	Type	RSRZ
1	A	300	ALA	3.3
1	D	55	LEU	3.3
1	B	131	SER	3.3
1	D	216	PRO	3.3
1	A	211	VAL	3.3
1	A	127	GLY	3.3
1	D	222	TRP	3.3
1	D	346	CYS	3.2
1	B	301	LEU	3.2
1	D	17	GLY	3.2
1	D	300	ALA	3.2
1	B	62	GLU	3.2
1	C	130	VAL	3.2
1	A	218	ASN	3.2
1	B	134	LEU	3.2
1	A	296	ASN	3.1
1	D	82	GLU	3.1
1	A	304	ALA	3.1
1	B	159	VAL	3.1
1	A	193	ASN	3.1
1	B	275	ILE	3.1
1	B	341	PRO	3.1
1	A	347	GLY	3.1
1	B	121	LEU	3.0
1	A	132	VAL	3.0
1	A	134	LEU	3.0
1	A	303	ALA	3.0
1	C	61	PRO	2.9
1	C	132	VAL	2.9
2	F	73	THR	2.9
1	A	297	VAL	2.9
1	C	59	ASN	2.9
1	C	152	ASN	2.9
1	D	259	GLU	2.8
1	C	1	MET	2.8
1	B	296	ASN	2.8
1	C	85	ASN	2.8
1	C	133	ILE	2.8
1	A	120	ILE	2.8
1	D	185	ILE	2.8
1	A	198	THR	2.7
1	B	273	GLU	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	298	ALA	2.7
1	D	261	GLN	2.7
2	E	75	ASN	2.7
2	G	71	MET	2.7
1	D	30	GLU	2.7
1	D	197	TYR	2.7
1	C	210	VAL	2.7
1	D	210	VAL	2.7
1	D	127	GLY	2.7
1	A	263	VAL	2.6
1	B	55	LEU	2.6
1	C	134	LEU	2.6
1	A	5	LEU	2.6
1	B	126	ILE	2.6
1	D	218	ASN	2.6
1	A	129	HIS	2.6
1	A	277	HIS	2.6
1	C	300	ALA	2.6
1	D	147	ASN	2.6
1	D	49	SER	2.5
2	F	71	MET	2.5
1	B	85	ASN	2.5
1	C	272	LYS	2.5
1	B	345	VAL	2.5
1	D	31	ASN	2.5
1	C	271	GLU	2.5
1	B	129	HIS	2.5
1	C	67	ASN	2.5
1	D	90	ASN	2.5
1	C	13	ARG	2.5
1	B	213	ALA	2.5
1	B	304	ALA	2.5
1	B	212	SER	2.5
1	A	17	GLY	2.4
1	C	218	ASN	2.4
1	A	9	VAL	2.4
1	B	118	VAL	2.4
1	D	306	VAL	2.4
1	A	234	TYR	2.4
1	D	153	THR	2.4
1	D	260	CYS	2.4
1	A	342	VAL	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	195	ASN	2.4
1	B	246	PHE	2.4
1	A	293	PRO	2.4
1	A	166	VAL	2.4
1	A	341	PRO	2.4
1	C	155	LEU	2.4
1	D	57	THR	2.4
1	D	191	ALA	2.4
1	A	301	LEU	2.3
1	B	236	ASN	2.3
1	D	3	TYR	2.3
1	C	52	THR	2.3
1	B	149	GLN	2.3
1	C	86	SER	2.3
1	C	270	SER	2.3
1	A	121	LEU	2.3
1	D	85	ASN	2.3
1	B	90	ASN	2.3
1	D	58	ALA	2.2
1	D	59	ASN	2.2
1	A	307	ILE	2.2
1	B	295	ASN	2.2
1	A	302	CYS	2.2
1	D	134	LEU	2.2
1	A	28	TYR	2.2
1	B	144	LEU	2.2
1	D	254	LYS	2.2
2	F	77	SNN	2.2
1	D	275	ILE	2.2
1	D	227	CYS	2.2
1	C	119	VAL	2.2
1	D	255	THR	2.1
1	C	166	VAL	2.1
1	D	29	VAL	2.1
1	D	302	CYS	2.1
1	B	292	ALA	2.1
1	D	159	VAL	2.1
1	A	237	ALA	2.1
1	A	216	PRO	2.1
1	B	293	PRO	2.1
1	D	108	LEU	2.1
1	B	143	ILE	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	29	VAL	2.1
1	D	272	LYS	2.1
1	C	235	ILE	2.1
1	C	222	TRP	2.1
1	A	87	THR	2.0
1	B	208	ILE	2.0
1	B	248	PRO	2.0
1	D	278	LYS	2.0
1	D	64	ASP	2.0
1	A	135	ALA	2.0
1	A	14	TYR	2.0
1	D	164	ASP	2.0
1	A	194	ILE	2.0

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

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	RSR	LLDF	B-factors(Å ²)	Q < 0.9
2	SNN	F	77	8/8	0.33	1.73	63,65,67,68	0
2	SNN	E	77	8/8	0.27	0.59	63,65,65,67	0

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. 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	RSR	LLDF	B-factors(\AA^2)	Q<0.9
4	SO4	B	351	5/5	0.14	-0.89	66,67,68,69	0
3	ZN	A	500	1/1	0.07	-1.49	50,50,50,50	0
3	ZN	C	500	1/1	0.06	-1.76	37,37,37,37	0
3	ZN	B	500	1/1	0.04	-1.87	45,45,45,45	0
3	ZN	D	500	1/1	0.04	-2.55	71,71,71,71	0
4	SO4	C	351	5/5	0.11	-4.29	85,85,85,85	0

6.5 Other polymers ⓘ

There are no such residues in this entry.