



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 08:59 AM GMT

PDB ID : 3GTE  
Title : Crystal Structure of Dicamba Monooxygenase with Non-heme Iron  
Authors : Rydel, T.J.; Sturman, E.J.; Moshiri, F.; Brown, G.R.; Qi, Y.  
Deposited on : 2009-03-27  
Resolution : 1.95 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](mailto: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.

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

MolProbity : 4.02b-467  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
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) : trunk26865

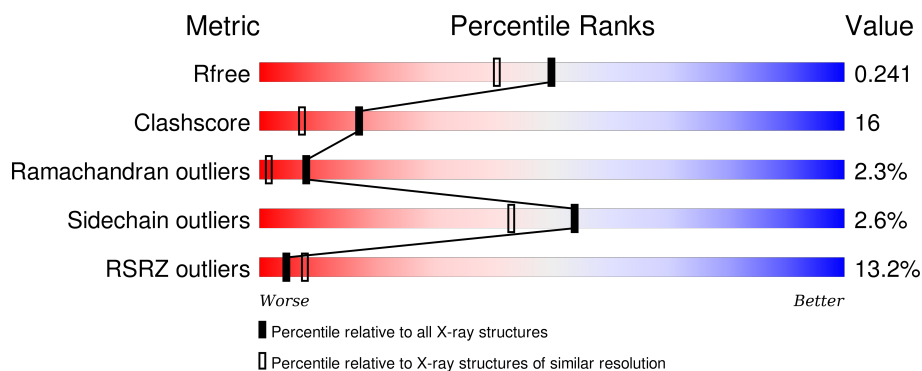
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.95 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))
$R_{free}$	91344	1833 (1.96-1.96)
Clashscore	102246	1953 (1.96-1.96)
Ramachandran outliers	100387	1936 (1.96-1.96)
Sidechain outliers	100360	1936 (1.96-1.96)
RSRZ outliers	91569	1835 (1.96-1.96)

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  $\geq 3$ , 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	349	<div> <div>14%</div> <div>70%</div> <div>25%</div> <div>..</div> </div>
1	B	349	<div> <div>11%</div> <div>73%</div> <div>22%</div> <div>..</div> </div>
1	C	349	<div> <div>13%</div> <div>69%</div> <div>22%</div> <div>.. 5%</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
4	ACT	A	651	-	-	X	-

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8411 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 DdmC.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	341	Total	C	N	O	S	0	0	0
			2641	1665	474	489	13			
1	B	341	Total	C	N	O	S	0	0	0
			2641	1665	474	489	13			
1	C	332	Total	C	N	O	S	0	0	0
			2569	1623	460	473	13			

There are 33 discrepancies between the modelled and reference sequences:

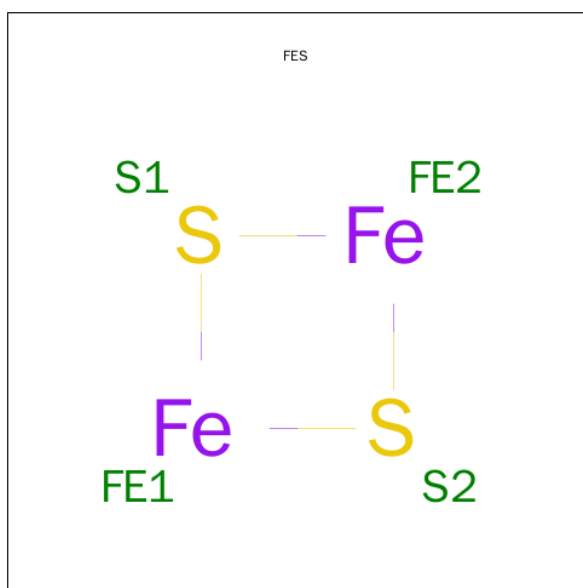
Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	EXPRESSION TAG	UNP Q5S3I3
A	2	ALA	MET	ENGINEERED	UNP Q5S3I3
A	341	ARG	-	EXPRESSION TAG	UNP Q5S3I3
A	342	LEU	-	EXPRESSION TAG	UNP Q5S3I3
A	343	GLU	-	EXPRESSION TAG	UNP Q5S3I3
A	344	HIS	-	EXPRESSION TAG	UNP Q5S3I3
A	345	HIS	-	EXPRESSION TAG	UNP Q5S3I3
A	346	HIS	-	EXPRESSION TAG	UNP Q5S3I3
A	347	HIS	-	EXPRESSION TAG	UNP Q5S3I3
A	348	HIS	-	EXPRESSION TAG	UNP Q5S3I3
A	349	HIS	-	EXPRESSION TAG	UNP Q5S3I3
B	1	MET	-	EXPRESSION TAG	UNP Q5S3I3
B	2	ALA	MET	ENGINEERED	UNP Q5S3I3
B	341	ARG	-	EXPRESSION TAG	UNP Q5S3I3
B	342	LEU	-	EXPRESSION TAG	UNP Q5S3I3
B	343	GLU	-	EXPRESSION TAG	UNP Q5S3I3
B	344	HIS	-	EXPRESSION TAG	UNP Q5S3I3
B	345	HIS	-	EXPRESSION TAG	UNP Q5S3I3
B	346	HIS	-	EXPRESSION TAG	UNP Q5S3I3
B	347	HIS	-	EXPRESSION TAG	UNP Q5S3I3
B	348	HIS	-	EXPRESSION TAG	UNP Q5S3I3
B	349	HIS	-	EXPRESSION TAG	UNP Q5S3I3
C	1	MET	-	EXPRESSION TAG	UNP Q5S3I3

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Chain	Residue	Modelled	Actual	Comment	Reference
C	2	ALA	MET	ENGINEERED	UNP Q5S3I3
C	341	ARG	-	EXPRESSION TAG	UNP Q5S3I3
C	342	LEU	-	EXPRESSION TAG	UNP Q5S3I3
C	343	GLU	-	EXPRESSION TAG	UNP Q5S3I3
C	344	HIS	-	EXPRESSION TAG	UNP Q5S3I3
C	345	HIS	-	EXPRESSION TAG	UNP Q5S3I3
C	346	HIS	-	EXPRESSION TAG	UNP Q5S3I3
C	347	HIS	-	EXPRESSION TAG	UNP Q5S3I3
C	348	HIS	-	EXPRESSION TAG	UNP Q5S3I3
C	349	HIS	-	EXPRESSION TAG	UNP Q5S3I3

- Molecule 2 is FE2/S2 (INORGANIC) CLUSTER (three-letter code: FES) (formula: Fe<sub>2</sub>S<sub>2</sub>).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Fe S 4 2 2	0	0
2	B	1	Total Fe S 4 2 2	0	0
2	C	1	Total Fe S 4 2 2	0	0

- Molecule 3 is FE (III) ION (three-letter code: FE) (formula: Fe).

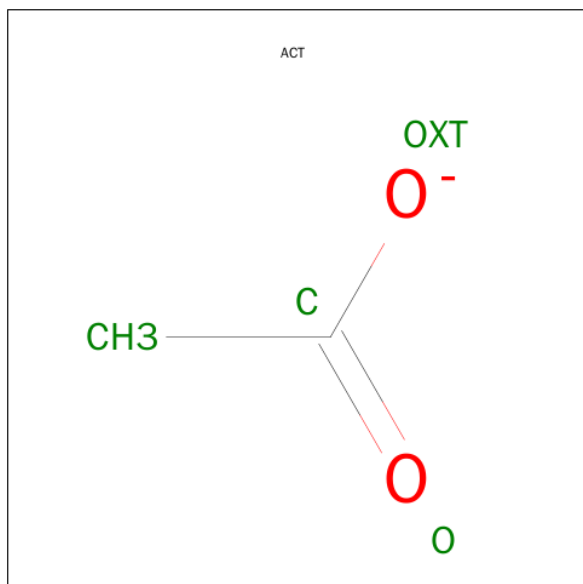
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total Fe 1 1	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Fe	0	0
			1	1		

- Molecule 4 is ACETATE ION (three-letter code: ACT) (formula:  $C_2H_3O_2$ ).



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

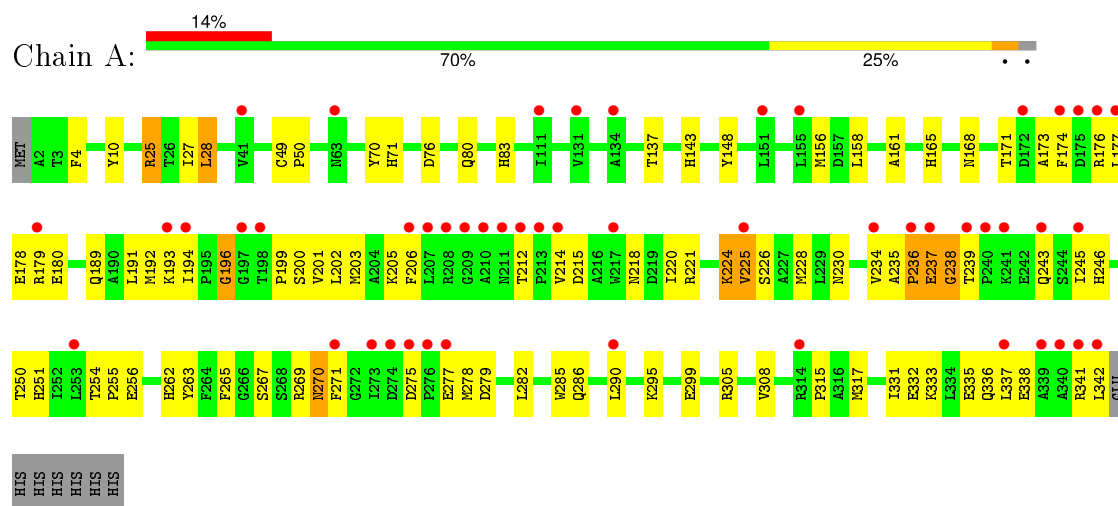
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	165	Total	O	0	0
			165	165		
5	B	177	Total	O	0	0
			177	177		
5	C	188	Total	O	0	0
			188	188		

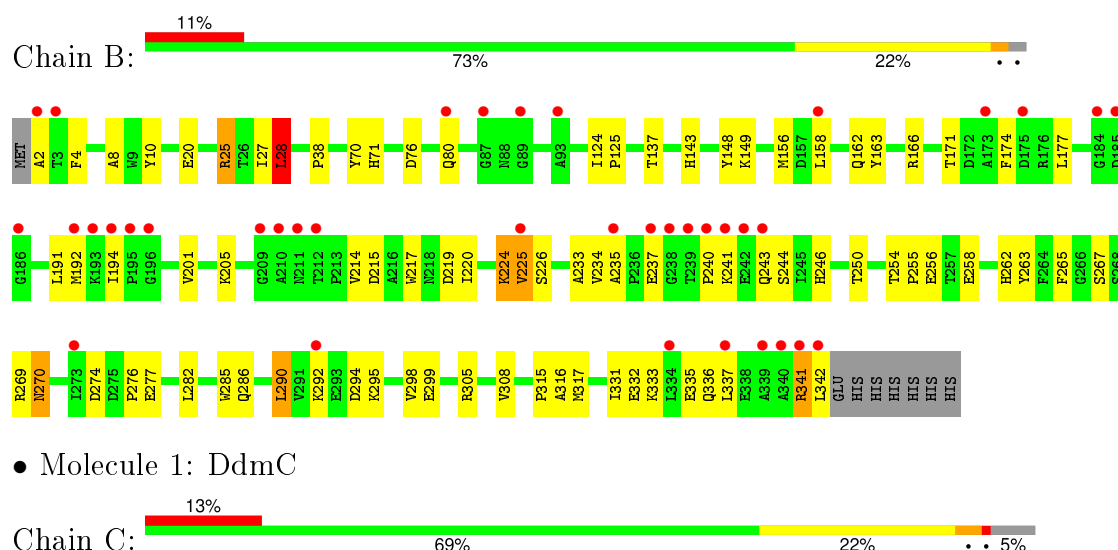
### 3 Residue-property plots [i](#)

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

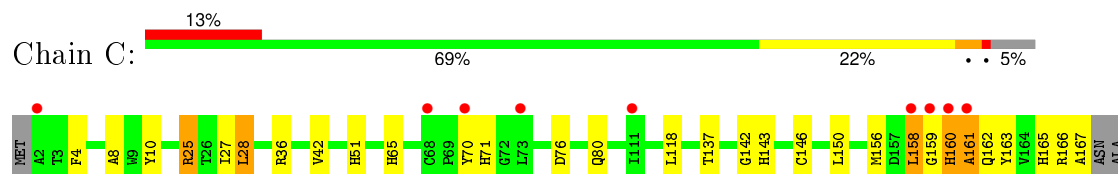
#### • Molecule 1: DdmC

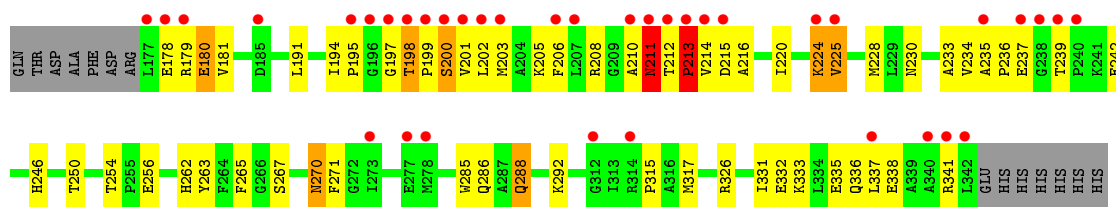


#### • Molecule 1: DdmC



#### • Molecule 1: DdmC







## 4 Data and refinement statistics

Property	Value	Source
Space group	P 32	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	81.11Å 81.11Å 158.81Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 1.95 36.12 – 1.95	Depositor EDS
% Data completeness (in resolution range)	83.6 (20.00-1.95) 84.0 (36.12-1.95)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.64 (at 1.95Å)	Xtriage
Refinement program	CNX2002	Depositor
R, $R_{free}$	0.235 , 0.262 0.216 , 0.241	Depositor DCC
$R_{free}$ test set	7209 reflections (11.28%)	DCC
Wilson B-factor (Å <sup>2</sup> )	30.0	Xtriage
Anisotropy	0.363	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 60.4	EDS
Estimated twinning fraction	0.004 for -h,-k,l 0.039 for h,-h-k,-l 0.024 for -k,-h,-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 71185 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	8411	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.96% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: ACT, FES, FE

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	0.28	0/2706	0.60	0/3685
1	B	0.29	0/2706	0.61	1/3685 (0.0%)
1	C	0.30	0/2632	0.63	0/3583
All	All	0.29	0/8044	0.61	1/10953 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	28	LEU	CA-CB-CG	5.04	126.88	115.30

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	2641	0	2577	88	0
1	B	2641	0	2577	77	0
1	C	2569	0	2515	89	0
2	A	4	0	0	1	0
2	B	4	0	0	1	0
2	C	4	0	0	1	0
3	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	1	0	0	0	0
4	A	8	0	6	5	0
4	B	8	0	6	1	0
5	A	165	0	0	2	0
5	B	177	0	0	6	0
5	C	188	0	0	7	0
All	All	8411	0	7681	254	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 16.

All (254) 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:A:214:VAL:HG11	1:A:234:VAL:HG23	1.47	0.95
1:C:202:LEU:HD23	1:C:234:VAL:HG11	1.54	0.88
1:C:8:ALA:HB2	1:C:224:LYS:HE2	1.55	0.87
1:C:198:THR:HB	1:C:199:PRO:HD3	1.56	0.85
1:A:214:VAL:HG13	1:A:235:ALA:O	1.77	0.84
1:A:178:GLU:HG3	1:A:193:LYS:HE2	1.61	0.82
1:A:236:PRO:HG2	1:A:239:THR:HG21	1.64	0.79
1:C:211:ASN:ND2	1:C:212:THR:H	1.81	0.78
1:C:205:LYS:HZ2	1:C:205:LYS:HB3	1.50	0.76
1:B:294:ASP:O	1:B:298:VAL:HG22	1.86	0.75
1:C:246:HIS:H	1:C:270:ASN:HD21	1.35	0.75
1:C:214:VAL:HG12	1:C:236:PRO:HA	1.68	0.75
1:A:286:GLN:HA	1:A:290:LEU:HD23	1.67	0.74
1:B:174:PHE:HA	1:B:177:LEU:HD13	1.71	0.73
1:A:214:VAL:HG11	1:A:234:VAL:CG2	2.17	0.72
1:C:205:LYS:HB3	1:C:205:LYS:NZ	2.05	0.72
1:B:171:THR:HG21	1:B:194:ILE:HD13	1.72	0.71
1:B:295:LYS:O	1:B:299:GLU:HG3	1.91	0.70
1:A:282:LEU:O	1:A:286:GLN:HG3	1.92	0.69
1:C:212:THR:OG1	1:C:213:PRO:HD2	1.94	0.68
1:C:233:ALA:HB2	1:C:246:HIS:HB3	1.76	0.67
1:A:251:HIS:CE1	4:A:651:ACT:H1	2.29	0.67
1:C:10:TYR:CE2	1:C:225:VAL:HG11	2.30	0.67
1:C:211:ASN:HD22	1:C:211:ASN:N	1.92	0.67
1:A:4:PHE:HB3	1:A:28:LEU:HD13	1.77	0.67
1:B:158:LEU:HD12	1:B:158:LEU:H	1.60	0.66
1:A:285:TRP:HE1	4:A:651:ACT:H3	1.60	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:331:ILE:O	1:A:335:GLU:HG3	1.96	0.66
1:C:4:PHE:HB3	1:C:28:LEU:HD13	1.77	0.66
1:A:245:ILE:HG23	1:A:271:PHE:CD2	2.31	0.66
1:B:331:ILE:O	1:B:335:GLU:HG3	1.97	0.65
1:B:246:HIS:H	1:B:270:ASN:HD21	1.45	0.65
1:B:143:HIS:HE1	1:B:256:GLU:OE1	1.79	0.64
1:A:27:ILE:HD12	1:A:254:THR:HG21	1.79	0.64
1:B:177:LEU:HD23	1:B:192:MET:HE3	1.79	0.64
1:A:285:TRP:HE1	4:A:651:ACT:CH3	2.10	0.64
1:C:142:GLY:HA3	1:C:285:TRP:CZ3	2.32	0.64
1:C:162:GLN:HG3	1:C:288:GLN:HE21	1.63	0.64
1:B:4:PHE:HB3	1:B:28:LEU:HD13	1.80	0.63
1:C:205:LYS:HE2	1:C:271:PHE:HZ	1.62	0.63
1:B:274:ASP:HB3	5:B:1038:HOH:O	1.99	0.63
1:A:214:VAL:HG12	1:A:215:ASP:N	2.13	0.62
1:A:251:HIS:HE1	4:A:651:ACT:H1	1.62	0.62
1:B:214:VAL:HG21	1:B:234:VAL:HG22	1.82	0.62
1:C:143:HIS:HE1	1:C:256:GLU:OE1	1.82	0.62
1:C:224:LYS:HE3	5:C:908:HOH:O	1.99	0.62
1:C:214:VAL:HB	1:C:235:ALA:O	2.00	0.62
1:C:165:HIS:CE1	1:C:205:LYS:HZ1	2.17	0.62
1:A:290:LEU:HD22	1:A:290:LEU:N	2.15	0.62
1:B:282:LEU:O	1:B:286:GLN:HG3	2.00	0.62
1:A:212:THR:O	1:A:214:VAL:HG23	1.99	0.61
1:A:342:LEU:HD23	1:A:342:LEU:H	1.64	0.61
1:C:331:ILE:O	1:C:335:GLU:HG3	2.00	0.61
1:C:246:HIS:H	1:C:270:ASN:ND2	1.98	0.60
1:C:163:TYR:O	1:C:166:ARG:HG3	2.01	0.60
1:A:338:GLU:O	1:A:342:LEU:HD23	2.01	0.60
1:C:162:GLN:HG3	1:C:288:GLN:NE2	2.17	0.60
1:B:286:GLN:HA	1:B:290:LEU:HD22	1.83	0.60
1:B:162:GLN:HE22	1:B:171:THR:H	1.50	0.59
1:A:214:VAL:CG1	1:A:215:ASP:N	2.65	0.59
1:A:246:HIS:H	1:A:270:ASN:HD21	1.49	0.59
1:B:177:LEU:CD1	1:B:194:ILE:HG12	2.33	0.58
1:B:235:ALA:HB2	1:B:244:SER:OG	2.04	0.58
1:A:83:HIS:HD2	5:A:827:HOH:O	1.87	0.58
1:C:205:LYS:HZ2	1:C:205:LYS:CB	2.15	0.57
1:A:143:HIS:HE1	1:A:256:GLU:OE1	1.87	0.57
1:C:211:ASN:ND2	1:C:211:ASN:N	2.53	0.57
1:A:178:GLU:HG3	1:A:193:LYS:HB2	1.87	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:177:LEU:HG	1:B:192:MET:SD	2.44	0.57
1:A:295:LYS:O	1:A:299:GLU:HG3	2.04	0.57
1:B:246:HIS:H	1:B:270:ASN:ND2	2.03	0.57
1:C:210:ALA:HB3	5:C:1009:HOH:O	2.05	0.57
1:B:71:HIS:HB2	2:B:501:FES:S1	2.45	0.56
1:B:240:PRO:HG2	1:B:243:GLN:HG2	1.88	0.56
1:A:315:PRO:HB2	1:A:317:MET:CE	2.36	0.56
1:B:174:PHE:O	1:B:177:LEU:HD13	2.05	0.56
1:B:174:PHE:HA	1:B:177:LEU:CD1	2.35	0.56
1:C:195:PRO:HA	1:C:215:ASP:OD1	2.06	0.56
1:C:8:ALA:CB	1:C:224:LYS:HE2	2.32	0.56
1:C:167:ALA:O	1:C:200:SER:HB3	2.06	0.56
1:A:199:PRO:HB2	1:A:203:MET:HB2	1.88	0.55
1:C:27:ILE:HD12	1:C:254:THR:HG21	1.88	0.55
1:C:10:TYR:CZ	1:C:225:VAL:HG11	2.41	0.55
1:B:158:LEU:N	1:B:158:LEU:HD12	2.20	0.55
1:A:71:HIS:HB2	2:A:501:FES:S1	2.47	0.55
1:A:214:VAL:CG1	1:A:234:VAL:HG23	2.31	0.55
1:B:25:ARG:HD2	1:B:262:HIS:CE1	2.41	0.55
1:C:167:ALA:HB2	1:C:203:MET:HG2	1.89	0.55
1:C:181:VAL:HG11	1:C:326:ARG:HD2	1.89	0.54
1:B:177:LEU:HD11	1:B:194:ILE:HG12	1.87	0.54
1:C:214:VAL:HG12	1:C:236:PRO:CA	2.37	0.54
1:A:25:ARG:HD2	1:A:262:HIS:CE1	2.42	0.54
1:B:156:MET:HA	1:B:156:MET:HE2	1.90	0.54
1:A:10:TYR:CE2	1:A:225:VAL:HG11	2.43	0.54
1:C:233:ALA:CB	1:C:246:HIS:HB3	2.38	0.54
1:B:217:TRP:CE2	1:B:241:LYS:HB3	2.43	0.53
1:A:173:ALA:HB3	1:A:194:ILE:CG2	2.39	0.53
1:A:290:LEU:HD22	1:A:290:LEU:H	1.73	0.53
1:C:71:HIS:HB2	2:C:501:FES:S1	2.48	0.53
1:B:156:MET:HE2	1:B:220:ILE:HG21	1.91	0.53
1:C:315:PRO:HB2	1:C:317:MET:HE1	1.91	0.53
1:A:315:PRO:HB2	1:A:317:MET:HE1	1.90	0.52
1:C:332:GLU:O	1:C:336:GLN:HG3	2.09	0.52
1:B:269:ARG:HD3	1:B:282:LEU:HD11	1.92	0.52
1:C:265:PHE:CE2	1:C:285:TRP:HE3	2.28	0.52
1:B:158:LEU:HD13	5:B:1000:HOH:O	2.09	0.52
1:B:276:PRO:HG2	1:B:277:GLU:OE2	2.09	0.52
1:B:285:TRP:NE1	4:B:653:ACT:H1	2.24	0.52
1:C:143:HIS:HD2	5:C:1115:HOH:O	1.93	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:315:PRO:HB2	1:C:317:MET:CE	2.39	0.52
1:A:173:ALA:HB3	1:A:194:ILE:HG22	1.92	0.52
1:C:118:LEU:HD11	1:C:341:ARG:HH12	1.75	0.52
1:C:206:PHE:CE2	1:C:246:HIS:HA	2.46	0.51
1:B:333:LYS:O	1:B:337:LEU:HG	2.11	0.51
1:B:10:TYR:CE2	1:B:225:VAL:HG11	2.44	0.51
1:B:27:ILE:HD12	1:B:254:THR:HG21	1.91	0.51
1:B:38:PRO:HD2	5:B:1113:HOH:O	2.09	0.51
1:C:206:PHE:HE2	1:C:246:HIS:HA	1.75	0.51
1:A:189:GLN:HG2	1:A:221:ARG:HA	1.93	0.51
1:A:286:GLN:HA	1:A:290:LEU:CD2	2.40	0.50
1:C:65:HIS:HA	5:C:842:HOH:O	2.10	0.50
1:B:2:ALA:HB1	1:B:258:GLU:OE1	2.10	0.50
1:C:156:MET:HE2	1:C:220:ILE:HD13	1.92	0.50
1:B:315:PRO:HB2	1:B:317:MET:HE1	1.94	0.50
1:C:160:HIS:O	1:C:161:ALA:O	2.30	0.50
1:C:236:PRO:O	1:C:239:THR:HG23	2.12	0.50
1:A:171:THR:HG23	1:A:174:PHE:HB2	1.93	0.50
1:B:20:GLU:HG3	5:B:1181:HOH:O	2.12	0.49
1:A:156:MET:HA	1:A:156:MET:HE2	1.93	0.49
1:B:295:LYS:HE2	5:B:1146:HOH:O	2.11	0.49
1:C:205:LYS:HZ3	1:C:206:PHE:HE1	1.60	0.49
1:B:156:MET:HE1	1:B:220:ILE:HG12	1.95	0.49
1:C:198:THR:HB	1:C:199:PRO:CD	2.36	0.49
1:C:286:GLN:NE2	1:C:292:LYS:H	2.11	0.49
1:C:162:GLN:HG3	1:C:288:GLN:O	2.13	0.49
1:B:8:ALA:HB2	1:B:224:LYS:HE2	1.95	0.49
1:C:42:VAL:HG23	5:C:1013:HOH:O	2.13	0.49
1:B:201:VAL:O	1:B:205:LYS:HG3	2.13	0.49
1:A:191:LEU:HD22	1:A:191:LEU:N	2.28	0.49
1:C:270:ASN:C	1:C:270:ASN:HD22	2.15	0.48
1:C:333:LYS:O	1:C:337:LEU:HG	2.13	0.48
1:A:275:ASP:HB3	1:A:278:MET:HB3	1.95	0.48
1:A:156:MET:HE3	1:A:228:MET:SD	2.53	0.48
1:B:315:PRO:HB2	1:B:317:MET:CE	2.42	0.48
1:B:292:LYS:O	1:B:292:LYS:HG3	2.13	0.48
1:B:156:MET:CE	1:B:156:MET:HA	2.43	0.48
1:B:174:PHE:CA	1:B:177:LEU:HD13	2.40	0.47
1:A:270:ASN:HD22	1:A:270:ASN:C	2.18	0.47
1:C:201:VAL:HG23	1:C:202:LEU:N	2.29	0.47
1:C:211:ASN:ND2	1:C:212:THR:N	2.56	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:224:LYS:HG2	1:C:225:VAL:HG23	1.97	0.47
1:A:171:THR:HG21	1:A:194:ILE:HD13	1.97	0.47
1:A:137:THR:HA	1:A:267:SER:O	2.15	0.47
1:B:240:PRO:HG2	1:B:243:GLN:CG	2.44	0.47
1:B:149:LYS:NZ	1:B:335:GLU:OE2	2.37	0.47
1:A:237:GLU:HB3	1:A:238:GLY:H	1.49	0.47
1:A:269:ARG:NH2	1:A:279:ASP:OD2	2.42	0.47
1:B:177:LEU:HD23	1:B:192:MET:CE	2.43	0.47
1:A:224:LYS:HG2	1:A:225:VAL:HG23	1.96	0.47
1:A:332:GLU:O	1:A:336:GLN:HG3	2.15	0.47
1:A:178:GLU:CG	1:A:193:LYS:HB2	2.44	0.47
1:A:180:GLU:HB2	1:A:191:LEU:HB2	1.97	0.47
1:A:237:GLU:O	1:A:239:THR:N	2.47	0.47
1:B:282:LEU:HD13	5:B:899:HOH:O	2.14	0.47
1:B:76:ASP:OD2	1:B:80:GLN:HB3	2.14	0.47
1:A:177:LEU:HD21	1:A:179:ARG:HH21	1.80	0.47
1:A:246:HIS:H	1:A:270:ASN:ND2	2.11	0.46
1:A:275:ASP:OD1	1:A:277:GLU:HB2	2.15	0.46
1:C:242:GLU:CD	1:C:242:GLU:H	2.18	0.46
1:B:224:LYS:HG2	1:B:225:VAL:HG23	1.97	0.46
1:C:194:ILE:HB	1:C:216:ALA:HB3	1.97	0.46
1:C:179:ARG:O	1:C:180:GLU:HB2	2.16	0.46
1:A:333:LYS:O	1:A:337:LEU:HG	2.15	0.46
1:A:202:LEU:HG	1:A:203:MET:HE2	1.97	0.46
1:B:214:VAL:HG23	1:B:235:ALA:C	2.36	0.45
1:A:148:TYR:CG	1:A:255:PRO:HG3	2.52	0.45
1:C:199:PRO:O	1:C:200:SER:HB2	2.16	0.45
1:B:233:ALA:HB2	1:B:246:HIS:HB3	1.98	0.45
1:A:156:MET:HA	1:A:156:MET:CE	2.46	0.45
1:A:224:LYS:O	1:A:225:VAL:HB	2.17	0.45
1:C:201:VAL:HG23	1:C:202:LEU:H	1.81	0.45
1:A:156:MET:HE2	1:A:220:ILE:HG21	1.98	0.45
1:C:76:ASP:OD2	1:C:80:GLN:HB3	2.17	0.45
1:A:200:SER:HB2	5:A:916:HOH:O	2.17	0.45
1:C:212:THR:O	1:C:214:VAL:HG13	2.17	0.45
1:C:233:ALA:HB2	1:C:246:HIS:CB	2.44	0.45
1:B:224:LYS:O	1:B:225:VAL:HB	2.16	0.45
1:A:168:ASN:ND2	1:A:202:LEU:HB2	2.32	0.45
1:A:161:ALA:HA	1:A:165:HIS:CD2	2.52	0.45
1:A:250:THR:O	1:A:265:PHE:HA	2.17	0.45
1:A:270:ASN:ND2	1:A:270:ASN:C	2.70	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:220:ILE:HA	1:C:230:ASN:HA	1.99	0.45
1:C:167:ALA:HB2	1:C:203:MET:CG	2.48	0.44
1:C:250:THR:O	1:C:265:PHE:HA	2.17	0.44
1:A:305:ARG:O	1:A:308:VAL:HG22	2.18	0.44
1:A:76:ASP:OD2	1:A:80:GLN:HB3	2.18	0.44
1:C:211:ASN:HD22	1:C:212:THR:H	1.63	0.44
1:C:270:ASN:C	1:C:270:ASN:ND2	2.70	0.44
1:A:275:ASP:HB3	1:A:278:MET:CB	2.48	0.44
1:C:180:GLU:HB3	1:C:191:LEU:HB2	2.00	0.44
1:A:158:LEU:H	1:A:158:LEU:HD22	1.83	0.44
1:C:211:ASN:HD22	1:C:211:ASN:H	1.65	0.44
1:B:137:THR:HA	1:B:267:SER:O	2.18	0.44
1:B:332:GLU:O	1:B:336:GLN:HG3	2.18	0.44
1:C:137:THR:HA	1:C:267:SER:O	2.17	0.43
1:B:163:TYR:O	1:B:166:ARG:HB2	2.18	0.43
1:C:25:ARG:HD2	1:C:262:HIS:CE1	2.52	0.43
1:B:158:LEU:CD1	1:B:158:LEU:H	2.28	0.43
1:B:158:LEU:HD11	1:B:220:ILE:HG21	2.01	0.43
1:B:10:TYR:CZ	1:B:225:VAL:HG11	2.53	0.43
1:A:176:ARG:NH2	1:A:196:GLY:O	2.52	0.43
1:A:192:MET:HB2	1:A:218:ASN:O	2.17	0.43
1:A:235:ALA:HB1	1:A:236:PRO:HD2	1.99	0.43
1:B:225:VAL:HG12	1:B:226:SER:N	2.33	0.43
1:C:156:MET:CE	1:C:228:MET:SD	3.06	0.43
1:A:269:ARG:HH11	1:A:282:LEU:HD12	1.83	0.43
1:C:156:MET:CE	1:C:220:ILE:HD13	2.49	0.43
1:A:201:VAL:O	1:A:205:LYS:HG3	2.18	0.43
1:B:191:LEU:HD23	1:B:219:ASP:HB3	2.00	0.43
4:A:650:ACT:H3	4:A:651:ACT:H2	2.00	0.43
1:A:225:VAL:HG12	1:A:226:SER:N	2.34	0.43
1:A:168:ASN:ND2	1:A:285:TRP:CE3	2.86	0.42
1:B:214:VAL:HG22	1:B:215:ASP:N	2.34	0.42
1:A:220:ILE:HA	1:A:230:ASN:HA	2.00	0.42
1:B:250:THR:O	1:B:265:PHE:HA	2.20	0.42
1:B:305:ARG:O	1:B:308:VAL:HG22	2.19	0.42
1:B:166:ARG:HB2	1:B:166:ARG:HE	1.64	0.42
1:B:316:ALA:O	1:C:51:HIS:HD2	2.02	0.42
1:A:236:PRO:O	1:A:237:GLU:C	2.57	0.42
1:C:156:MET:CE	1:C:156:MET:HA	2.50	0.42
1:C:178:GLU:OE1	1:C:179:ARG:NE	2.53	0.42
1:B:337:LEU:O	1:B:341:ARG:N	2.47	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:267:SER:OG	1:A:282:LEU:HD13	2.20	0.41
1:A:203:MET:O	1:A:206:PHE:HB2	2.20	0.41
1:A:156:MET:HE1	1:A:220:ILE:HG12	2.02	0.41
1:A:341:ARG:HA	1:A:341:ARG:HD2	1.71	0.41
1:B:217:TRP:CZ2	1:B:241:LYS:HB3	2.56	0.41
1:A:305:ARG:HA	1:A:308:VAL:HG22	2.02	0.41
1:C:195:PRO:O	1:C:197:GLY:N	2.43	0.41
1:B:156:MET:CE	1:B:220:ILE:HD13	2.50	0.41
1:C:338:GLU:HG3	5:C:1100:HOH:O	2.21	0.41
1:C:224:LYS:O	1:C:225:VAL:HB	2.19	0.41
1:C:36:ARG:NE	5:C:1013:HOH:O	2.27	0.41
1:A:214:VAL:HG13	1:A:235:ALA:C	2.39	0.41
1:B:270:ASN:C	1:B:270:ASN:HD22	2.23	0.41
1:C:146:CYS:HB2	1:C:150:LEU:HD12	2.02	0.41
1:B:124:ILE:HA	1:B:125:PRO:HD3	1.96	0.41
1:C:28:LEU:N	1:C:28:LEU:HD23	2.37	0.40
1:B:234:VAL:O	1:B:244:SER:OG	2.39	0.40
1:B:148:TYR:CG	1:B:255:PRO:HG3	2.56	0.40
1:A:49:CYS:HA	1:A:50:PRO:HD3	1.90	0.40
1:A:27:ILE:HD12	1:A:254:THR:CG2	2.49	0.40
1:C:10:TYR:OH	1:C:225:VAL:HG11	2.22	0.40
1:C:158:LEU:C	1:C:158:LEU:HD13	2.42	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	339/349 (97%)	313 (92%)	19 (6%)	7 (2%)	9	2
1	B	339/349 (97%)	317 (94%)	18 (5%)	4 (1%)	16	5
1	C	328/349 (94%)	296 (90%)	20 (6%)	12 (4%)	4	0

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	1006/1047 (96%)	926 (92%)	57 (6%)	23 (2%)	8 1

All (23) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	237	GLU
1	C	161	ALA
1	A	196	GLY
1	A	225	VAL
1	A	237	GLU
1	A	238	GLY
1	B	225	VAL
1	B	341	ARG
1	C	158	LEU
1	C	200	SER
1	C	225	VAL
1	A	224	LYS
1	B	224	LYS
1	C	180	GLU
1	C	198	THR
1	C	208	ARG
1	C	211	ASN
1	C	224	LYS
1	A	243	GLN
1	C	237	GLU
1	C	213	PRO
1	A	236	PRO
1	C	159	GLY

### 5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	276/284 (97%)	271 (98%)	5 (2%)	66 60
1	B	276/284 (97%)	269 (98%)	7 (2%)	55 45

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	269/284 (95%)	260 (97%)	9 (3%)	45	32
All	All	821/852 (96%)	800 (97%)	21 (3%)	54	43

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	ARG
1	A	28	LEU
1	A	70	TYR
1	A	263	TYR
1	A	270	ASN
1	B	25	ARG
1	B	28	LEU
1	B	70	TYR
1	B	263	TYR
1	B	270	ASN
1	B	290	LEU
1	B	342	LEU
1	C	25	ARG
1	C	28	LEU
1	C	70	TYR
1	C	160	HIS
1	C	211	ASN
1	C	213	PRO
1	C	263	TYR
1	C	270	ASN
1	C	288	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	63	ASN
1	A	83	HIS
1	A	143	HIS
1	A	168	ASN
1	A	211	ASN
1	A	270	ASN
1	A	288	GLN
1	B	63	ASN
1	B	143	HIS
1	B	162	GLN

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Mol	Chain	Res	Type
1	B	223	ASN
1	B	243	GLN
1	B	270	ASN
1	C	63	ASN
1	C	143	HIS
1	C	211	ASN
1	C	223	ASN
1	C	270	ASN
1	C	286	GLN
1	C	288	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

Of 9 ligands modelled in this entry, 2 are monoatomic - leaving 7 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
2	FES	A	501	1	0,4,4	0.00	-	0,4,4	0.00	-
4	ACT	A	650	3	1,3,3	0.43	0	0,3,3	0.00	-
4	ACT	A	651	-	1,3,3	1.59	0	0,3,3	0.00	-
2	FES	B	501	1	0,4,4	0.00	-	0,4,4	0.00	-

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
4	ACT	B	652	3	1,3,3	0.82	0	0,3,3	0.00	-
4	ACT	B	653	-	1,3,3	0.70	0	0,3,3	0.00	-
2	FES	C	501	1	0,4,4	0.00	-	0,4,4	0.00	-

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	FES	A	501	1	-	0/0/4/4	0/1/1/1
4	ACT	A	650	3	-	0/0/0/0	0/0/0/0
4	ACT	A	651	-	-	0/0/0/0	0/0/0/0
2	FES	B	501	1	-	0/0/4/4	0/1/1/1
4	ACT	B	652	3	-	0/0/0/0	0/0/0/0
4	ACT	B	653	-	-	0/0/0/0	0/0/0/0
2	FES	C	501	1	-	0/0/4/4	0/1/1/1

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.

6 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	FES	1	0
4	A	650	ACT	1	0
4	A	651	ACT	5	0
2	B	501	FES	1	0
4	B	653	ACT	1	0
2	C	501	FES	1	0

## 5.7 Other polymers ⓘ

There are no such residues in this entry.

## 5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

## 6 Fit of model and data

### 6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	341/349 (97%)	0.85	50 (14%) <b>3</b> <b>5</b>	22, 39, 73, 80	0
1	B	341/349 (97%)	0.66	38 (11%) <b>7</b> <b>11</b>	24, 37, 63, 85	0
1	C	332/349 (95%)	0.94	46 (13%) <b>4</b> <b>6</b>	23, 35, 73, 82	0
All	All	1014/1047 (96%)	0.81	134 (13%) <b>4</b> <b>8</b>	22, 37, 70, 85	0

All (134) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	177	LEU	15.3
1	C	200	SER	11.0
1	B	339	ALA	10.2
1	A	211	ASN	9.3
1	A	342	LEU	9.0
1	C	212	THR	8.3
1	C	211	ASN	8.1
1	C	207	LEU	8.0
1	C	197	GLY	7.8
1	C	210	ALA	6.8
1	C	199	PRO	6.8
1	B	340	ALA	6.4
1	A	340	ALA	6.4
1	B	210	ALA	6.2
1	C	198	THR	6.1
1	A	209	GLY	5.8
1	B	211	ASN	5.5
1	C	237	GLU	5.5
1	A	208	ARG	5.5
1	A	239	THR	5.1
1	C	340	ALA	5.1
1	C	203	MET	4.7
1	A	276	PRO	4.6

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Mol	Chain	Res	Type	RSRZ
1	C	337	LEU	4.6
1	A	274	ASP	4.6
1	A	213	PRO	4.6
1	A	197	GLY	4.4
1	C	158	LEU	4.4
1	B	192	MET	4.4
1	C	159	GLY	4.3
1	A	210	ALA	4.3
1	A	337	LEU	4.3
1	A	243	GLN	4.3
1	B	239	THR	4.1
1	B	212	THR	4.1
1	C	202	LEU	4.1
1	A	225	VAL	4.1
1	C	312	GLY	4.0
1	C	235	ALA	4.0
1	C	201	VAL	3.9
1	B	237	GLU	3.9
1	A	212	THR	3.9
1	B	225	VAL	3.8
1	A	273	ILE	3.8
1	B	3	THR	3.8
1	A	339	ALA	3.8
1	C	206	PHE	3.7
1	B	93	ALA	3.7
1	A	194	ILE	3.7
1	A	214	VAL	3.6
1	A	134	ALA	3.6
1	C	273	ILE	3.5
1	C	342	LEU	3.5
1	A	131	VAL	3.5
1	B	238	GLY	3.4
1	C	213	PRO	3.4
1	A	236	PRO	3.4
1	C	341	ARG	3.3
1	C	160	HIS	3.3
1	A	172	ASP	3.3
1	B	2	ALA	3.2
1	C	178	GLU	3.2
1	A	175	ASP	3.2
1	A	174	PHE	3.2
1	C	2	ALA	3.2

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Mol	Chain	Res	Type	RSRZ
1	C	239	THR	3.2
1	B	195	PRO	3.2
1	C	70	TYR	3.1
1	A	206	PHE	3.1
1	C	215	ASP	3.1
1	A	193	LYS	3.0
1	B	242	GLU	3.0
1	B	341	ARG	3.0
1	A	240	PRO	3.0
1	B	185	ASP	3.0
1	B	158	LEU	2.9
1	B	209	GLY	2.9
1	A	198	THR	2.9
1	A	271	PHE	2.9
1	A	151	LEU	2.9
1	A	207	LEU	2.9
1	A	217	TRP	2.8
1	B	184	GLY	2.8
1	B	175	ASP	2.8
1	C	224	LYS	2.7
1	A	245	ILE	2.7
1	C	225	VAL	2.7
1	B	240	PRO	2.6
1	B	186	GLY	2.6
1	B	89	GLY	2.6
1	C	277	GLU	2.6
1	C	68	CYS	2.5
1	B	292	LYS	2.5
1	A	176	ARG	2.5
1	A	241	LYS	2.5
1	B	196	GLY	2.5
1	C	179	ARG	2.5
1	C	314	ARG	2.5
1	B	235	ALA	2.5
1	C	185	ASP	2.5
1	A	177	LEU	2.4
1	B	342	LEU	2.4
1	A	341	ARG	2.4
1	B	243	GLN	2.4
1	B	193	LYS	2.4
1	B	273	ILE	2.4
1	A	155	LEU	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	290	LEU	2.3
1	A	234	VAL	2.3
1	A	179	ARG	2.3
1	A	314	ARG	2.3
1	C	238	GLY	2.3
1	C	161	ALA	2.3
1	B	337	LEU	2.2
1	C	214	VAL	2.2
1	B	173	ALA	2.2
1	C	278	MET	2.2
1	A	111	ILE	2.2
1	B	80	GLN	2.2
1	C	111	ILE	2.2
1	A	237	GLU	2.1
1	A	277	GLU	2.1
1	B	334	LEU	2.1
1	A	41	VAL	2.1
1	C	240	PRO	2.1
1	A	253	LEU	2.1
1	B	241	LYS	2.1
1	C	196	GLY	2.1
1	C	195	PRO	2.1
1	A	275	ASP	2.1
1	C	73	LEU	2.0
1	A	63	ASN	2.0
1	B	87	GLY	2.0
1	B	194	ILE	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
4	ACT	A	651	4/4	0.95	0.14	0.25	39,42,42,43	0
2	FES	A	501	4/4	0.99	0.11	-0.30	24,25,26,26	0
4	ACT	A	650	4/4	0.93	0.13	-0.52	37,38,38,40	0
2	FES	B	501	4/4	0.99	0.09	-0.67	30,31,31,31	0
4	ACT	B	652	4/4	0.96	0.09	-0.81	34,35,35,36	0
3	FE	A	502	1/1	0.99	0.13	-0.82	39,39,39,39	0
2	FES	C	501	4/4	0.99	0.15	-0.84	26,27,27,29	0
3	FE	B	502	1/1	0.99	0.08	-1.06	31,31,31,31	0
4	ACT	B	653	4/4	0.98	0.08	-2.28	35,37,37,38	0

## 6.5 Other polymers

There are no such residues in this entry.