



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 4, 2026 – 06:04 PM EDT

PDB ID : 3GTE / pdb_00003gte
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

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.50

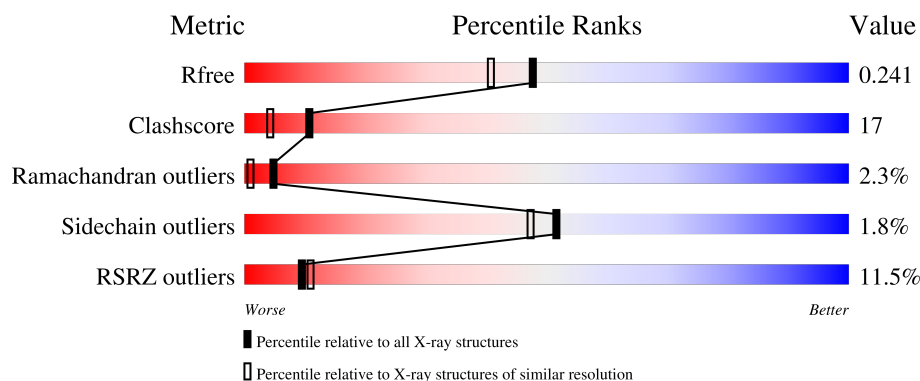
1 Overall quality at a glance

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}	180053	3494 (1.96-1.96)
Clashscore	190562	3612 (1.96-1.96)
Ramachandran outliers	187476	3587 (1.96-1.96)
Sidechain outliers	187428	3587 (1.96-1.96)
RSRZ outliers	180081	3495 (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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	11% 68% 25% 5%
1	B	349	8% 72% 23% • •
1	C	349	14% 68% 22% • • 5%

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:

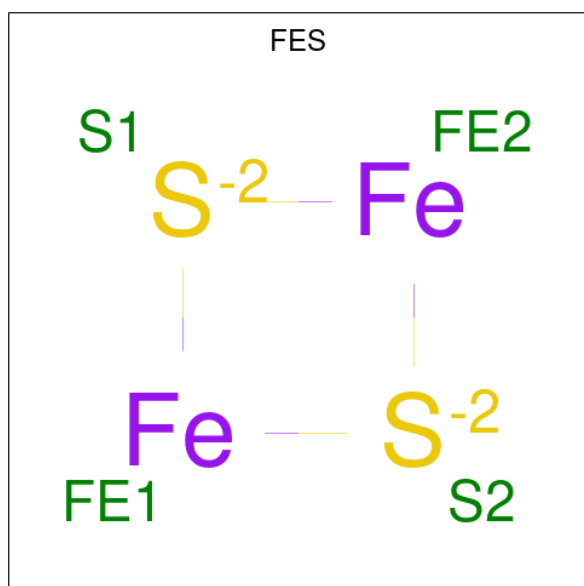
Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	expression tag	UNP Q5S3I3
A	2	ALA	MET	engineered mutation	UNP Q5S3I3
A	341	ARG	-	expression tag	UNP Q5S3I3
A	342	LEU	-	expression tag	UNP Q5S3I3
A	343	GLU	-	expression tag	UNP Q5S3I3
A	344	HIS	-	expression tag	UNP Q5S3I3
A	345	HIS	-	expression 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 mutation	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 mutation	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 (CCD ID: FES) (formula: Fe_2S_2).



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 (CCD ID: FE) (formula: Fe).

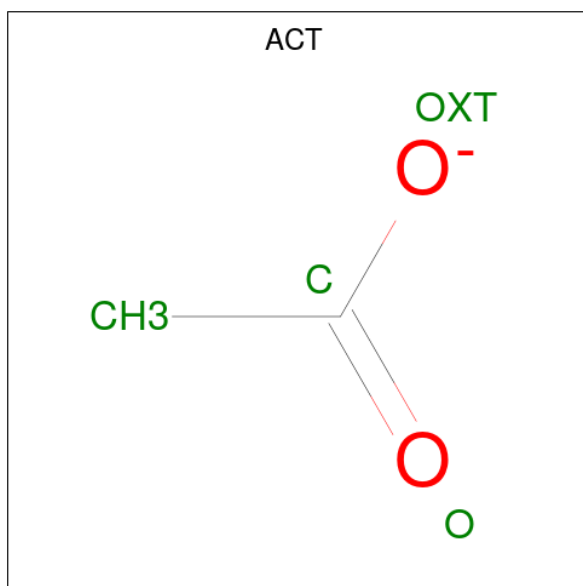
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Fe 1 1	0	0

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

- Molecule 4 is ACETATE ION (CCD ID: 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		

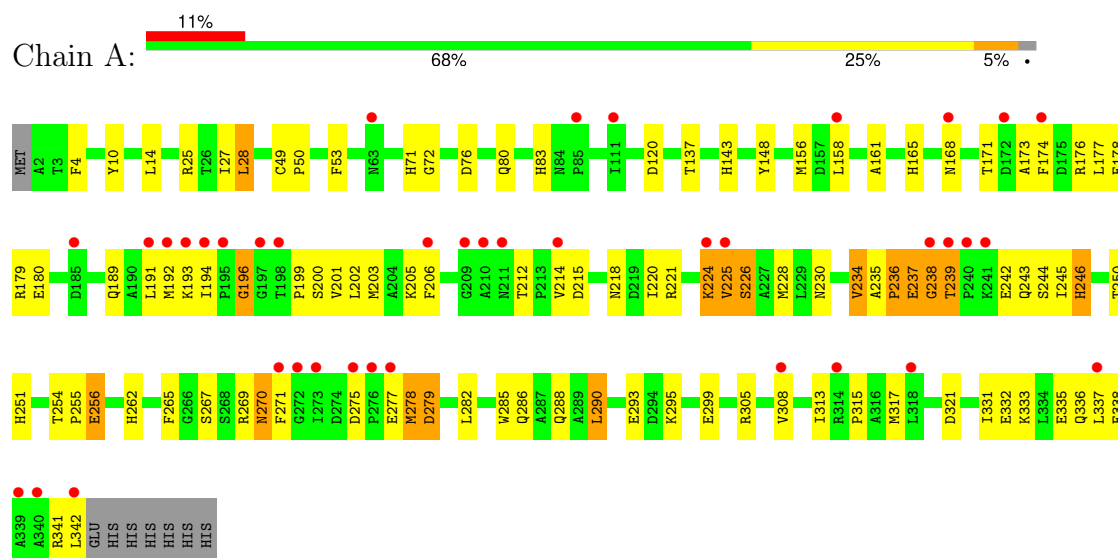
- Molecule 5 is water.

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		

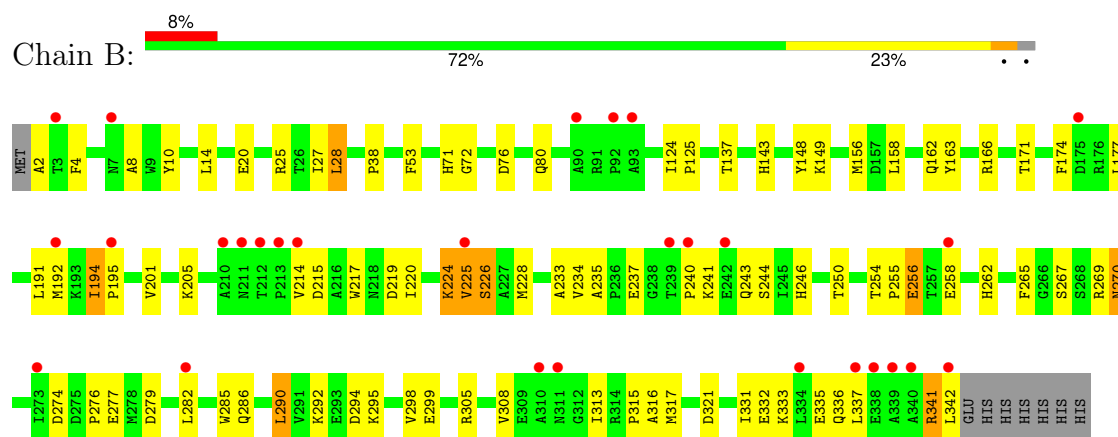
3 Residue-property plots

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

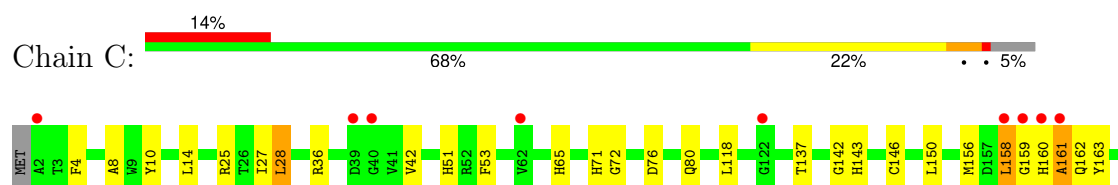
• Molecule 1: DdmC

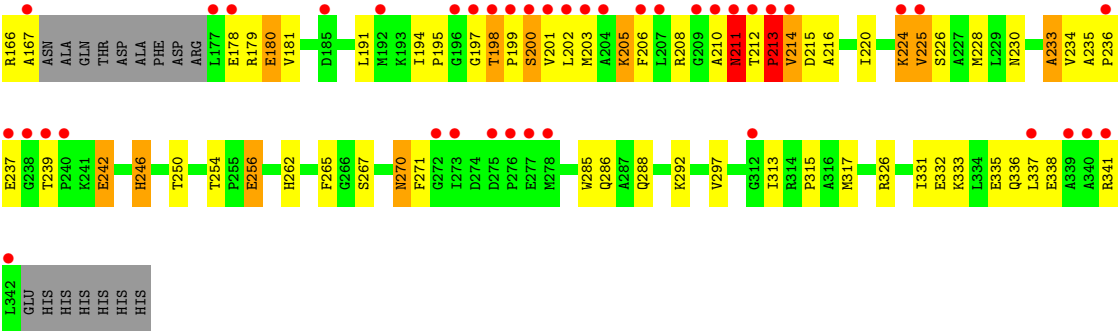


• Molecule 1: DdmC



• Molecule 1: DdmC





4 Data and refinement statistics

Property	Value	Source
Space group	P 32	Depositor
Cell constants a, b, c, α , β , γ	81.11Å 81.11Å 158.81Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 1.95 20.00 – 1.95	Depositor EDS
% Data completeness (in resolution range)	83.6 (20.00-1.95) 84.0 (20.00-1.95)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.64 (at 1.95Å)	Xtriage
Refinement program	CNX	Depositor
R, R_{free}	0.235 , 0.262 0.216 , 0.241	Depositor DCC
R_{free} test set	7209 reflections (10.13%)	wwPDB-VP
Wilson B-factor (Å ²)	30.0	Xtriage
Anisotropy	0.363	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 51.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.004 for -h,-k,l 0.039 for h,-h-k,-l 0.024 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8411	wwPDB-VP
Average B, all atoms (Å ²)	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.*

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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: FES, FE, ACT

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	0.35	0/2706	0.91	12/3685 (0.3%)
1	B	0.37	0/2706	0.93	11/3685 (0.3%)
1	C	0.40	0/2632	0.93	10/3583 (0.3%)
All	All	0.37	0/8044	0.92	33/10953 (0.3%)

There are no bond length outliers.

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	226	SER	N-CA-C	7.21	118.92	110.41
1	C	226	SER	N-CA-C	7.10	118.78	110.41
1	A	226	SER	N-CA-C	6.93	118.58	110.41
1	B	194	ILE	CA-C-N	6.85	126.61	119.76
1	B	194	ILE	C-N-CA	6.85	126.61	119.76
1	B	72	GLY	N-CA-C	6.57	124.10	115.36
1	C	72	GLY	N-CA-C	6.50	124.34	115.30
1	B	256	GLU	N-CA-C	-6.32	103.47	112.13
1	C	297	VAL	N-CA-C	6.28	116.44	110.53
1	A	72	GLY	N-CA-C	6.21	123.93	115.30
1	A	278	MET	N-CA-C	-6.19	104.61	111.36
1	A	246	HIS	N-CA-C	6.17	118.33	109.14
1	A	313	ILE	N-CA-C	6.10	116.90	108.84
1	C	246	HIS	N-CA-C	6.01	117.91	108.96
1	C	214	VAL	N-CA-C	5.96	117.29	108.65
1	A	256	GLU	N-CA-C	-5.89	104.06	112.13
1	C	313	ILE	N-CA-C	5.72	117.01	108.54
1	C	233	ALA	N-CA-C	5.64	117.65	109.07
1	B	313	ILE	N-CA-C	5.59	116.22	108.84
1	A	120	ASP	CA-C-N	5.41	125.75	119.47
1	A	120	ASP	C-N-CA	5.41	125.75	119.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	14	LEU	N-CA-C	-5.41	103.36	110.39
1	C	256	GLU	N-CA-C	-5.35	104.42	111.96
1	B	321	ASP	N-CA-C	5.30	119.85	113.17
1	A	279	ASP	N-CA-C	-5.26	105.55	111.28
1	A	53	PHE	N-CA-C	5.17	119.54	113.23
1	B	195	PRO	N-CA-C	5.14	119.26	111.19
1	A	14	LEU	N-CA-C	-5.12	103.23	110.29
1	A	321	ASP	N-CA-C	5.07	119.55	113.17
1	B	53	PHE	N-CA-C	5.07	119.41	113.23
1	B	279	ASP	N-CA-C	-5.04	105.79	111.28
1	C	53	PHE	N-CA-C	5.03	119.37	113.23
1	C	14	LEU	N-CA-C	-5.03	103.36	110.29

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	2641	0	2577	89	0
1	B	2641	0	2577	80	0
1	C	2569	0	2515	91	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
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	260	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 17.

All (260) 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:C:205:LYS:HB3	1:C:205:LYS:HZ3	1.19	1.05
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:156:MET:HE3	1:C:228:MET:SD	2.16	0.85
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:156:MET:HE3	1:A:228:MET:SD	2.17	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:B:156:MET:HE2	1:B:156:MET:HA	1.64	0.78
1:C:211:ASN:ND2	1:C:212:THR:H	1.81	0.78
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:177:LEU:HD23	1:B:192:MET:HE3	1.71	0.73
1:B:174:PHE:HA	1:B:177:LEU:HD13	1.71	0.73
1:A:156:MET:HE2	1:A:156:MET:HA	1.72	0.72
1:A:214:VAL:HG11	1:A:234:VAL:CG2	2.17	0.72
1:B:171:THR:HG21	1:B:194:ILE:HD13	1.72	0.71
1:B:156:MET:HE3	1:B:228:MET:SD	2.30	0.71
1:B:295:LYS:O	1:B:299:GLU:HG3	1.91	0.70
1:B:156:MET:HE1	1:B:220:ILE:HG12	1.72	0.70
1:C:156:MET:HE2	1:C:156:MET:HA	1.72	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:A:156:MET:HE1	1:A:220:ILE:HG12	1.76	0.67
1:C:10:TYR:CE2	1:C:225:VAL:HG11	2.30	0.67
1:A:4:PHE:HB3	1:A:28:LEU:HD13	1.77	0.67
1:C:211:ASN:N	1:C:211:ASN:HD22	1.92	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
1:A:331:ILE:O	1:A:335:GLU:HG3	1.96	0.66
1:C:205:LYS:HB3	1:C:205:LYS:NZ	2.05	0.66
1:A:245:ILE:HG23	1:A:271:PHE:CD2	2.31	0.66
1:C:4:PHE:HB3	1:C:28:LEU:HD13	1.77	0.66
1:B:331:ILE:O	1:B:335:GLU:HG3	1.97	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:C:205:LYS:HZ3	1:C:205:LYS:CB	2.03	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:A:290:LEU:N	1:A:290:LEU:HD22	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:163:TYR:O	1:C:166:ARG:HG3	2.01	0.60
1:C:246:HIS:H	1:C:270:ASN:ND2	1.98	0.60
1:A:338:GLU:O	1:A:342:LEU:HD23	2.01	0.60
1:B:286:GLN:HA	1:B:290:LEU:HD22	1.83	0.60
1:C:162:GLN:HG3	1:C:288:GLN:NE2	2.17	0.60
1:B:162:GLN:HE22	1:B:171:THR:H	1.50	0.59
1:C:156:MET:HE1	1:C:220:ILE:HG12	1.84	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:A:143:HIS:HE1	1:A:256:GLU:OE1	1.87	0.57
1:A:178:GLU:HG3	1:A:193:LYS:HB2	1.87	0.57
1:A:295:LYS:O	1:A:299:GLU:HG3	2.04	0.57
1:B:177:LEU:HG	1:B:192:MET:SD	2.44	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:A:315:PRO:HB2	1:A:317:MET:CE	2.36	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:240:PRO:HG2	1:B:243:GLN:HG2	1.88	0.56
1:B:174:PHE:O	1:B:177:LEU:HD13	2.05	0.56
1:C:205:LYS:HZ1	1:C:206:PHE:HE1	1.54	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:HD12	1:B:158:LEU:N	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:A:25:ARG:HD2	1:A:262:HIS:CE1	2.42	0.54
1:C:214:VAL:HG12	1:C:236:PRO:CA	2.37	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:C:270:ASN:C	1:C:270:ASN:HD22	2.15	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:A:315:PRO:HB2	1:A:317:MET:HE3	1.90	0.53
1:C:71:HIS:HB2	2:C:501:FES:S1	2.48	0.53
1:C:315:PRO:HB2	1:C:317:MET:HE3	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:A:270:ASN:HD22	1:A:270:ASN:C	2.18	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
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: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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:333:LYS:O	1:B:337:LEU:HG	2.11	0.51
1:C:206:PHE:CE2	1:C:246:HIS:HA	2.46	0.51
1:B:38:PRO:HD2	5:B:1113:HOH:O	2.09	0.51
1:C:242:GLU:CD	1:C:242:GLU:H	2.18	0.51
1:C:206:PHE:HE2	1:C:246:HIS:HA	1.75	0.51
1:C:211:ASN:ND2	1:C:211:ASN:N	2.53	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:B:2:ALA:HB1	1:B:258:GLU:OE1	2.10	0.50
1:C:65:HIS:HA	5:C:842:HOH:O	2.10	0.50
1:B:156:MET:HE1	1:B:220:ILE:CG1	2.40	0.50
1:B:214:VAL:HG23	1:B:235:ALA:C	2.36	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:270:ASN:C	1:A:270:ASN:ND2	2.70	0.49
1:C:270:ASN:ND2	1:C:270:ASN:C	2.70	0.49
1:B:295:LYS:HE2	5:B:1146:HOH:O	2.11	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:A:156:MET:HE2	1:A:156:MET:CA	2.42	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:C:162:GLN:HG3	1:C:288:GLN:O	2.13	0.49
1:A:191:LEU:HD22	1:A:191:LEU:N	2.28	0.49
1:B:201:VAL:O	1:B:205:LYS:HG3	2.13	0.49
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:B:315:PRO:HB2	1:B:317:MET:CE	2.42	0.48
1:B:315:PRO:HB2	1:B:317:MET:HE3	1.96	0.48
1:B:292:LYS:O	1:B:292:LYS:HG3	2.13	0.48
1:B:174:PHE:CA	1:B:177:LEU:HD13	2.40	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
1:A:171:THR:HG21	1:A:194:ILE:HD13	1.97	0.47
1:C:224:LYS:HG2	1:C:225:VAL:HG23	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:224:LYS:HG2	1:A:225:VAL:HG23	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:156:MET:HE2	1:B:156:MET:CA	2.39	0.47
1:B:177:LEU:HD23	1:B:192:MET:CE	2.43	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:177:LEU:HD21	1:A:179:ARG:HH21	1.80	0.47
1:A:237:GLU:O	1:A:239:THR:N	2.47	0.47
1:B:76:ASP:OD2	1:B:80:GLN:HB3	2.14	0.47
1:B:282:LEU:HD13	5:B:899:HOH:O	2.14	0.47
1:A:236:PRO:O	1:A:237:GLU:C	2.57	0.46
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:B:224:LYS:HG2	1:B:225:VAL:HG23	1.97	0.46
1:C:179:ARG:O	1:C:180:GLU:HB2	2.16	0.46
1:C:194:ILE:HB	1:C:216:ALA:HB3	1.97	0.46
1:A:333:LYS:O	1:A:337:LEU:HG	2.15	0.46
1:A:148:TYR:CG	1:A:255:PRO:HG3	2.52	0.45
1:A:214:VAL:HG13	1:A:235:ALA:C	2.39	0.45
1:B:270:ASN:C	1:B:270:ASN:HD22	2.23	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: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:200:SER:HB2	5:A:916:HOH:O	2.17	0.45
1:C:76:ASP:OD2	1:C:80:GLN:HB3	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:C:158:LEU:HD13	1:C:158:LEU:C	2.42	0.45
1:A:161:ALA:HA	1:A:165:HIS:CD2	2.52	0.45
1:A:168:ASN:ND2	1:A:202:LEU:HB2	2.32	0.45
1:A:250:THR:O	1:A:265:PHE:HA	2.17	0.45
1:B:270:ASN:ND2	1:B:270:ASN:C	2.74	0.45
1:C:156:MET:HE2	1:C:156:MET:CA	2.42	0.45
1:C:220:ILE:HA	1:C:230:ASN:HA	1.99	0.45
1:A:305:ARG:O	1:A:308:VAL:HG22	2.18	0.44
1:C:156:MET:HE1	1:C:220:ILE:CG1	2.47	0.44
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:A:158:LEU:HD22	1:A:158:LEU:H	1.83	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: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:211:ASN:HD22	1:C:211:ASN:H	1.65	0.44
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:C:137:THR:HA	1:C:267:SER:O	2.17	0.43
1:B:158:LEU:H	1:B:158:LEU:CD1	2.28	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:B:10:TYR:CZ	1:B:225:VAL:HG11	2.53	0.43
1:B:158:LEU:HD11	1:B:220:ILE:HG21	2.01	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:A:269:ARG:HH11	1:A:282:LEU:HD12	1.83	0.43
1:A:201:VAL:O	1:A:205:LYS:HG3	2.18	0.43
1:C:156:MET:CE	1:C:220:ILE:HD13	2.49	0.43
1:B:191:LEU:HD23	1:B:219:ASP:HB3	2.00	0.43
1:A:225:VAL:HG12	1:A:226:SER:N	2.34	0.43
4:A:650:ACT:H3	4:A:651:ACT:H2	2.00	0.43
1:A:168:ASN:ND2	1:A:285:TRP:CE3	2.86	0.42
1:A:220:ILE:HA	1:A:230:ASN:HA	2.00	0.42
1:B:214:VAL:HG22	1:B:215:ASP:N	2.34	0.42
1:A:242:GLU:C	1:A:244:SER:H	2.27	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: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
1:A:203:MET:O	1:A:206:PHE:HB2	2.20	0.41
1:A:267:SER:OG	1:A:282:LEU:HD13	2.20	0.41
1:A:288:GLN:HE21	1:A:293:GLU:CD	2.28	0.41
1:A:341:ARG:HD2	1:A:341:ARG:HA	1.71	0.41
1:A:305:ARG:HA	1:A:308:VAL:HG22	2.02	0.41
1:B:217:TRP:CZ2	1:B:241:LYS:HB3	2.56	0.41
1:C:156:MET:HE1	1:C:220:ILE:CD1	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:195:PRO:O	1:C:197:GLY:N	2.43	0.41
1:B:174:PHE:C	1:B:177:LEU:HD13	2.45	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:36:ARG:NE	5:C:1013:HOH:O	2.27	0.41
1:C:224:LYS:O	1:C:225:VAL:HB	2.19	0.41
1:A:177:LEU:HD21	1:A:192:MET:HE3	2.03	0.41
1:B:124:ILE:HA	1:B:125:PRO:HD3	1.96	0.41
1:C:146:CYS:HB2	1:C:150:LEU:HD12	2.02	0.41
1:B:234:VAL:O	1:B:244:SER:OG	2.39	0.40
1:C:28:LEU:N	1:C:28:LEU:HD23	2.37	0.40
1:A:49:CYS:HA	1:A:50:PRO:HD3	1.90	0.40
1:B:148:TYR:CG	1:B:255:PRO:HG3	2.56	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

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	339/349 (97%)	313 (92%)	19 (6%)	7 (2%)	5	1
1	B	339/349 (97%)	317 (94%)	18 (5%)	4 (1%)	10	4
1	C	328/349 (94%)	296 (90%)	20 (6%)	12 (4%)	2	0
All	All	1006/1047 (96%)	926 (92%)	57 (6%)	23 (2%)	5	1

All (23) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	237	GLU
1	C	161	ALA

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Mol	Chain	Res	Type
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%)	51	47
1	B	276/284 (97%)	272 (99%)	4 (1%)	59	56
1	C	269/284 (95%)	263 (98%)	6 (2%)	45	40
All	All	821/852 (96%)	806 (98%)	15 (2%)	51	47

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	LEU

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Mol	Chain	Res	Type
1	A	234	VAL
1	A	239	THR
1	A	270	ASN
1	A	290	LEU
1	B	28	LEU
1	B	270	ASN
1	B	290	LEU
1	B	342	LEU
1	C	28	LEU
1	C	205	LYS
1	C	211	ASN
1	C	213	PRO
1	C	242	GLU
1	C	270	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	63	ASN
1	A	80	GLN
1	A	83	HIS
1	A	143	HIS
1	A	168	ASN
1	A	211	ASN
1	A	223	ASN
1	A	243	GLN
1	A	270	ASN
1	A	288	GLN
1	B	63	ASN
1	B	80	GLN
1	B	143	HIS
1	B	162	GLN
1	B	223	ASN
1	B	243	GLN
1	B	270	ASN
1	C	63	ASN
1	C	80	GLN
1	C	143	HIS
1	C	211	ASN
1	C	223	ASN
1	C	270	ASN
1	C	286	GLN

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Mol	Chain	Res	Type
1	C	288	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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	$\# Z > 2$
4	ACT	B	653	-	3,3,3	0.28	0	3,3,3	0.70	0
4	ACT	A	651	-	3,3,3	0.52	0	3,3,3	0.55	0
2	FES	C	501	1	0,4,4	-	-	-		
4	ACT	B	652	3	3,3,3	0.34	0	3,3,3	0.69	0
2	FES	A	501	1	0,4,4	-	-	-		
2	FES	B	501	1	0,4,4	-	-	-		
4	ACT	A	650	3	3,3,3	0.40	0	3,3,3	0.73	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	FES	C	501	1	-	-	0/1/1/1
2	FES	A	501	1	-	-	0/1/1/1
2	FES	B	501	1	-	-	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
4	B	653	ACT	1	0
4	A	651	ACT	5	0
2	C	501	FES	1	0
2	A	501	FES	1	0
2	B	501	FES	1	0
4	A	650	ACT	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2		OWAB(Å ²)	Q<0.9
1	A	341/349 (97%)	0.72	39 (11%)	10 11	22, 39, 73, 80	0
1	B	341/349 (97%)	0.70	28 (8%)	17 20	24, 37, 63, 85	0
1	C	332/349 (95%)	0.86	50 (15%)	5 6	23, 35, 73, 82	0
All	All	1014/1047 (96%)	0.76	117 (11%)	9 11	22, 37, 70, 85	0

All (117) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	207	LEU	6.9
1	C	200	SER	6.6
1	C	177	LEU	6.3
1	C	212	THR	5.8
1	B	339	ALA	5.7
1	C	203	MET	5.7
1	A	273	ILE	5.3
1	A	225	VAL	5.1
1	C	312	GLY	5.0
1	C	206	PHE	5.0
1	C	199	PRO	4.9
1	C	211	ASN	4.6
1	B	210	ALA	4.5
1	B	340	ALA	4.5
1	A	340	ALA	4.5
1	A	339	ALA	4.4
1	C	337	LEU	4.2
1	A	211	ASN	4.2
1	C	213	PRO	4.2
1	A	193	LYS	4.1
1	C	201	VAL	4.1
1	B	342	LEU	4.1
1	C	277	GLU	4.0

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Mol	Chain	Res	Type	RSRZ
1	C	159	GLY	4.0
1	B	3	THR	3.9
1	C	275	ASP	3.9
1	B	334	LEU	3.9
1	C	272	GLY	3.8
1	C	197	GLY	3.7
1	A	210	ALA	3.7
1	A	63	ASN	3.5
1	A	191	LEU	3.6
1	B	225	VAL	3.5
1	C	339	ALA	3.5
1	C	276	PRO	3.5
1	C	225	VAL	3.5
1	C	202	LEU	3.5
1	B	192	MET	3.5
1	A	172	ASP	3.4
1	C	342	LEU	3.4
1	C	240	PRO	3.4
1	C	209	GLY	3.4
1	C	340	ALA	3.4
1	B	239	THR	3.3
1	C	158	LEU	3.3
1	C	239	THR	3.2
1	A	337	LEU	3.1
1	C	160	HIS	3.1
1	A	271	PHE	3.1
1	B	240	PRO	3.1
1	A	198	THR	3.1
1	C	204	ALA	3.0
1	A	276	PRO	2.9
1	C	237	GLU	2.9
1	B	310	ALA	2.9
1	B	273	ILE	2.9
1	A	239	THR	2.9
1	A	342	LEU	2.9
1	C	273	ILE	2.9
1	A	158	LEU	2.8
1	B	337	LEU	2.8
1	C	196	GLY	2.8
1	B	212	THR	2.8
1	C	198	THR	2.8
1	A	308	VAL	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	277	GLU	2.7
1	A	240	PRO	2.6
1	B	258	GLU	2.6
1	C	178	GLU	2.6
1	A	111	ILE	2.5
1	A	195	PRO	2.5
1	B	211	ASN	2.5
1	A	314	ARG	2.5
1	B	93	ALA	2.5
1	C	224	LYS	2.5
1	C	40	GLY	2.5
1	C	2	ALA	2.5
1	C	214	VAL	2.5
1	C	238	GLY	2.4
1	A	168	ASN	2.4
1	C	185	ASP	2.4
1	C	341	ARG	2.4
1	A	197	GLY	2.4
1	C	62	VAL	2.4
1	A	241	LYS	2.4
1	B	175	ASP	2.3
1	B	311	ASN	2.3
1	A	224	LYS	2.3
1	C	122	GLY	2.3
1	A	238	GLY	2.3
1	A	214	VAL	2.3
1	A	192	MET	2.3
1	B	92	PRO	2.2
1	A	194	ILE	2.2
1	C	39	ASP	2.2
1	A	272	GLY	2.2
1	B	213	PRO	2.2
1	C	236	PRO	2.2
1	C	167	ALA	2.2
1	A	206	PHE	2.1
1	C	278	MET	2.1
1	C	210	ALA	2.1
1	A	174	PHE	2.1
1	A	275	ASP	2.1
1	B	242	GLU	2.1
1	A	318	LEU	2.1
1	B	282	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	90	ALA	2.1
1	C	161	ALA	2.1
1	B	338	GLU	2.1
1	A	185	ASP	2.1
1	C	192	MET	2.0
1	A	209	GLY	2.0
1	B	7	ASN	2.0
1	A	85	PRO	2.0
1	B	195	PRO	2.0
1	B	214	VAL	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 oligosaccharides 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. 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	B-factors(Å ²)	Q<0.9
4	ACT	A	650	4/4	0.90	0.12	37,38,38,40	0
4	ACT	A	651	4/4	0.91	0.09	39,42,42,43	0
4	ACT	B	652	4/4	0.91	0.10	34,35,35,36	0
4	ACT	B	653	4/4	0.95	0.12	35,37,37,38	0
2	FES	B	501	4/4	0.98	0.04	30,31,31,31	0
3	FE	A	502	1/1	0.98	0.04	39,39,39,39	0
2	FES	C	501	4/4	0.99	0.03	26,27,27,29	0
3	FE	B	502	1/1	0.99	0.04	31,31,31,31	0
2	FES	A	501	4/4	1.00	0.04	24,25,26,26	0

6.5 Other polymers [i](#)

There are no such residues in this entry.