



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 17, 2020 – 10:18 AM JST

PDB ID : 7C1H
Title : Crystal structure of the starter condensation domain of rhizomide synthetase RzmA
Authors : Zhong, L.; Diao, X.; Zhang, N.; Li, F.W.; Zhou, H.B.; Chen, H.N.; Ren, X.; Zhang, Y.; Wu, D.; Bian, X.
Deposited on : 2020-05-04
Resolution : 2.30 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.14.6
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.14.6

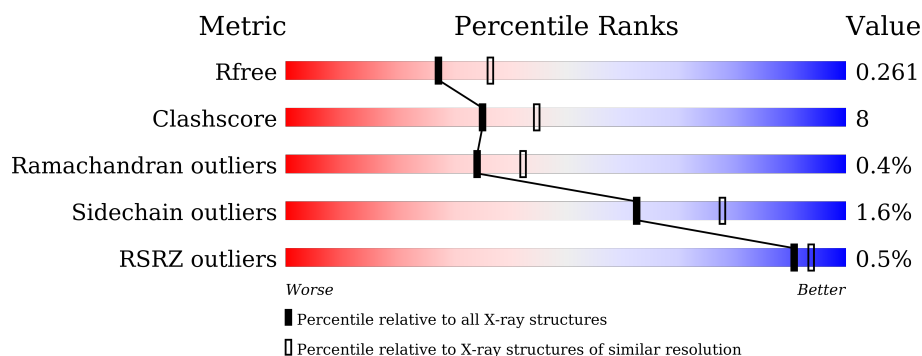
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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}	130704	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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 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	434	<div> <div>80%</div> <div>14%</div> <div>6%</div> </div>
1	B	434	<div> <div>70%</div> <div>20%</div> <div>10%</div> </div>
1	C	434	<div> <div>77%</div> <div>16%</div> <div>6%</div> </div>
1	D	434	<div> <div>%</div> <div>72%</div> <div>18%</div> <div>9%</div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 13327 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 Non-ribosomal peptide synthetase modules.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	410	Total	C	N	O	S	0	0	0
			3245	2057	566	601	21			
1	B	391	Total	C	N	O	S	0	0	0
			3107	1974	539	574	20			
1	C	408	Total	C	N	O	S	0	0	0
			3236	2053	565	597	21			
1	D	394	Total	C	N	O	S	0	0	0
			3125	1986	542	577	20			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP E5ATN9
A	-1	SER	-	expression tag	UNP E5ATN9
A	0	HIS	-	expression tag	UNP E5ATN9
B	-2	GLY	-	expression tag	UNP E5ATN9
B	-1	SER	-	expression tag	UNP E5ATN9
B	0	HIS	-	expression tag	UNP E5ATN9
C	-2	GLY	-	expression tag	UNP E5ATN9
C	-1	SER	-	expression tag	UNP E5ATN9
C	0	HIS	-	expression tag	UNP E5ATN9
D	-2	GLY	-	expression tag	UNP E5ATN9
D	-1	SER	-	expression tag	UNP E5ATN9
D	0	HIS	-	expression tag	UNP E5ATN9

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	172	Total	O	0	0
			172	172		
2	B	146	Total	O	0	0
			146	146		

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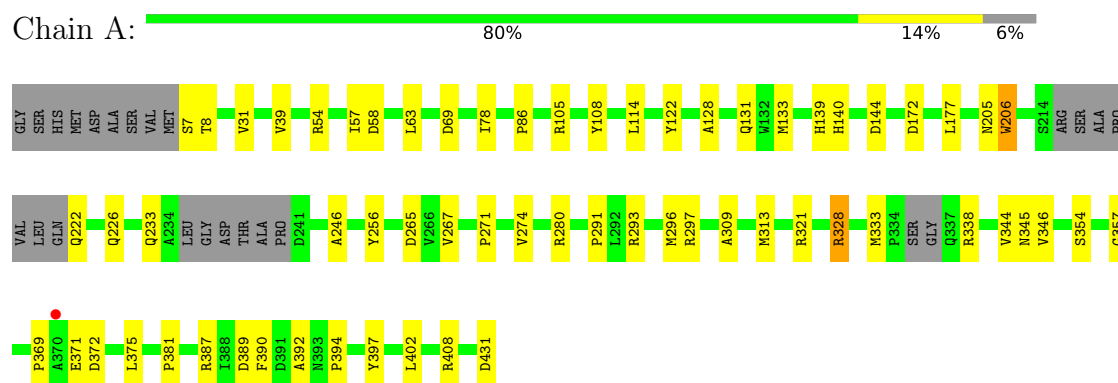
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	C	173	Total 173	O 173	0	0
2	D	123	Total 123	O 123	0	0

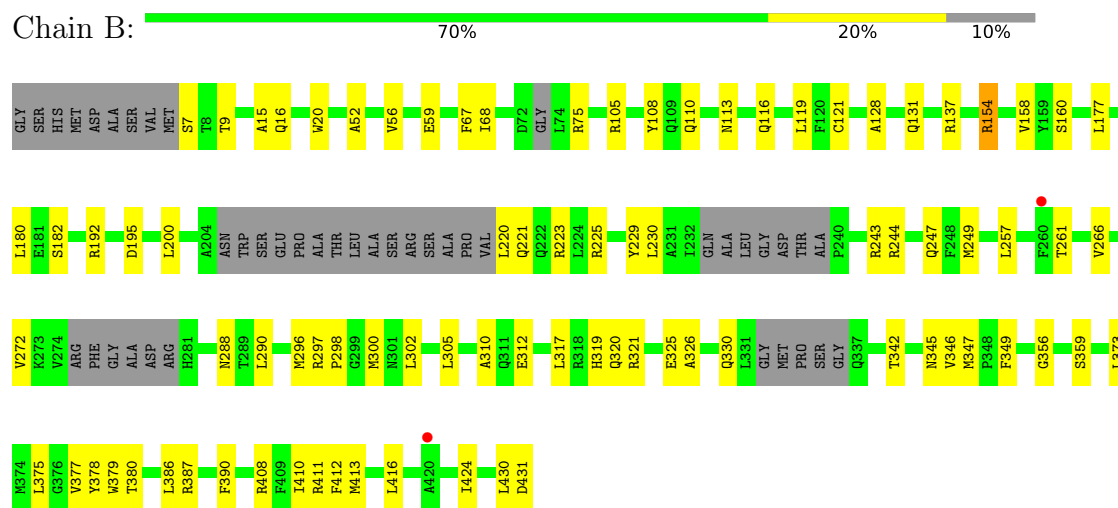
3 Residue-property plots [i](#)

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.

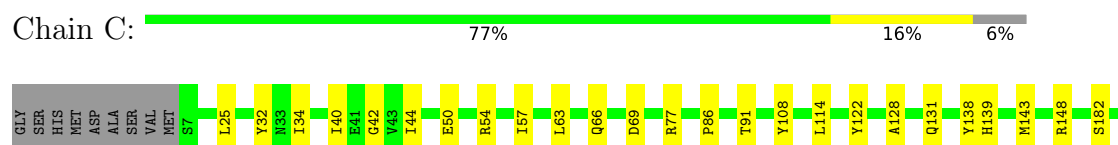
- Molecule 1: Non-ribosomal peptide synthetase modules

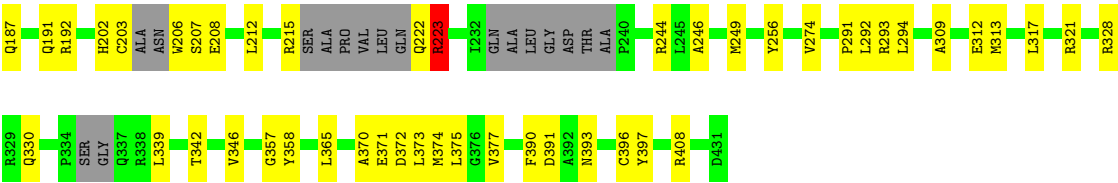


- Molecule 1: Non-ribosomal peptide synthetase modules



- Molecule 1: Non-ribosomal peptide synthetase modules





• Molecule 1: Non-ribosomal peptide synthetase modules



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	60.05Å 88.76Å 97.88Å 67.30° 89.90° 89.92°	Depositor
Resolution (Å)	49.97 – 2.30 48.44 – 2.30	Depositor EDS
% Data completeness (in resolution range)	86.8 (49.97-2.30) 43.0 (48.44-2.30)	Depositor EDS
R_{merge}	0.03	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.13 (at 2.29Å)	Xtriage
Refinement program	PHENIX 1.14	Depositor
R, R_{free}	0.227 , 0.253 0.223 , 0.261	Depositor DCC
R_{free} test set	1030 reflections (2.77%)	wwPDB-VP
Wilson B-factor (Å ²)	25.9	Xtriage
Anisotropy	0.072	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 18.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.499 for h,-k,-l	Xtriage
Reported twinning fraction	0.500 for h,-k,-l	Depositor
Outliers	0 of 37161 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	13327	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.69% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	0.26	0/3322	0.41	0/4520
1	B	0.25	0/3178	0.42	0/4321
1	C	0.24	0/3313	0.41	0/4505
1	D	0.25	0/3198	0.43	0/4350
All	All	0.25	0/13011	0.42	0/17696

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3245	0	3157	34	0
1	B	3107	0	3033	64	0
1	C	3236	0	3153	52	0
1	D	3125	0	3054	61	0
2	A	172	0	0	5	0
2	B	146	0	0	17	0
2	C	173	0	0	13	0
2	D	123	0	0	5	0
All	All	13327	0	12397	206	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 8.

All (206) 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:317:LEU:HA	2:C:612:HOH:O	1.78	0.81
1:B:302:LEU:HD22	1:B:416:LEU:HB3	1.61	0.81
1:D:230:LEU:HD21	1:D:232:ILE:HD13	1.69	0.74
1:D:232:ILE:HG21	2:D:600:HOH:O	1.88	0.74
1:B:379:TRP:HB2	1:B:386:LEU:HD12	1.70	0.73
1:B:220:LEU:HB2	2:B:585:HOH:O	1.88	0.72
1:B:20:TRP:HZ3	1:B:67:PHE:HB3	1.57	0.70
1:C:223:ARG:HD2	1:C:370:ALA:N	2.07	0.70
1:C:374:MET:HB2	1:C:391:ASP:HB2	1.75	0.69
1:A:69:ASP:OD2	1:B:243:ARG:NH1	2.27	0.68
1:B:302:LEU:HD13	1:B:416:LEU:HD22	1.75	0.68
1:C:309:ALA:O	1:C:313:MET:HG3	1.93	0.68
1:B:330:GLN:NE2	2:B:502:HOH:O	2.24	0.67
1:C:91:THR:OG1	2:C:501:HOH:O	2.12	0.67
1:B:128:ALA:HB3	1:B:131:GLN:HB3	1.77	0.67
1:D:379:TRP:HB2	1:D:386:LEU:HD12	1.77	0.66
1:D:20:TRP:HH2	1:D:115:THR:HG22	1.61	0.65
1:B:182:SER:OG	1:B:321:ARG:NH2	2.30	0.65
1:D:231:ALA:C	1:D:232:ILE:HD12	2.18	0.64
1:D:128:ALA:HB3	1:D:131:GLN:HB3	1.80	0.63
1:D:232:ILE:CG2	2:D:600:HOH:O	2.45	0.63
1:C:244:ARG:NH1	2:C:511:HOH:O	2.32	0.62
1:D:142:MET:SD	2:D:510:HOH:O	2.56	0.62
1:C:182:SER:OG	1:C:321:ARG:NH2	2.33	0.62
1:B:317:LEU:HA	1:B:320:GLN:HE21	1.65	0.62
1:B:290:LEU:HD11	1:B:320:GLN:HA	1.81	0.61
1:C:244:ARG:NH1	2:C:514:HOH:O	2.33	0.61
1:D:230:LEU:CD2	1:D:232:ILE:HD13	2.30	0.61
1:A:128:ALA:HB3	1:A:131:GLN:HB3	1.83	0.61
1:A:205:ASN:ND2	2:A:514:HOH:O	2.32	0.61
1:B:266:VAL:HG22	1:B:296:MET:HE1	1.83	0.60
1:D:231:ALA:HB2	1:D:385:GLN:HA	1.82	0.60
1:D:68:ILE:HG13	1:D:75:ARG:HB2	1.82	0.60
1:D:328:ARG:O	1:D:328:ARG:NH1	2.29	0.59
1:B:16:GLN:NE2	2:B:513:HOH:O	2.33	0.59
1:C:371:GLU:O	2:C:502:HOH:O	2.17	0.59
1:B:230:LEU:HD21	1:B:386:LEU:HD23	1.84	0.59
1:B:346:VAL:HG22	1:B:377:VAL:HB	1.85	0.59
1:D:264:GLN:HA	1:D:296:MET:HB2	1.84	0.59
1:A:140:HIS:HD2	2:A:595:HOH:O	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:219:VAL:HG22	1:D:220:LEU:HD12	1.84	0.58
1:B:244:ARG:NH2	2:B:520:HOH:O	2.36	0.57
1:D:231:ALA:O	1:D:232:ILE:HD12	2.05	0.57
1:D:7:SER:N	2:D:511:HOH:O	2.36	0.57
1:C:69:ASP:OD2	1:D:243:ARG:NH1	2.37	0.57
1:A:78:ILE:HG13	1:A:177:LEU:HD11	1.87	0.57
1:D:302:LEU:HD22	1:D:416:LEU:HB3	1.86	0.57
1:A:140:HIS:CD2	2:A:595:HOH:O	2.56	0.57
1:B:430:LEU:HA	2:B:544:HOH:O	2.03	0.57
1:A:246:ALA:HA	1:A:346:VAL:HG21	1.87	0.56
1:B:378:TYR:HB2	1:B:387:ARG:HB3	1.86	0.56
1:C:330:GLN:NE2	2:C:523:HOH:O	2.38	0.56
1:B:272:VAL:HA	2:B:527:HOH:O	2.06	0.56
1:B:113:ASN:HA	2:B:570:HOH:O	2.05	0.55
1:A:274:VAL:HG11	1:A:321:ARG:HG3	1.88	0.55
1:B:411:ARG:HD3	2:B:599:HOH:O	2.05	0.55
1:D:339:LEU:H	1:D:339:LEU:HD12	1.73	0.54
1:A:226:GLN:HG3	1:A:402:LEU:HD23	1.90	0.54
1:D:199:TRP:HB3	1:D:327:LEU:HD13	1.90	0.54
1:D:408:ARG:HD3	1:D:431:ASP:OD1	2.07	0.54
1:C:293:ARG:NH1	2:C:525:HOH:O	2.40	0.54
1:D:219:VAL:HG22	1:D:220:LEU:H	1.73	0.53
1:D:259:ARG:HD3	1:D:424:ILE:HG22	1.91	0.53
1:C:223:ARG:HB3	1:C:223:ARG:CZ	2.37	0.53
1:D:17:THR:O	1:D:21:LEU:HG	2.08	0.53
1:A:222:GLN:HB3	1:A:394:PRO:HG2	1.90	0.53
1:B:257:LEU:O	1:B:261:THR:OG1	2.20	0.53
1:C:393:ASN:HB3	1:C:396:CYS:HB2	1.90	0.52
1:B:290:LEU:HB2	2:B:518:HOH:O	2.08	0.52
1:A:265:ASP:HB3	1:A:293:ARG:NE	2.24	0.52
1:D:302:LEU:HB2	1:D:416:LEU:HD22	1.91	0.52
1:C:40:ILE:HG21	1:C:44:ILE:HD11	1.92	0.52
1:B:113:ASN:HB3	1:B:116:GLN:HB2	1.92	0.51
1:B:408:ARG:HE	1:B:431:ASP:HB3	1.74	0.51
1:C:202:HIS:HE1	2:C:563:HOH:O	1.93	0.51
1:B:229:TYR:HB3	1:C:50:GLU:HG3	1.91	0.51
1:A:206:TRP:HH2	1:A:291:PRO:HB2	1.76	0.51
1:A:338:ARG:HH12	1:A:371:GLU:HG2	1.76	0.51
1:D:326:ALA:O	1:D:330:GLN:HB2	2.11	0.51
1:B:410:ILE:HA	1:B:413:MET:HE3	1.92	0.50
1:B:177:LEU:HD23	1:B:180:LEU:HD12	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:207:SER:O	2:C:503:HOH:O	2.18	0.50
1:D:31:VAL:HA	1:D:220:LEU:HG	1.93	0.50
1:D:424:ILE:H	1:D:424:ILE:HD12	1.77	0.50
1:C:34:ILE:HD11	1:C:143:MET:HB3	1.93	0.50
1:D:259:ARG:NH1	1:D:424:ILE:O	2.45	0.50
1:B:412:PHE:CE2	1:B:416:LEU:HD11	2.47	0.49
1:B:7:SER:N	2:B:528:HOH:O	2.45	0.49
1:B:249:MET:HG3	1:B:377:VAL:HG21	1.93	0.49
1:C:42:GLY:HA3	1:C:358:TYR:CD2	2.48	0.49
1:D:375:LEU:HD13	1:D:390:PHE:HE1	1.78	0.49
1:A:172:ASP:HA	1:A:280:ARG:NH1	2.28	0.49
1:A:57:ILE:HA	1:A:63:LEU:HD12	1.94	0.49
1:A:256:TYR:HE1	1:A:408:ARG:HG2	1.77	0.48
1:D:17:THR:HG23	1:D:74:LEU:HD12	1.95	0.48
1:B:200:LEU:HD11	2:B:525:HOH:O	2.13	0.48
1:B:244:ARG:NH2	2:B:529:HOH:O	2.45	0.48
1:B:220:LEU:HD12	2:B:585:HOH:O	2.13	0.48
1:B:326:ALA:O	1:B:330:GLN:HG2	2.14	0.48
1:C:212:LEU:HB2	1:C:397:TYR:HE2	1.77	0.48
1:D:273:LYS:HD2	1:D:281:HIS:ND1	2.29	0.48
1:C:66:GLN:OE1	1:C:77:ARG:NH1	2.47	0.48
1:C:223:ARG:HH21	1:C:391:ASP:HA	1.79	0.47
1:C:330:GLN:HA	1:D:318:ARG:NH2	2.29	0.47
1:B:160:SER:OG	1:B:356:GLY:N	2.46	0.47
1:D:258:TYR:HB2	1:D:296:MET:HG3	1.95	0.47
1:D:301:ASN:OD1	1:D:304:SER:HB2	2.14	0.47
1:D:15:ALA:HB2	1:D:325:GLU:HG2	1.95	0.47
1:B:59:GLU:OE1	1:B:154:ARG:NE	2.30	0.47
1:C:128:ALA:HB3	1:C:131:GLN:HB3	1.97	0.47
1:A:114:LEU:HD21	1:A:139:HIS:CE1	2.50	0.47
1:C:222:GLN:O	1:C:223:ARG:HG3	2.15	0.47
1:A:7:SER:OG	1:A:8:THR:N	2.48	0.47
1:D:230:LEU:HD23	1:D:231:ALA:N	2.30	0.47
1:D:372:ASP:HA	1:D:393:ASN:HB2	1.97	0.46
1:C:292:LEU:HD11	1:C:312:GLU:HB3	1.97	0.46
1:C:202:HIS:CE1	2:C:563:HOH:O	2.67	0.46
1:D:322:TYR:HE2	1:D:327:LEU:HB2	1.80	0.46
1:A:381:PRO:HG3	2:A:658:HOH:O	2.14	0.46
1:B:349:PHE:H	1:B:380:THR:HG23	1.80	0.46
1:A:39:VAL:HG22	1:A:133:MET:HG2	1.97	0.46
1:B:154:ARG:O	1:B:158:VAL:HG22	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:272:VAL:HG12	1:B:288:ASN:O	2.16	0.45
1:C:108:TYR:CE2	1:C:365:LEU:HD22	2.51	0.45
1:C:138:TYR:HE2	1:C:148:ARG:HG3	1.81	0.45
1:D:20:TRP:O	1:D:24:GLN:HG2	2.17	0.45
1:A:267:VAL:HG22	1:A:293:ARG:HD3	1.98	0.45
1:C:328:ARG:HG2	1:C:339:LEU:HG	1.98	0.45
1:A:381:PRO:CG	2:A:658:HOH:O	2.64	0.45
1:B:195:ASP:OD1	1:B:319:HIS:ND1	2.41	0.45
1:C:57:ILE:HA	1:C:63:LEU:HD12	1.99	0.45
1:C:138:TYR:CE2	1:C:148:ARG:HG3	2.51	0.45
1:D:322:TYR:CE2	1:D:327:LEU:HB2	2.52	0.45
1:B:297:ARG:HB3	1:B:300:MET:HE3	1.98	0.45
1:B:192:ARG:NH1	2:B:524:HOH:O	2.41	0.45
1:B:325:GLU:H	1:B:325:GLU:CD	2.20	0.45
1:D:144:ASP:N	1:D:144:ASP:OD1	2.49	0.45
1:C:375:LEU:HD13	1:C:390:PHE:HE1	1.81	0.44
1:A:431:ASP:N	1:A:431:ASP:OD1	2.50	0.44
1:D:408:ARG:HD3	1:D:431:ASP:CG	2.37	0.44
1:D:405:HIS:CE1	1:D:408:ARG:HH21	2.36	0.44
1:B:305:LEU:HA	1:B:305:LEU:HD23	1.75	0.44
1:C:249:MET:HG3	1:C:377:VAL:HG21	2.00	0.44
1:C:256:TYR:HE1	1:C:408:ARG:HG2	1.83	0.44
1:B:15:ALA:HB2	1:B:325:GLU:HG2	1.99	0.44
1:A:31:VAL:HG22	1:A:369:PRO:HB3	1.99	0.44
1:C:346:VAL:HG22	1:C:377:VAL:HB	2.00	0.44
1:D:347:MET:O	1:D:378:TYR:HA	2.18	0.44
1:B:121:CYS:HB3	1:B:137:ARG:HB3	2.00	0.43
1:C:50:GLU:OE1	1:C:54:ARG:NH2	2.48	0.43
1:B:56:VAL:HG22	1:B:154:ARG:HD2	2.00	0.43
1:C:32:TYR:HB2	2:C:565:HOH:O	2.18	0.43
1:C:86:PRO:HD2	1:C:122:TYR:O	2.18	0.43
1:D:114:LEU:HD21	1:D:139:HIS:CE1	2.53	0.43
1:B:288:ASN:ND2	1:B:320:GLN:O	2.45	0.43
1:C:246:ALA:HA	1:C:346:VAL:HG21	2.01	0.43
1:D:232:ILE:HG23	2:D:596:HOH:O	2.17	0.43
1:B:247:GLN:OE1	1:B:310:ALA:HB2	2.19	0.43
1:C:274:VAL:HG11	1:C:321:ARG:HG3	1.99	0.43
1:B:225:ARG:NH2	2:B:519:HOH:O	2.52	0.43
1:D:229:TYR:CZ	1:D:387:ARG:HD2	2.53	0.43
1:B:52:ALA:HB1	1:B:158:VAL:HG23	2.00	0.43
1:C:342:THR:HA	1:C:373:LEU:O	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:291:PRO:O	2:C:504:HOH:O	2.20	0.43
1:D:310:ALA:HA	1:D:313:MET:HE2	2.00	0.43
1:B:105:ARG:HA	1:B:108:TYR:CZ	2.54	0.43
1:B:347:MET:O	1:B:378:TYR:HA	2.19	0.43
1:D:301:ASN:N	1:D:301:ASN:OD1	2.52	0.43
1:A:328:ARG:HG2	1:A:333:MET:SD	2.59	0.42
1:B:119:LEU:HA	1:B:119:LEU:HD23	1.74	0.42
1:D:355:PHE:O	1:D:358:TYR:HD1	2.02	0.42
1:C:114:LEU:HD21	1:C:139:HIS:NE2	2.35	0.42
1:C:25:LEU:O	1:D:307:GLN:NE2	2.41	0.42
1:C:191:GLN:NE2	2:C:538:HOH:O	2.52	0.42
1:A:309:ALA:O	1:A:313:MET:HG3	2.19	0.42
1:B:345:ASN:ND2	1:B:347:MET:SD	2.92	0.42
1:D:23:GLN:OE1	1:D:139:HIS:NE2	2.42	0.42
1:D:415:VAL:HG11	1:D:427:ILE:HG23	2.01	0.42
1:A:86:PRO:HD2	1:A:122:TYR:O	2.18	0.42
1:D:328:ARG:HA	1:D:328:ARG:HD2	1.78	0.42
1:B:375:LEU:HD13	1:B:390:PHE:HE1	1.83	0.42
1:B:300:MET:HG3	1:B:424:ILE:HD12	2.00	0.42
1:B:68:ILE:HG13	1:B:75:ARG:HB2	2.02	0.42
1:A:387:ARG:NH1	1:A:389:ASP:OD2	2.47	0.42
1:B:116:GLN:NE2	2:B:538:HOH:O	2.52	0.42
1:B:342:THR:HG23	1:B:373:LEU:HD22	2.01	0.42
1:D:80:THR:OG1	1:D:80:THR:O	2.32	0.41
1:A:144:ASP:OD1	1:A:144:ASP:N	2.53	0.41
1:A:271:PRO:HG2	1:A:345:ASN:OD1	2.19	0.41
1:C:203:CYS:HA	1:C:206:TRP:CD1	2.54	0.41
1:C:187:GLN:HA	1:C:192:ARG:HD2	2.02	0.41
1:D:183:ASP:O	1:D:187:GLN:HG2	2.21	0.41
1:B:9:THR:CG2	1:B:75:ARG:HB3	2.51	0.41
1:C:317:LEU:O	1:C:317:LEU:HD13	2.20	0.41
1:A:344:VAL:HG22	1:A:375:LEU:HD23	2.02	0.41
1:D:232:ILE:HD11	1:D:414:GLN:HE21	1.85	0.41
1:D:29:SER:OG	1:D:219:VAL:HG23	2.21	0.41
1:A:392:ALA:HB1	1:A:397:TYR:CD1	2.55	0.40
1:C:114:LEU:HD21	1:C:139:HIS:CE1	2.55	0.40
1:A:375:LEU:HD13	1:A:390:PHE:HE1	1.85	0.40
1:B:312:GLU:HA	2:B:532:HOH:O	2.21	0.40
1:C:294:LEU:HD21	1:C:309:ALA:HA	2.04	0.40
1:D:240:PRO:HB2	1:D:241:ASP:H	1.58	0.40
1:B:221:GLN:NE2	1:B:223:ARG:HD3	2.37	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:105:ARG:HA	1:A:108:TYR:CZ	2.56	0.40
1:D:342:THR:HA	1:D:373:LEU:O	2.21	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	402/434 (93%)	388 (96%)	12 (3%)	2 (0%)	29	35
1	B	379/434 (87%)	369 (97%)	9 (2%)	1 (0%)	41	50
1	C	398/434 (92%)	384 (96%)	11 (3%)	3 (1%)	19	23
1	D	384/434 (88%)	368 (96%)	15 (4%)	1 (0%)	41	50
All	All	1563/1736 (90%)	1509 (96%)	47 (3%)	7 (0%)	34	42

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	372	ASP
1	C	223	ARG
1	A	357	GLY
1	C	357	GLY
1	C	372	ASP
1	B	298	PRO
1	D	298	PRO

5.3.2 Protein sidechains [i](#)

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	336/354 (95%)	328 (98%)	8 (2%)	49	66
1	B	324/354 (92%)	321 (99%)	3 (1%)	78	89
1	C	336/354 (95%)	333 (99%)	3 (1%)	78	89
1	D	326/354 (92%)	319 (98%)	7 (2%)	53	70
All	All	1322/1416 (93%)	1301 (98%)	21 (2%)	62	78

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

Mol	Chain	Res	Type
1	A	54	ARG
1	A	58	ASP
1	A	206	TRP
1	A	233	GLN
1	A	296	MET
1	A	297	ARG
1	A	328	ARG
1	A	354	SER
1	B	110	GLN
1	B	154	ARG
1	B	359	SER
1	C	208	GLU
1	C	215	ARG
1	C	223	ARG
1	D	105	ARG
1	D	226	GLN
1	D	265	ASP
1	D	327	LEU
1	D	340	PHE
1	D	359	SER
1	D	400	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	140	HIS
1	B	226	GLN
1	B	247	GLN
1	D	414	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	410/434 (94%)	-0.33	1 (0%) 95 96	14, 28, 52, 66	0
1	B	391/434 (90%)	-0.16	2 (0%) 91 94	21, 38, 59, 72	0
1	C	408/434 (94%)	-0.29	0 100 100	14, 28, 53, 83	0
1	D	394/434 (90%)	-0.13	5 (1%) 77 81	20, 38, 59, 74	0
All	All	1603/1736 (92%)	-0.23	8 (0%) 91 94	14, 33, 56, 83	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	260	PHE	3.4
1	B	420	ALA	3.2
1	A	370	ALA	2.7
1	D	232	ILE	2.6
1	D	70	SER	2.6
1	D	302	LEU	2.5
1	D	420	ALA	2.3
1	D	258	TYR	2.1

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 monosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.