



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 11:03 PM GMT

PDB ID : 1W96
Title : CRYSTAL STRUCTURE OF BIOTIN CARBOXYLASE DOMAIN OF
ACETYL-COENZYME A CARBOXYLASE FROM SACCHAROMYCES
CEREVISIAE IN COMPLEX WITH SORAPHEN A
Authors : Shen, Y.; Volrath, S.L.; Weatherly, S.C.; Elich, T.D.; Tong, L.
Deposited on : 2004-10-06
Resolution : 1.80 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (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

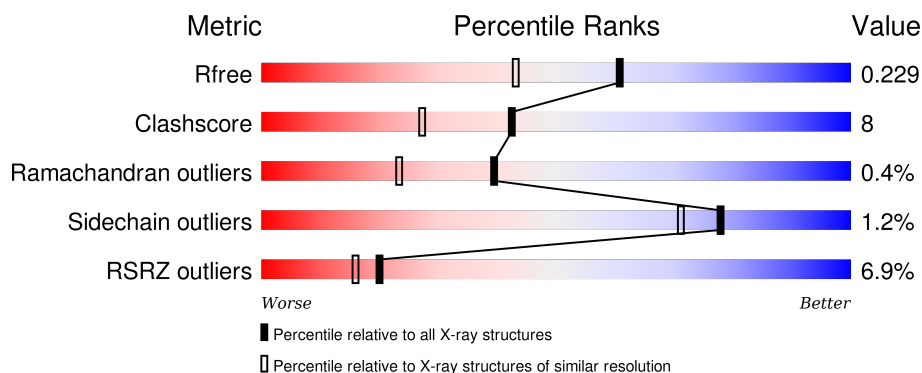
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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	4533 (1.80-1.80)
Clashscore	102246	5383 (1.80-1.80)
Ramachandran outliers	100387	5320 (1.80-1.80)
Sidechain outliers	100360	5319 (1.80-1.80)
RSRZ outliers	91569	4547 (1.80-1.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	554	<div> <div>6%</div> <div> <div></div> <div>81%</div> <div>17%</div> <div>..</div> </div> </div>
1	B	554	<div> <div>6%</div> <div> <div></div> <div>84%</div> <div>13%</div> <div>..</div> </div> </div>
1	C	554	<div> <div>7%</div> <div> <div></div> <div>80%</div> <div>16%</div> <div>..</div> </div> </div>

2 Entry composition [i](#)

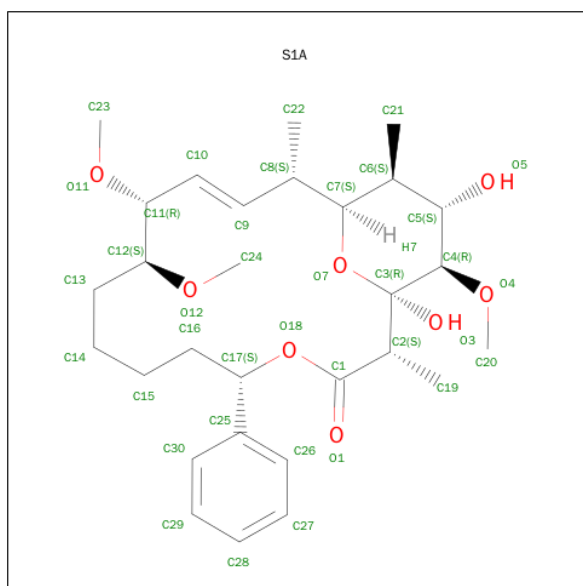
There are 3 unique types of molecules in this entry. The entry contains 14282 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 ACETYL-COENZYME A CARBOXYLASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	549	Total	C	N	O	S	0	0	0
			4290	2697	751	824	18			
1	B	542	Total	C	N	O	S	0	0	0
			4249	2674	746	811	18			
1	C	537	Total	C	N	O	S	0	0	1
			4187	2632	740	797	18			

- Molecule 2 is SORAPHEN A (three-letter code: S1A) (formula: C₂₉H₄₄O₈).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			37	29	8		
2	B	1	Total	C	O	0	0
			37	29	8		
2	C	1	Total	C	O	0	0
			37	29	8		

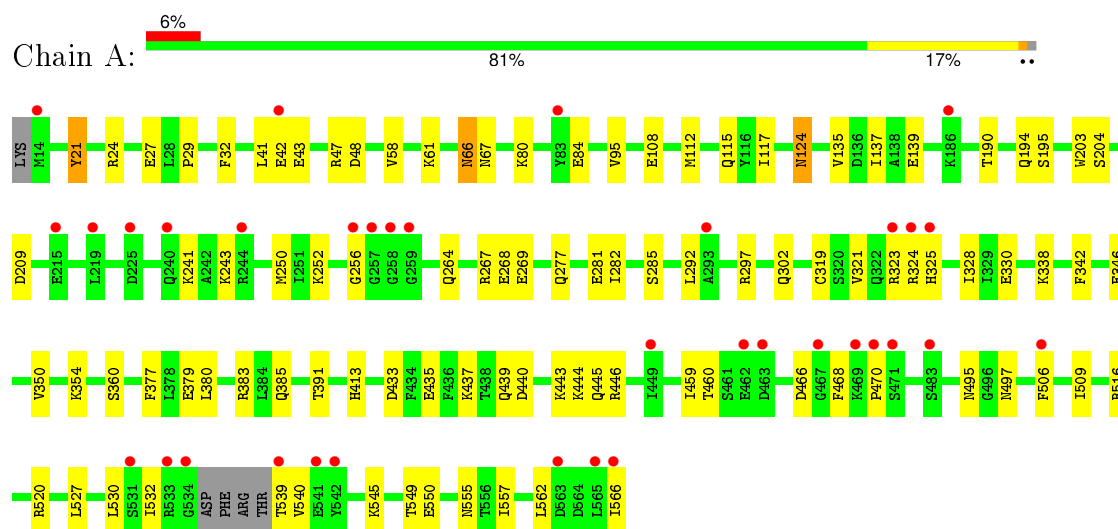
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	454	Total 454	O 454	0	0
3	B	510	Total 510	O 510	0	0
3	C	481	Total 481	O 481	0	0

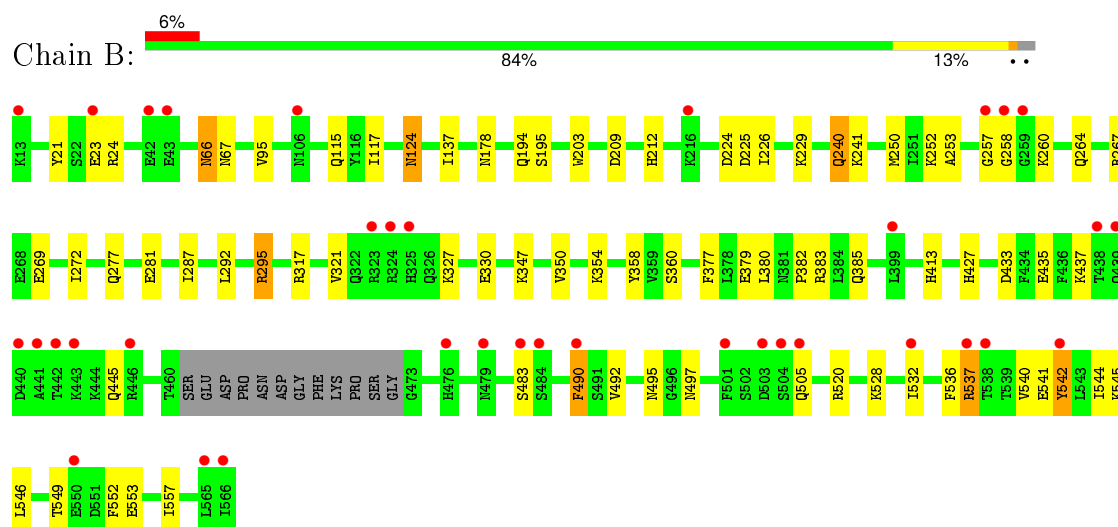
3 Residue-property plots

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.

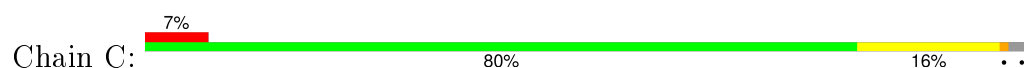
• Molecule 1: ACETYL-COENZYME A CARBOXYLASE

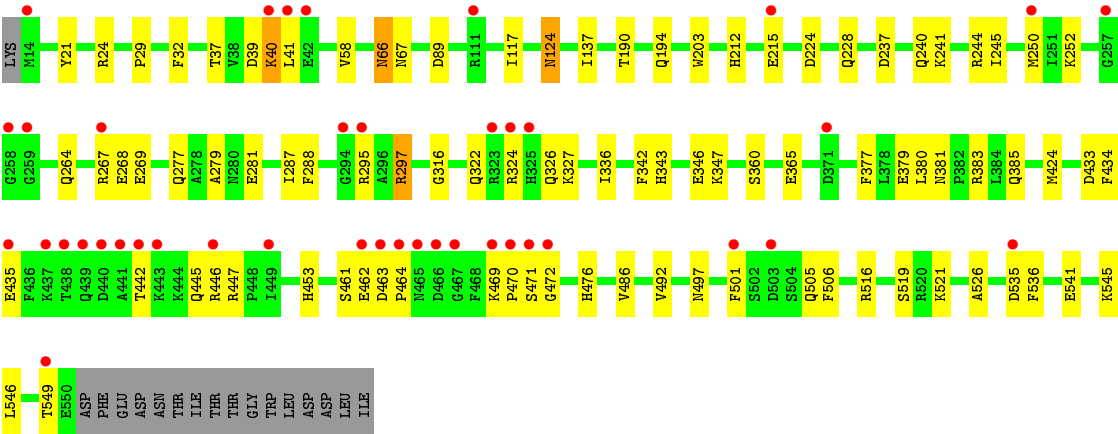


• Molecule 1: ACETYL-COENZYME A CARBOXYLASE



• Molecule 1: ACETYL-COENZYME A CARBOXYLASE





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	63.83Å 96.52Å 139.95Å 90.00° 96.82° 90.00°	Depositor
Resolution (Å)	29.20 – 1.80 29.20 – 1.79	Depositor EDS
% Data completeness (in resolution range)	96.3 (29.20-1.80) 95.8 (29.20-1.79)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.17 (at 1.79Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.195 , 0.230 0.195 , 0.229	Depositor DCC
R_{free} test set	11642 reflections (8.38%)	DCC
Wilson B-factor (Å ²)	19.2	Xtriage
Anisotropy	0.251	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 55.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 156180 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	14282	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.41% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: S1A

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	0.30	0/4381	0.56	0/5927
1	B	0.32	0/4338	0.56	0/5867
1	C	0.31	0/4277	0.56	0/5786
All	All	0.31	0/12996	0.56	0/17580

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

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	4290	0	4199	74	0
1	B	4249	0	4173	63	0
1	C	4187	0	4112	78	0
2	A	37	0	44	0	0
2	B	37	0	44	0	0
2	C	37	0	44	0	0
3	A	454	0	0	9	0
3	B	510	0	0	12	0
3	C	481	0	0	11	0
All	All	14282	0	12616	210	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 (210) 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:66:ASN:HD22	1:C:67:ASN:H	1.18	0.91
1:A:433:ASP:H	1:A:445:GLN:HE22	1.14	0.90
1:B:66:ASN:HD22	1:B:67:ASN:H	1.16	0.88
1:A:66:ASN:HD22	1:A:67:ASN:H	1.17	0.87
1:B:433:ASP:H	1:B:445:GLN:HE22	1.23	0.86
1:C:433:ASP:H	1:C:445:GLN:HE22	1.24	0.85
1:A:319:CYS:HB3	1:A:330:GLU:OE2	1.78	0.83
1:A:250:MET:HE1	1:A:264:GLN:HE21	1.48	0.77
1:A:43:GLU:HA	3:A:2056:HOH:O	1.82	0.77
1:A:443:LYS:HD2	1:C:244:ARG:NH2	2.00	0.76
1:C:462:GLU:HG2	1:C:464:PRO:HD3	1.70	0.74
1:A:47:ARG:HD2	3:A:2056:HOH:O	1.89	0.72
1:A:433:ASP:H	1:A:445:GLN:NE2	1.87	0.72
1:C:279:ALA:HB2	1:C:287:ILE:HD11	1.72	0.70
1:B:23:GLU:HG3	1:B:24:ARG:N	2.07	0.69
1:B:253:ALA:HA	1:B:287:ILE:HD13	1.73	0.69
1:B:250:MET:SD	1:B:264:GLN:HG2	2.32	0.69
1:B:433:ASP:H	1:B:445:GLN:NE2	1.91	0.69
1:B:528:LYS:O	1:B:532:ILE:HD13	1.92	0.69
1:C:383:ARG:HH12	1:C:385:GLN:HG2	1.59	0.68
1:B:117:ILE:HD12	1:B:137:ILE:HG23	1.75	0.67
1:A:269:GLU:H	1:A:269:GLU:CD	1.98	0.66
1:C:124:ASN:HD21	1:C:497:ASN:HD22	1.45	0.65
1:C:250:MET:SD	1:C:264:GLN:HG2	2.38	0.64
1:C:250:MET:HE1	1:C:264:GLN:HG2	1.80	0.63
1:C:117:ILE:HD12	1:C:137:ILE:HG23	1.80	0.63
1:B:224:ASP:HB2	1:C:215:GLU:OE1	1.99	0.63
1:B:545:LYS:O	1:B:549:THR:HG23	1.99	0.62
1:B:537:ARG:O	1:B:541:GLU:HB2	2.00	0.62
1:C:267:ARG:HB3	1:C:269:GLU:OE2	2.00	0.62
1:C:433:ASP:H	1:C:445:GLN:NE2	1.93	0.62
1:C:442:THR:HG23	3:C:2406:HOH:O	2.00	0.62
1:C:463:ASP:HB2	1:C:505:GLN:NE2	2.15	0.61
1:A:321:VAL:HG12	3:A:2280:HOH:O	2.01	0.61
1:B:212:HIS:HD2	3:C:2270:HOH:O	1.84	0.60
1:B:23:GLU:HG3	1:B:24:ARG:H	1.65	0.60
1:B:540:VAL:O	1:B:544:ILE:HG13	2.02	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:40:LYS:NZ	1:C:40:LYS:HB2	2.17	0.60
1:A:532:ILE:O	1:A:532:ILE:HG22	2.02	0.59
1:C:277:GLN:O	1:C:281:GLU:HG3	2.02	0.59
1:A:117:ILE:HD12	1:A:137:ILE:HG23	1.84	0.59
3:B:2279:HOH:O	1:C:212:HIS:HD2	1.85	0.59
1:B:492:VAL:HG12	1:B:505:GLN:HG2	1.84	0.58
1:A:267:ARG:HB3	1:A:269:GLU:OE2	2.04	0.58
1:A:446:ARG:HG2	1:A:446:ARG:HH11	1.68	0.58
1:C:424:MET:SD	1:C:445:GLN:HG2	2.44	0.58
1:B:240:GLN:HG3	1:B:241:LYS:N	2.18	0.58
1:A:250:MET:SD	1:A:264:GLN:HG2	2.44	0.57
1:B:490:PHE:CE1	1:B:505:GLN:HB2	2.39	0.57
1:C:461:SER:HB3	1:C:536:PHE:CZ	2.40	0.56
1:A:460:THR:O	1:A:539:THR:HG23	2.05	0.56
1:C:383:ARG:NH1	1:C:385:GLN:HG2	2.20	0.56
1:B:483:SER:HB3	3:B:2482:HOH:O	2.06	0.56
1:B:95:VAL:HG22	1:B:115:GLN:HG3	1.87	0.56
1:A:277:GLN:O	1:A:281:GLU:HG3	2.05	0.56
1:C:501:PHE:HA	3:C:2447:HOH:O	2.06	0.56
1:B:194:GLN:NE2	1:B:203:TRP:HE1	2.04	0.55
1:A:190:THR:O	1:A:194:GLN:HG3	2.05	0.55
1:B:542:TYR:HA	1:B:545:LYS:HD2	1.86	0.55
1:B:66:ASN:HD22	1:B:67:ASN:N	1.96	0.55
1:B:541:GLU:O	1:B:545:LYS:HG3	2.06	0.55
1:A:545:LYS:O	1:A:549:THR:HG23	2.07	0.55
1:C:463:ASP:HB2	1:C:505:GLN:HE21	1.71	0.54
1:B:250:MET:SD	1:B:292:LEU:HD13	2.48	0.54
1:A:256:GLY:HA2	1:A:282:ILE:HG13	1.90	0.54
1:A:194:GLN:NE2	1:A:203:TRP:HE1	2.06	0.54
1:C:240:GLN:HG3	1:C:241:LYS:N	2.23	0.54
1:B:194:GLN:HE22	1:B:203:TRP:HE1	1.55	0.53
1:B:277:GLN:O	1:B:281:GLU:HG3	2.09	0.53
1:C:66:ASN:HD22	1:C:67:ASN:N	1.98	0.53
1:A:323:ARG:HB2	1:A:562:LEU:HD23	1.91	0.53
1:C:250:MET:CE	1:C:264:GLN:HG2	2.39	0.53
1:C:41:LEU:HD12	3:C:2063:HOH:O	2.08	0.53
1:A:80:LYS:O	1:A:84:GLU:HG3	2.09	0.53
1:A:460:THR:HG22	1:A:506:PHE:HD2	1.74	0.52
1:C:462:GLU:HG2	1:C:464:PRO:CD	2.38	0.52
1:A:470:PRO:HG3	3:A:2397:HOH:O	2.08	0.52
1:B:124:ASN:HB3	1:B:495:ASN:O	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:324:ARG:HB2	3:C:2337:HOH:O	2.09	0.52
1:C:447:ARG:HG3	3:C:2414:HOH:O	2.09	0.52
1:C:194:GLN:NE2	1:C:203:TRP:HE1	2.08	0.51
1:C:269:GLU:CD	1:C:269:GLU:H	2.13	0.51
1:C:279:ALA:HB2	1:C:287:ILE:CD1	2.40	0.51
1:B:350:VAL:CG1	1:B:354:LYS:HE2	2.41	0.51
1:B:124:ASN:HD21	1:B:497:ASN:HD22	1.59	0.50
1:C:194:GLN:HE22	1:C:203:TRP:HE1	1.59	0.50
1:A:27:GLU:CD	1:B:229:LYS:HE2	2.31	0.50
1:B:225:ASP:OD1	1:B:226:ILE:HG12	2.11	0.50
1:A:440:ASP:OD1	1:A:444:LYS:HE3	2.11	0.50
1:A:328:ILE:HB	3:A:2280:HOH:O	2.12	0.50
1:C:250:MET:HB3	1:C:252:LYS:HZ3	1.77	0.50
1:B:536:PHE:HB2	1:B:540:VAL:HG13	1.94	0.50
1:C:461:SER:HB3	1:C:536:PHE:CE2	2.47	0.50
1:C:237:ASP:O	1:C:240:GLN:HG2	2.11	0.50
1:C:343:HIS:O	1:C:347:LYS:HG3	2.11	0.50
1:A:66:ASN:HD22	1:A:67:ASN:N	1.97	0.49
1:A:243:LYS:HE3	1:A:268:GLU:OE1	2.13	0.49
1:A:297:ARG:NH2	1:A:555:ASN:O	2.45	0.49
1:C:377:PHE:CZ	1:C:379:GLU:HA	2.47	0.49
1:C:37:THR:OG1	1:C:40:LYS:HG2	2.12	0.49
1:A:350:VAL:CG1	1:A:354:LYS:HE2	2.43	0.49
1:A:527:LEU:HD22	1:A:540:VAL:HB	1.94	0.49
1:B:536:PHE:O	1:B:540:VAL:HG22	2.12	0.49
1:C:492:VAL:HG22	3:C:2447:HOH:O	2.12	0.49
1:A:413:HIS:HE1	3:A:2360:HOH:O	1.95	0.49
1:B:269:GLU:CD	1:B:269:GLU:H	2.15	0.49
1:C:279:ALA:CB	1:C:287:ILE:HD11	2.40	0.48
1:C:486:VAL:HG11	1:C:526:ALA:CB	2.43	0.48
1:B:321:VAL:CG2	1:B:557:ILE:HD12	2.42	0.48
1:C:462:GLU:C	1:C:464:PRO:HD3	2.33	0.48
1:C:471:SER:HB2	1:C:476:HIS:CE1	2.48	0.47
1:A:194:GLN:HE22	1:A:203:TRP:HE1	1.61	0.47
1:A:377:PHE:CZ	1:A:379:GLU:HA	2.49	0.47
1:C:316:GLY:HA3	1:C:342:PHE:HZ	1.79	0.47
1:B:549:THR:O	1:B:553:GLU:HG3	2.15	0.47
1:A:282:ILE:HG22	1:A:285:SER:HB2	1.97	0.47
1:A:204:SER:HB2	1:A:241:LYS:HD3	1.97	0.47
1:A:459:ILE:HD13	1:A:509:ILE:HG13	1.97	0.46
1:C:252:LYS:HD2	1:C:288:PHE:CZ	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:521:LYS:C	1:C:521:LYS:HD3	2.36	0.46
1:A:124:ASN:HD21	1:A:497:ASN:HD22	1.64	0.46
1:C:268:GLU:HG3	3:C:2297:HOH:O	2.14	0.46
1:A:108:GLU:O	1:A:112:MET:HG3	2.16	0.46
1:A:27:GLU:OE1	1:B:229:LYS:HE2	2.16	0.46
1:B:490:PHE:CD1	1:B:536:PHE:HE2	2.33	0.46
1:B:350:VAL:HG12	1:B:354:LYS:HE2	1.97	0.46
1:B:295:ARG:HG3	3:B:2322:HOH:O	2.15	0.46
1:B:413:HIS:HE1	3:B:2013:HOH:O	1.98	0.46
1:A:250:MET:SD	1:A:292:LEU:HD13	2.56	0.46
1:A:319:CYS:HB3	1:A:330:GLU:CD	2.37	0.45
1:A:435:GLU:HB2	1:A:437:LYS:HG3	1.98	0.45
1:A:42:GLU:HG2	1:A:43:GLU:N	2.32	0.45
1:B:21:TYR:CD2	1:B:24:ARG:HD2	2.51	0.45
1:A:250:MET:HB3	1:A:252:LYS:NZ	2.32	0.45
1:C:541:GLU:CG	1:C:545:LYS:HE3	2.46	0.45
1:C:453:HIS:HD2	1:C:519:SER:OG	1.99	0.45
1:C:380:LEU:C	1:C:380:LEU:HD23	2.36	0.45
1:C:297:ARG:HG3	1:C:336:ILE:HD13	1.99	0.45
1:B:117:ILE:CD1	1:B:137:ILE:HG23	2.45	0.45
1:A:557:ILE:C	1:A:557:ILE:HD12	2.37	0.45
1:A:324:ARG:O	1:A:325:HIS:HB2	2.16	0.45
1:C:365:GLU:OE2	1:C:381:ASN:OD1	2.34	0.45
1:C:446:ARG:HG2	1:C:446:ARG:HH11	1.81	0.44
1:C:546:LEU:O	1:C:549:THR:HG22	2.18	0.44
1:B:260:LYS:HE3	3:B:2304:HOH:O	2.16	0.44
1:C:190:THR:O	1:C:194:GLN:HG3	2.18	0.44
1:A:21:TYR:CE2	1:A:24:ARG:HD2	2.51	0.44
1:B:347:LYS:NZ	3:B:2358:HOH:O	2.51	0.44
1:A:446:ARG:HG2	1:A:446:ARG:NH1	2.32	0.44
1:B:267:ARG:HB3	1:B:269:GLU:OE2	2.17	0.44
1:A:550:GLU:HB2	3:A:2441:HOH:O	2.17	0.44
1:C:462:GLU:HG2	1:C:464:PRO:CG	2.48	0.44
1:B:542:TYR:O	1:B:546:LEU:HG	2.17	0.44
1:A:466:ASP:OD1	1:A:468:PHE:HB2	2.17	0.44
1:B:377:PHE:CZ	1:B:379:GLU:HA	2.53	0.44
1:A:516:ARG:NH1	1:A:520:ARG:HH21	2.16	0.43
1:B:435:GLU:HB2	1:B:437:LYS:HG3	2.00	0.43
1:B:178:ASN:ND2	3:B:2226:HOH:O	2.50	0.43
1:B:272:ILE:HG12	3:B:2308:HOH:O	2.18	0.43
1:A:48:ASP:OD1	3:A:2056:HOH:O	2.21	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:195:SER:HA	1:B:209:ASP:HB2	2.00	0.43
1:C:326:GLN:OE1	1:C:326:GLN:HA	2.17	0.43
1:C:469:LYS:HA	1:C:470:PRO:HD3	1.89	0.43
1:C:21:TYR:CD2	1:C:24:ARG:HD2	2.55	0.42
1:A:95:VAL:HG22	1:A:115:GLN:CG	2.49	0.42
1:B:520:ARG:NH2	3:B:2481:HOH:O	2.52	0.42
1:A:29:PRO:HD2	1:A:32:PHE:CD2	2.53	0.42
1:A:95:VAL:HG22	1:A:115:GLN:HG2	2.01	0.42
1:C:41:LEU:HD23	1:C:58:VAL:HG21	2.02	0.42
1:A:302:GLN:HG2	1:A:391:THR:HG21	2.01	0.42
1:A:439:GLN:NE2	1:C:245:ILE:HG13	2.33	0.42
1:B:250:MET:HE3	1:B:252:LYS:HZ3	1.83	0.42
1:A:530:LEU:HD23	1:A:540:VAL:HG11	2.01	0.42
1:B:552:PHE:CD1	1:B:557:ILE:HD13	2.55	0.42
1:B:317:ARG:CZ	3:B:2335:HOH:O	2.66	0.42
1:A:323:ARG:O	1:A:324:ARG:HB2	2.19	0.42
1:A:124:ASN:HB3	1:A:495:ASN:O	2.19	0.42
1:C:342:PHE:O	1:C:346:GLU:HG3	2.19	0.42
1:A:195:SER:HA	1:A:209:ASP:HB2	2.01	0.42
1:A:383:ARG:NH1	1:A:385:GLN:HG2	2.35	0.42
1:C:453:HIS:HE1	3:C:2338:HOH:O	2.03	0.42
1:A:61:LYS:HE2	1:A:95:VAL:CG2	2.50	0.42
1:A:460:THR:HG22	1:A:506:PHE:CD2	2.53	0.41
1:C:541:GLU:HG2	1:C:545:LYS:HE3	2.02	0.41
1:A:566:ILE:O	1:A:566:ILE:HG22	2.20	0.41
1:B:327:LYS:HD3	1:B:330:GLU:OE2	2.20	0.41
1:B:380:LEU:C	1:B:380:LEU:HD23	2.41	0.41
1:B:483:SER:HB3	3:B:2455:HOH:O	2.21	0.41
1:A:380:LEU:HD23	1:A:380:LEU:C	2.41	0.41
1:C:433:ASP:C	1:C:435:GLU:H	2.24	0.41
1:A:135:VAL:O	1:A:139:GLU:HG3	2.20	0.41
1:C:250:MET:HB3	1:C:252:LYS:NZ	2.36	0.41
1:C:322:GLN:HG2	1:C:327:LYS:HA	2.02	0.41
1:C:224:ASP:O	1:C:228:GLN:HG2	2.21	0.41
1:C:250:MET:CE	1:C:264:GLN:HE21	2.34	0.41
1:A:41:LEU:HD23	1:A:58:VAL:HG21	2.03	0.41
1:C:39:ASP:HB2	3:C:2052:HOH:O	2.20	0.41
1:C:472:GLY:HA2	3:C:2427:HOH:O	2.21	0.41
1:B:358:TYR:OH	1:B:382:PRO:HA	2.21	0.40
1:A:338:LYS:HG2	3:A:2293:HOH:O	2.21	0.40
1:B:21:TYR:CE2	1:B:24:ARG:HD2	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:501:PHE:CD1	1:C:501:PHE:C	2.95	0.40
1:A:350:VAL:HG12	1:A:354:LYS:HE2	2.03	0.40
1:C:453:HIS:CD2	1:C:516:ARG:HA	2.57	0.40
1:C:29:PRO:HD2	1:C:32:PHE:CD2	2.56	0.40
1:B:383:ARG:NH1	1:B:385:GLN:HG2	2.37	0.40
1:A:342:PHE:O	1:A:346:GLU:HG3	2.21	0.40
1:B:427:HIS:HD2	3:B:2011:HOH:O	2.05	0.40
1:C:295:ARG:HD2	1:C:295:ARG:N	2.36	0.40
1:C:506:PHE:N	1:C:506:PHE:CD1	2.88	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	545/554 (98%)	523 (96%)	21 (4%)	1 (0%)	52	35
1	B	538/554 (97%)	519 (96%)	16 (3%)	3 (1%)	30	14
1	C	535/554 (97%)	516 (96%)	17 (3%)	2 (0%)	39	23
All	All	1618/1662 (97%)	1558 (96%)	54 (3%)	6 (0%)	39	23

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	258	GLY
1	A	360	SER
1	B	360	SER
1	C	360	SER
1	C	434	PHE
1	B	257	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	462/467 (99%)	459 (99%)	3 (1%)	90	88
1	B	457/467 (98%)	450 (98%)	7 (2%)	72	62
1	C	450/467 (96%)	444 (99%)	6 (1%)	76	68
All	All	1369/1401 (98%)	1353 (99%)	16 (1%)	78	71

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	21	TYR
1	A	66	ASN
1	A	124	ASN
1	B	66	ASN
1	B	124	ASN
1	B	240	GLN
1	B	295	ARG
1	B	490	PHE
1	B	537	ARG
1	B	542	TYR
1	C	40	LYS
1	C	66	ASN
1	C	89	ASP
1	C	124	ASN
1	C	297	ARG
1	C	535	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (53) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	31	HIS
1	A	66	ASN
1	A	106	ASN
1	A	124	ASN
1	A	152	HIS

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Mol	Chain	Res	Type
1	A	194	GLN
1	A	228	GLN
1	A	264	GLN
1	A	280	ASN
1	A	370	HIS
1	A	381	ASN
1	A	413	HIS
1	A	445	GLN
1	A	453	HIS
1	A	495	ASN
1	B	31	HIS
1	B	66	ASN
1	B	106	ASN
1	B	124	ASN
1	B	152	HIS
1	B	194	GLN
1	B	212	HIS
1	B	264	GLN
1	B	280	ASN
1	B	370	HIS
1	B	381	ASN
1	B	413	HIS
1	B	445	GLN
1	B	453	HIS
1	B	476	HIS
1	B	479	ASN
1	B	495	ASN
1	B	497	ASN
1	B	505	GLN
1	C	31	HIS
1	C	66	ASN
1	C	106	ASN
1	C	124	ASN
1	C	126	ASN
1	C	152	HIS
1	C	194	GLN
1	C	212	HIS
1	C	228	GLN
1	C	240	GLN
1	C	264	GLN
1	C	370	HIS
1	C	439	GLN

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Mol	Chain	Res	Type
1	C	445	GLN
1	C	453	HIS
1	C	476	HIS
1	C	485	ASN
1	C	495	ASN
1	C	505	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 ⓘ

3 ligands 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	S1A	A	1567	-	38,39,39	0.99	1 (2%)	42,55,55	1.27	6 (14%)
2	S1A	B	1567	-	38,39,39	1.05	4 (10%)	42,55,55	1.25	6 (14%)
2	S1A	C	1550	-	38,39,39	1.01	1 (2%)	42,55,55	1.28	7 (16%)

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	S1A	A	1567	-	-	0/44/67/67	0/1/3/3
2	S1A	B	1567	-	-	0/44/67/67	0/1/3/3
2	S1A	C	1550	-	-	0/44/67/67	0/1/3/3

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1567	S1A	C10-C9	2.04	1.38	1.32
2	B	1567	S1A	C5-C4	2.14	1.58	1.53
2	A	1567	S1A	C10-C9	2.27	1.39	1.32
2	B	1567	S1A	C3-C4	2.28	1.57	1.53
2	C	1550	S1A	C10-C9	2.36	1.39	1.32
2	B	1567	S1A	C25-C17	2.49	1.55	1.51

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1550	S1A	O18-C1-O1	-3.68	116.31	123.89
2	A	1567	S1A	O18-C1-O1	-3.59	116.49	123.89
2	B	1567	S1A	O18-C1-O1	-3.46	116.77	123.89
2	A	1567	S1A	C8-C7-C6	-3.04	110.51	115.69
2	C	1550	S1A	C8-C7-C6	-2.85	110.83	115.69
2	B	1567	S1A	C8-C7-C6	-2.79	110.93	115.69
2	B	1567	S1A	C16-C17-C25	-2.61	108.22	112.98
2	A	1567	S1A	C16-C17-C25	-2.53	108.37	112.98
2	C	1550	S1A	C16-C17-C25	-2.38	108.64	112.98
2	C	1550	S1A	C20-O4-C4	-2.36	109.88	114.39
2	B	1567	S1A	C20-O4-C4	-2.29	110.01	114.39
2	A	1567	S1A	C20-O4-C4	-2.23	110.12	114.39
2	C	1550	S1A	C11-C10-C9	-2.19	119.84	124.48
2	A	1567	S1A	C11-C10-C9	-2.17	119.89	124.48
2	B	1567	S1A	C11-C10-C9	-2.07	120.09	124.48
2	C	1550	S1A	C17-O18-C1	-2.01	112.63	116.90
2	A	1567	S1A	O18-C1-C2	2.36	113.76	110.83
2	C	1550	S1A	O18-C1-C2	2.45	113.86	110.83
2	B	1567	S1A	O18-C1-C2	2.61	114.07	110.83

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	549/554 (99%)	0.22	35 (6%) 23 18	10, 24, 47, 57	0
1	B	542/554 (97%)	0.22	36 (6%) 22 17	9, 21, 45, 55	0
1	C	537/554 (96%)	0.29	41 (7%) 17 13	12, 23, 48, 60	0
All	All	1628/1662 (97%)	0.24	112 (6%) 20 16	9, 23, 47, 60	0

All (112) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	325	HIS	8.0
1	B	257	GLY	7.8
1	B	566	ILE	7.5
1	C	258	GLY	7.1
1	C	323	ARG	6.1
1	C	472	GLY	5.8
1	C	294	GLY	5.8
1	C	324	ARG	5.7
1	A	257	GLY	5.6
1	B	258	GLY	5.6
1	C	549	THR	5.3
1	A	324	ARG	5.1
1	B	503	ASP	5.0
1	C	438	THR	4.8
1	A	566	ILE	4.7
1	A	534	GLY	4.6
1	A	14	MET	4.6
1	A	467	GLY	4.6
1	C	469	LYS	4.5
1	B	439	GLN	4.5
1	C	462	GLU	4.3
1	C	464	PRO	4.3
1	C	42	GLU	4.1

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Mol	Chain	Res	Type	RSRZ
1	B	324	ARG	4.1
1	C	439	GLN	4.1
1	B	501	PHE	4.0
1	C	463	ASP	4.0
1	B	438	THR	4.0
1	C	295	ARG	3.9
1	C	470	PRO	3.8
1	C	466	ASP	3.8
1	A	533	ARG	3.7
1	B	216	LYS	3.6
1	B	490	PHE	3.6
1	A	258	GLY	3.5
1	C	267	ARG	3.5
1	B	440	ASP	3.5
1	C	14	MET	3.3
1	B	565	LEU	3.2
1	B	484	SER	3.2
1	C	471	SER	3.2
1	C	257	GLY	3.1
1	C	371	ASP	3.1
1	B	23	GLU	3.0
1	A	462	GLU	3.0
1	C	41	LEU	3.0
1	A	471	SER	3.0
1	C	446	ARG	3.0
1	B	479	ASN	2.9
1	C	259	GLY	2.9
1	C	442	THR	2.9
1	B	476	HIS	2.9
1	A	244	ARG	2.8
1	C	441	ALA	2.8
1	A	542	TYR	2.8
1	B	542	TYR	2.8
1	B	441	ALA	2.8
1	A	256	GLY	2.7
1	C	465	ASN	2.7
1	C	40	LYS	2.7
1	B	106	ASN	2.7
1	A	463	ASP	2.7
1	C	449	ILE	2.7
1	A	323	ARG	2.7
1	B	483	SER	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	470	PRO	2.7
1	B	443	LYS	2.6
1	A	240	GLN	2.5
1	A	225	ASP	2.5
1	C	535	ASP	2.5
1	C	440	ASP	2.5
1	B	538	THR	2.4
1	C	501	PHE	2.4
1	A	219	LEU	2.4
1	B	505	GLN	2.4
1	C	435	GLU	2.4
1	A	186	LYS	2.3
1	C	111	ARG	2.3
1	A	449	ILE	2.3
1	B	537	ARG	2.3
1	C	503	ASP	2.3
1	A	83	TYR	2.3
1	A	42	GLU	2.2
1	A	541	GLU	2.2
1	A	259	GLY	2.2
1	B	325	HIS	2.2
1	B	259	GLY	2.2
1	C	467	GLY	2.2
1	A	531	SER	2.2
1	A	293	ALA	2.2
1	B	550	GLU	2.2
1	C	437	LYS	2.2
1	B	323	ARG	2.2
1	B	442	THR	2.2
1	A	563	ASP	2.1
1	A	483	SER	2.1
1	A	469	LYS	2.1
1	B	42	GLU	2.1
1	A	539	THR	2.1
1	A	506	PHE	2.1
1	A	325	HIS	2.1
1	C	215	GLU	2.1
1	A	215	GLU	2.1
1	B	532	ILE	2.1
1	C	443	LYS	2.1
1	C	250	MET	2.1
1	B	399	LEU	2.0

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Mol	Chain	Res	Type	RSRZ
1	B	504	SER	2.0
1	A	565	LEU	2.0
1	B	13	LYS	2.0
1	B	446	ARG	2.0
1	B	43	GLU	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, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q < 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	S1A	A	1567	37/37	0.93	0.14	0.72	13,17,24,26	0
2	S1A	B	1567	37/37	0.95	0.12	0.26	9,13,22,23	0
2	S1A	C	1550	37/37	0.94	0.12	-0.16	14,18,24,24	0

6.5 Other polymers [i](#)

There are no such residues in this entry.