



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 10, 2022 – 06:10 PM JST

PDB ID : 7DTQ
Title : Apo Crystal Structure of Octaketide Synthase from *A. arborescens*
Authors : Li, T.L.; Adhikari, K.; Li, Y.S.
Deposited on : 2021-01-06
Resolution : 1.75 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	:	2.25
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)	:	2.25

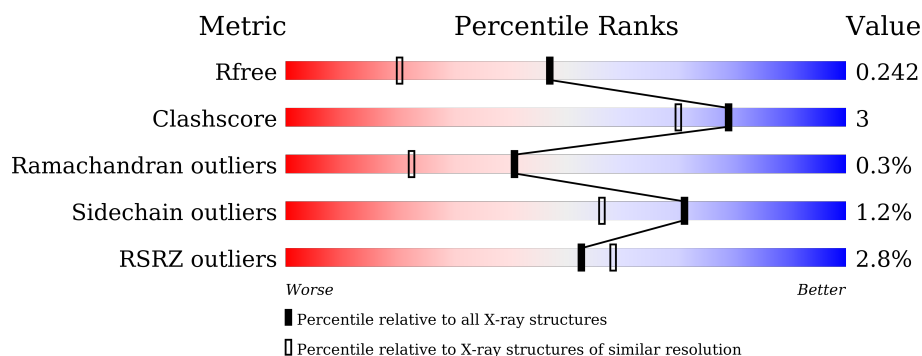
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.75 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	2340 (1.76-1.76)
Clashscore	141614	2466 (1.76-1.76)
Ramachandran outliers	138981	2437 (1.76-1.76)
Sidechain outliers	138945	2437 (1.76-1.76)
RSRZ outliers	127900	2298 (1.76-1.76)

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	403	<div> <div>3%</div> <div>90%</div> <div>6%</div> <div>••</div> </div>
1	B	403	<div> <div>2%</div> <div>90%</div> <div>6%</div> <div>•</div> </div>
1	C	403	<div> <div>2%</div> <div>91%</div> <div>5%</div> <div>•</div> </div>
1	D	403	<div> <div>4%</div> <div>89%</div> <div>7%</div> <div>•</div> </div>

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 12096 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.

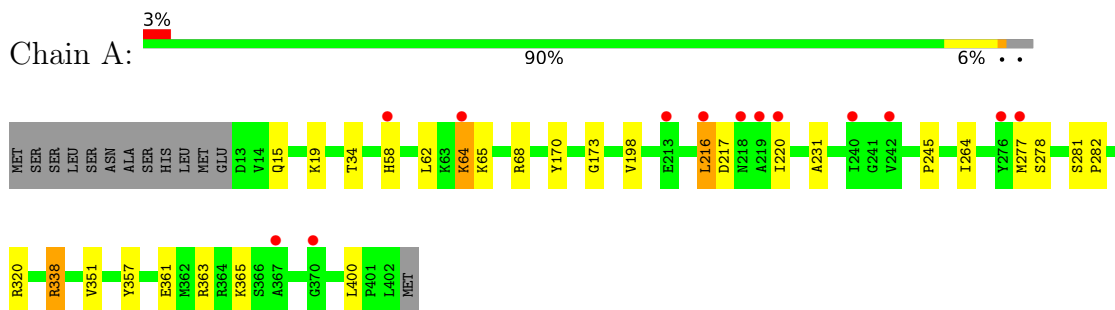
- Molecule 1 is a protein called Octaketide synthase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	390	Total	C	N	O	S	0	0	0
			3024	1920	517	562	25			
1	B	390	Total	C	N	O	S	0	0	0
			3024	1920	517	562	25			
1	C	390	Total	C	N	O	S	0	0	0
			3024	1920	517	562	25			
1	D	390	Total	C	N	O	S	0	0	0
			3024	1920	517	562	25			

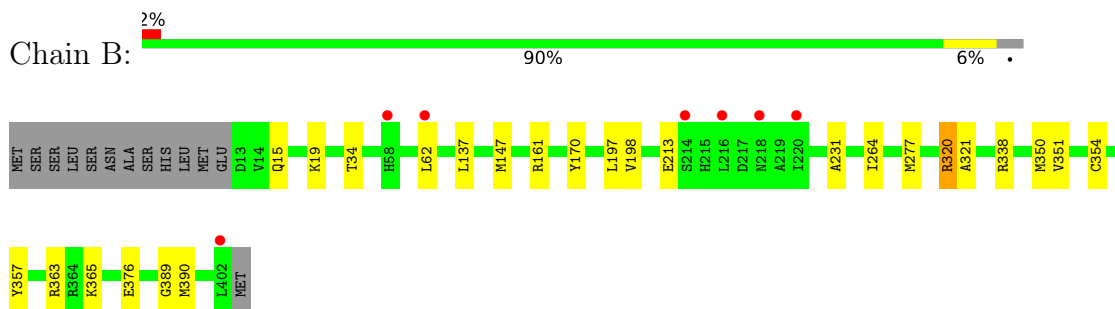
3 Residue-property plots [i](#)

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.

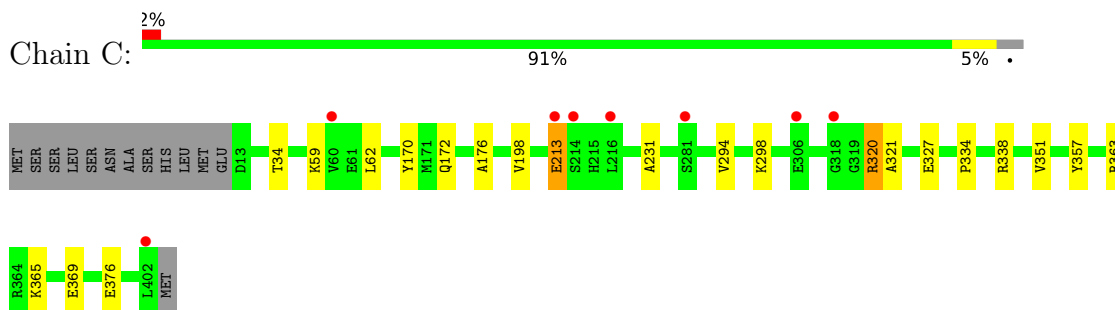
- Molecule 1: Octaketide synthase 1



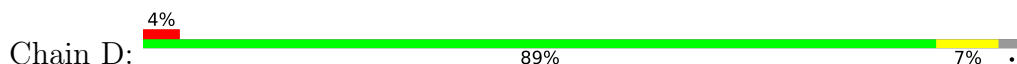
- Molecule 1: Octaketide synthase 1

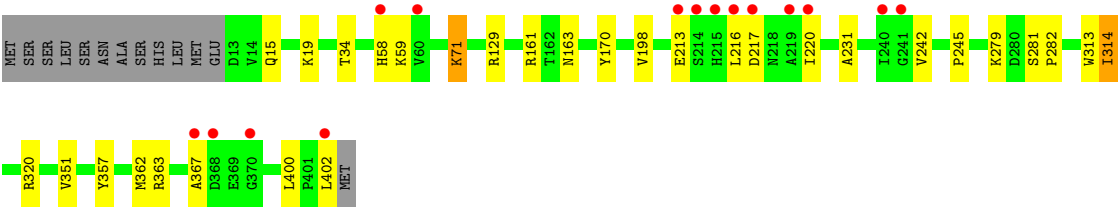


- Molecule 1: Octaketide synthase 1



- Molecule 1: Octaketide synthase 1





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	131.90Å 51.68Å 131.88Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	25.60 – 1.75 25.62 – 1.75	Depositor EDS
% Data completeness (in resolution range)	95.3 (25.60-1.75) 95.3 (25.62-1.75)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.86 (at 1.75Å)	Xtriage
Refinement program	PHENIX (1.11.1_2575: ???)	Depositor
R, R_{free}	0.225 , 0.242 0.225 , 0.242	Depositor DCC
R_{free} test set	2011 reflections (1.17%)	wwPDB-VP
Wilson B-factor (Å ²)	23.6	Xtriage
Anisotropy	0.127	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 18.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.003 for -l,k,h 0.436 for -h,-k,l 0.006 for l,-k,h	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	12096	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 3.97% 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

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	0.50	3/3091 (0.1%)	0.68	6/4187 (0.1%)
1	B	0.38	0/3091	0.60	2/4187 (0.0%)
1	C	0.40	1/3091 (0.0%)	0.64	2/4187 (0.0%)
1	D	0.43	0/3091	0.63	1/4187 (0.0%)
All	All	0.43	4/12364 (0.0%)	0.64	11/16748 (0.1%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	338	ARG	NE-CZ	10.12	1.46	1.33
1	A	64	LYS	CE-NZ	6.83	1.66	1.49
1	C	213	GLU	CD-OE2	6.31	1.32	1.25
1	A	216	LEU	CG-CD1	5.68	1.72	1.51

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	213	GLU	CA-CB-CG	-12.21	86.53	113.40
1	A	365	LYS	CD-CE-NZ	8.78	131.90	111.70
1	A	216	LEU	CB-CG-CD1	8.55	125.54	111.00
1	D	314	ILE	CG1-CB-CG2	7.80	128.55	111.40
1	A	365	LYS	CG-CD-CE	-6.32	92.95	111.90
1	A	64	LYS	CB-CG-CD	-6.23	95.40	111.60
1	B	161	ARG	NE-CZ-NH1	5.82	123.21	120.30
1	A	216	LEU	CA-CB-CG	-5.80	101.95	115.30
1	A	338	ARG	NE-CZ-NH2	-5.66	117.47	120.30
1	C	213	GLU	OE1-CD-OE2	5.51	129.91	123.30
1	B	161	ARG	NE-CZ-NH2	-5.13	117.74	120.30

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	3024	0	3005	18	0
1	B	3024	0	3005	12	0
1	C	3024	0	3005	16	0
1	D	3024	0	3005	24	0
All	All	12096	0	12020	67	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 3.

All (67) 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:59:LYS:NZ	1:C:213:GLU:OE1	2.10	0.84
1:D:216:LEU:HD13	1:D:216:LEU:O	1.85	0.77
1:D:217:ASP:O	1:D:220:ILE:HG22	1.88	0.73
1:A:58:HIS:NE2	1:C:327:GLU:OE2	2.25	0.69
1:B:15:GLN:HG3	1:B:19:LYS:HE2	1.74	0.69
1:A:62:LEU:HD11	1:A:216:LEU:HD11	1.76	0.67
1:D:313:TRP:O	1:D:314:ILE:HD13	1.95	0.66
1:D:314:ILE:CD1	1:D:362:MET:CE	2.76	0.64
1:C:59:LYS:CD	1:C:213:GLU:OE2	2.48	0.62
1:D:314:ILE:CD1	1:D:362:MET:HE3	2.30	0.61
1:B:264:ILE:HG12	1:B:277:MET:HG3	1.83	0.60
1:D:58:HIS:NE2	1:D:59:LYS:HE2	2.18	0.58
1:D:314:ILE:HD11	1:D:362:MET:SD	2.44	0.58
1:C:59:LYS:HD3	1:C:213:GLU:OE2	2.05	0.57
1:A:15:GLN:HE21	1:A:19:LYS:HE2	1.70	0.56
1:D:59:LYS:HD2	1:D:213:GLU:OE2	2.06	0.56
1:C:59:LYS:HD2	1:C:213:GLU:OE2	2.06	0.55
1:D:217:ASP:HA	1:D:220:ILE:HG22	1.92	0.52
1:D:314:ILE:HD11	1:D:362:MET:CE	2.39	0.52
1:D:314:ILE:HG22	1:D:314:ILE:O	2.10	0.51
1:B:34:THR:HB	1:B:357:TYR:CZ	2.46	0.50
1:A:217:ASP:HA	1:A:220:ILE:HG22	1.94	0.49
1:B:365:LYS:NZ	1:B:376:GLU:HG2	2.28	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:217:ASP:O	1:D:220:ILE:CG2	2.59	0.48
1:D:245:PRO:HB2	1:D:400:LEU:HD13	1.95	0.48
1:D:15:GLN:HG3	1:D:19:LYS:HE2	1.96	0.48
1:D:129:ARG:HH11	1:D:242:VAL:HG11	1.79	0.47
1:D:313:TRP:C	1:D:314:ILE:HD13	2.34	0.47
1:C:365:LYS:HE3	1:C:369:GLU:OE2	2.14	0.47
1:C:34:THR:HB	1:C:357:TYR:CZ	2.51	0.46
1:A:338:ARG:HH22	1:A:361:GLU:HG3	1.81	0.45
1:A:34:THR:HB	1:A:357:TYR:CZ	2.51	0.45
1:A:217:ASP:O	1:A:220:ILE:HG22	2.16	0.45
1:C:365:LYS:NZ	1:C:376:GLU:HG2	2.32	0.45
1:A:64:LYS:CE	1:A:68:ARG:HH12	2.30	0.45
1:C:320:ARG:HD3	1:C:321:ALA:N	2.31	0.45
1:B:137:LEU:HD13	1:B:197:LEU:HD23	1.99	0.45
1:B:62:LEU:HD13	1:B:213:GLU:HG3	2.00	0.44
1:D:34:THR:HB	1:D:357:TYR:CZ	2.52	0.43
1:B:320:ARG:HD3	1:B:321:ALA:N	2.33	0.43
1:A:245:PRO:HB2	1:A:400:LEU:HD13	2.00	0.43
1:B:350:MET:N	1:B:354:CYS:SG	2.91	0.43
1:D:71:LYS:HA	1:D:71:LYS:HD3	1.75	0.43
1:A:277:MET:HG2	1:A:278:SER:O	2.19	0.43
1:A:173:GLY:CA	1:B:147:MET:HE1	2.49	0.42
1:C:62:LEU:HD12	1:C:62:LEU:HA	1.81	0.42
1:D:217:ASP:C	1:D:220:ILE:HG22	2.39	0.42
1:C:198:VAL:O	1:C:231:ALA:HA	2.19	0.42
1:A:58:HIS:HE1	1:C:334:PRO:HB3	1.84	0.42
1:D:198:VAL:O	1:D:231:ALA:HA	2.19	0.42
1:D:245:PRO:HB2	1:D:400:LEU:CD1	2.50	0.42
1:A:62:LEU:CD1	1:A:216:LEU:HD11	2.46	0.42
1:A:281:SER:OG	1:A:282:PRO:HD3	2.20	0.41
1:C:294:VAL:CG1	1:C:298:LYS:HD2	2.49	0.41
1:B:198:VAL:O	1:B:231:ALA:HA	2.20	0.41
1:B:389:GLY:N	1:B:390:MET:HA	2.35	0.41
1:C:338:ARG:HH21	1:C:365:LYS:HD3	1.85	0.41
1:A:198:VAL:O	1:A:231:ALA:HA	2.19	0.41
1:D:281:SER:OG	1:D:282:PRO:HD3	2.21	0.41
1:A:64:LYS:HB2	1:A:64:LYS:HE2	1.89	0.41
1:A:264:ILE:HG12	1:A:277:MET:HG3	2.03	0.41
1:C:172:GLN:HB2	1:C:176:ALA:HB2	2.03	0.41
1:D:367:ALA:HB2	1:D:402:LEU:HD11	2.02	0.41
1:D:161:ARG:NH2	1:D:163:ASN:OD1	2.54	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:338:ARG:HH21	1:B:365:LYS:HD3	1.86	0.40
1:A:338:ARG:HG2	1:A:338:ARG:HH11	1.87	0.40
1:C:294:VAL:HG12	1:C:298:LYS:HD2	2.02	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	388/403 (96%)	379 (98%)	8 (2%)	1 (0%)	41	22
1	B	388/403 (96%)	380 (98%)	7 (2%)	1 (0%)	41	22
1	C	388/403 (96%)	379 (98%)	8 (2%)	1 (0%)	41	22
1	D	388/403 (96%)	380 (98%)	7 (2%)	1 (0%)	41	22
All	All	1552/1612 (96%)	1518 (98%)	30 (2%)	4 (0%)	41	22

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	351	VAL
1	D	351	VAL
1	B	351	VAL
1	C	351	VAL

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	326/338 (96%)	322 (99%)	4 (1%)	71	56
1	B	326/338 (96%)	323 (99%)	3 (1%)	78	67
1	C	326/338 (96%)	323 (99%)	3 (1%)	78	67
1	D	326/338 (96%)	321 (98%)	5 (2%)	65	49
All	All	1304/1352 (96%)	1289 (99%)	15 (1%)	71	56

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

Mol	Chain	Res	Type
1	A	65	LYS
1	A	170	TYR
1	A	320	ARG
1	A	363	ARG
1	B	170	TYR
1	B	320	ARG
1	B	363	ARG
1	C	170	TYR
1	C	320	ARG
1	C	363	ARG
1	D	71	LYS
1	D	170	TYR
1	D	279	LYS
1	D	320	ARG
1	D	363	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 [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 ⓘ

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	390/403 (96%)	0.29	13 (3%) 46 53	16, 24, 40, 50	0
1	B	390/403 (96%)	0.28	7 (1%) 68 76	16, 24, 45, 54	0
1	C	390/403 (96%)	0.26	8 (2%) 63 71	17, 24, 45, 55	0
1	D	390/403 (96%)	0.33	15 (3%) 40 47	16, 24, 44, 57	0
All	All	1560/1612 (96%)	0.29	43 (2%) 53 58	16, 24, 44, 57	0

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	216	LEU	6.4
1	A	216	LEU	5.5
1	B	216	LEU	5.4
1	D	240	ILE	5.3
1	A	220	ILE	5.0
1	C	213	GLU	4.0
1	D	58	HIS	3.9
1	A	218	ASN	3.9
1	B	220	ILE	3.8
1	D	241	GLY	3.8
1	D	370	GLY	3.7
1	D	219	ALA	3.7
1	D	214	SER	3.6
1	C	402	LEU	3.6
1	D	220	ILE	3.3
1	D	217	ASP	3.2
1	C	60	VAL	3.2
1	A	64	LYS	3.2
1	B	402	LEU	3.0
1	D	215	HIS	3.0
1	D	60	VAL	2.9

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	58	HIS	2.9
1	A	219	ALA	2.9
1	C	214	SER	2.8
1	A	213	GLU	2.7
1	D	213	GLU	2.7
1	D	368	ASP	2.6
1	A	240	ILE	2.6
1	C	281	SER	2.6
1	A	242	VAL	2.5
1	D	402	LEU	2.5
1	D	367	ALA	2.3
1	B	214	SER	2.3
1	C	306	GLU	2.3
1	A	277	MET	2.2
1	B	218	ASN	2.2
1	C	216	LEU	2.2
1	A	367	ALA	2.2
1	B	58	HIS	2.1
1	A	370	GLY	2.1
1	C	318	GLY	2.1
1	A	276	TYR	2.0
1	B	62	LEU	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 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.