



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 11:21 AM UTC

PDB ID : 8ES6 / pdb_00008es6
Title : Crystal structure of an unusual amidase ClbL from colibactin gene cluster
Authors : Tripathi, P.; Bruner, S.D.
Deposited on : 2022-10-13
Resolution : 1.90 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 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

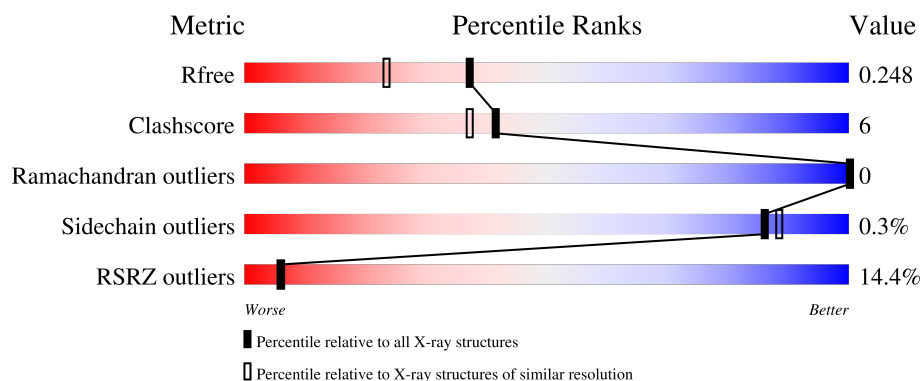
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	488	4% 80% 8% 12%
1	B	488	22% 71% 16% 12%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6969 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 Colibactin biosynthesis amidase ClbL.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	430	Total	C	N	O	S	0	3	0
			3376	2135	599	623	19			
1	B	428	Total	C	N	O	S	0	0	0
			3335	2109	594	615	17			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	488	GLY	-	expression tag	UNP Q0P7K2
B	488	GLY	-	expression tag	UNP Q0P7K2

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	160	Total	O	0	0
			160	160		
2	B	98	Total	O	0	0
			98	98		

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	103.92Å 56.37Å 145.02Å 90.00° 91.84° 90.00°	Depositor
Resolution (Å)	49.55 – 1.90 49.55 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.7 (49.55-1.90) 99.9 (49.55-1.90)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.42 (at 1.90Å)	Xtriage
Refinement program	PHENIX 1.20_4459	Depositor
R, R_{free}	0.206 , 0.249 0.207 , 0.248	Depositor DCC
R_{free} test set	3213 reflections (4.80%)	wwPDB-VP
Wilson B-factor (Å ²)	25.0	Xtriage
Anisotropy	0.701	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 37.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.017 for -h,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6969	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 6.93% 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: SEB

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	0.40	0/3452	0.62	1/4688 (0.0%)
1	B	0.31	0/3412	0.53	0/4635
All	All	0.36	0/6864	0.58	1/9323 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	395	VAL	N-CA-C	-5.02	107.94	112.96

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	3376	0	3290	29	0
1	B	3335	0	3246	56	0
2	A	160	0	0	0	0
2	B	98	0	0	0	0
All	All	6969	0	6536	85	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 6.

All (85) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:421:ASP:OD2	1:A:423:ASN:HB2	1.47	1.10
1:B:205:VAL:HG23	1:B:206:PRO:HD3	1.42	1.00
1:B:33:ARG:HH22	1:B:488:GLY:H	1.24	0.85
1:A:404[B]:MET:HE1	1:A:412:LYS:H	1.42	0.84
1:A:176:LEU:HB3	1:A:179:SEB:HE2	1.60	0.82
1:A:60:ASP:O	1:A:64:LEU:HD13	1.86	0.76
1:B:39:LYS:HG2	1:B:40:PRO:HD3	1.69	0.75
1:B:95:HIS:HA	1:B:98:LYS:HB2	1.68	0.75
1:A:177:PHE:HB2	1:A:179:SEB:HH1	1.75	0.68
1:B:83:PHE:O	1:B:122:LYS:NZ	2.28	0.66
1:A:275:GLN:HG3	1:A:279:GLN:NE2	2.11	0.66
1:B:82:SER:HB2	1:B:205:VAL:HG22	1.76	0.65
1:A:33:ARG:HH22	1:A:488:GLY:H	1.45	0.65
1:A:265:TRP:CG	1:A:300:MET:HG2	2.33	0.64
1:A:18:SER:OG	1:A:20:GLU:HG3	1.99	0.62
1:B:39:LYS:CG	1:B:40:PRO:HD3	2.29	0.62
1:B:368:ARG:HB3	1:B:370:THR:HG22	1.83	0.60
1:B:175:ASP:HA	1:B:179:SEB:HB3	1.82	0.60
1:B:3:GLU:HG2	1:B:29:HIS:HB2	1.85	0.59
1:A:421:ASP:OD2	1:A:423:ASN:CB	2.38	0.58
1:B:92:SER:HB2	1:B:205:VAL:HG21	1.83	0.58
1:B:404:MET:O	1:B:411:ARG:NE	2.24	0.57
1:A:197:GLY:HA2	1:A:242:TRP:CE2	2.39	0.57
1:B:411:ARG:O	1:B:412:LYS:HD2	2.06	0.55
1:A:265:TRP:CD1	1:A:300:MET:HG2	2.42	0.55
1:A:387:ARG:HG2	1:A:388:TYR:CE1	2.42	0.55
1:B:205:VAL:CG2	1:B:206:PRO:HD3	2.27	0.53
1:B:33:ARG:NH2	1:B:488:GLY:H	2.01	0.53
1:B:265:TRP:CG	1:B:300:MET:HG2	2.44	0.53
1:B:37:PHE:O	1:B:40:PRO:HD2	2.08	0.52
1:A:175:ASP:HA	1:A:179:SEB:HB3	1.91	0.52
1:B:97:LEU:HD22	1:B:100:ASN:HD22	1.74	0.52
1:B:368:ARG:HB3	1:B:370:THR:CG2	2.40	0.52
1:B:404:MET:HB3	1:B:411:ARG:HG2	1.91	0.52
1:A:312:ILE:HD11	1:A:373:LEU:HD23	1.93	0.51
1:B:193:ARG:NH2	1:B:433:THR:HA	2.26	0.51
1:A:130:ALA:C	1:A:179:SEB:HI2	2.35	0.51
1:A:275:GLN:HG3	1:A:279:GLN:HE21	1.76	0.51
1:B:135:TYR:HB2	1:B:141:THR:HG22	1.93	0.50
1:B:245:PRO:HB3	1:B:462:PHE:CE2	2.46	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:448:MET:HE1	1:B:485:GLU:CD	2.37	0.49
1:B:128:MET:HE3	1:B:206:PRO:HG3	1.93	0.49
1:B:175:ASP:CG	1:B:193:ARG:HG3	2.37	0.49
1:B:373:LEU:HD22	1:B:377:TYR:CZ	2.48	0.47
1:B:31:PHE:HE1	1:B:47:GLN:HB2	1.78	0.47
1:B:73:HIS:HA	1:B:116:GLY:HA3	1.96	0.47
1:B:92:SER:CB	1:B:205:VAL:HG21	2.45	0.47
1:B:241:PHE:CD1	1:B:241:PHE:N	2.84	0.46
1:A:201:VAL:HG12	1:A:201:VAL:O	2.16	0.46
1:A:235:MET:HE3	1:A:235:MET:HB3	1.79	0.46
1:B:82:SER:O	1:B:204:HIS:HD2	1.98	0.46
1:B:45:VAL:HG23	1:B:46:GLN:HG3	1.98	0.45
1:B:175:ASP:HB2	1:B:180:LEU:HB2	1.98	0.45
1:B:285:ILE:HG23	1:B:296:VAL:HG11	1.97	0.45
1:A:406:GLN:HB2	1:A:407:PRO:HD2	1.98	0.45
1:B:310:PHE:O	1:B:314:VAL:HG23	2.16	0.45
1:B:34:MET:CE	1:B:44:VAL:HG11	2.46	0.45
1:A:411:ARG:O	1:A:412:LYS:HD3	2.17	0.45
1:A:39:LYS:HB2	1:A:40:PRO:HD3	1.99	0.45
1:B:3:GLU:HG3	1:B:7:ARG:NH2	2.32	0.45
1:B:44:VAL:HG22	1:B:123:THR:HG22	2.00	0.44
1:B:181:ARG:HB3	1:B:441:VAL:HG21	2.00	0.43
1:A:31:PHE:CE1	1:A:120:MET:HB3	2.54	0.43
1:B:409:ARG:HA	1:B:409:ARG:HD2	1.75	0.43
1:B:195:SER:OG	1:B:438:PRO:HD3	2.18	0.43
1:A:241:PHE:N	1:A:241:PHE:CD1	2.86	0.43
1:A:379:ARG:O	1:A:383:THR:HG23	2.19	0.43
1:A:308:ALA:O	1:A:312:ILE:HG13	2.19	0.43
1:B:199:MET:HE3	1:B:199:MET:HB2	1.73	0.43
1:B:192:HIS:O	1:B:222:GLY:HA3	2.19	0.43
1:A:75:LEU:HA	1:A:76:PRO:HD3	1.92	0.42
1:B:81:GLU:HA	1:B:122:LYS:HE3	2.01	0.42
1:B:18:SER:OG	1:B:20:GLU:HG3	2.20	0.42
1:B:64:LEU:HD23	1:B:64:LEU:HA	1.89	0.41
1:B:103:THR:HG22	1:B:104:GLN:HG2	2.01	0.41
1:B:464:ASP:OD1	1:B:464:ASP:N	2.52	0.41
1:B:77:CYS:HA	1:B:168:THR:OG1	2.20	0.41
1:B:143:HIS:HB3	1:B:148:ARG:HG2	2.03	0.41
1:B:312:ILE:O	1:B:316:LEU:HD13	2.20	0.41
1:B:91:THR:HG21	1:B:138:LEU:HD13	2.03	0.41
1:B:197:GLY:HA2	1:B:242:TRP:NE1	2.36	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:387:ARG:O	1:A:387:ARG:HG3	2.20	0.40
1:B:234:MET:HE3	1:B:234:MET:HB3	1.99	0.40
1:A:177:PHE:H	1:A:179:SEB:CE	2.33	0.40
1:B:448:MET:HE1	1:B:485:GLU:OE2	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	426/488 (87%)	411 (96%)	15 (4%)	0	100	100
1	B	421/488 (86%)	401 (95%)	20 (5%)	0	100	100
All	All	847/976 (87%)	812 (96%)	35 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	353/392 (90%)	352 (100%)	1 (0%)	86	88
1	B	348/392 (89%)	347 (100%)	1 (0%)	86	88
All	All	701/784 (89%)	699 (100%)	2 (0%)	86	88

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	193	ARG
1	B	39	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	86	GLN
1	A	104	GLN
1	A	279	GLN
1	B	42	ASN
1	B	100	ASN
1	B	276	GLN
1	B	405	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues 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
1	SEB	B	179	1	4,5,17	0.61	0	1,5,23	0.62	0
1	SEB	A	179	1	15,16,17	2.58	2 (13%)	16,21,23	1.80	4 (25%)

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
1	SEB	B	179	1	-	0/2/4/15	-
1	SEB	A	179	1	-	5/10/13/15	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	179	SEB	CE-SD	-7.77	1.70	1.78
1	A	179	SEB	OG-SD	4.89	1.71	1.56

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	179	SEB	OG-CB-CA	-4.75	101.81	107.63
1	A	179	SEB	OD2-SD-CE	3.07	115.48	108.71
1	A	179	SEB	OG-SD-OD2	-2.71	97.69	107.39
1	A	179	SEB	OG-SD-CE	-2.28	97.88	104.18

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	179	SEB	SD-CE-CZ-CH2
1	A	179	SEB	SD-CE-CZ-CH1
1	A	179	SEB	CA-CB-OG-SD
1	A	179	SEB	CB-OG-SD-OD2
1	A	179	SEB	CB-OG-SD-CE

There are no ring outliers.

2 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	B	179	SEB	1	0
1	A	179	SEB	5	0

5.5 Carbohydrates

There are no oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

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	429/488 (87%)	0.19	18 (4%)	40 43	13, 25, 45, 80	3 (0%)
1	B	427/488 (87%)	1.14	105 (24%)	2 1	18, 38, 71, 85	0
All	All	856/976 (87%)	0.67	123 (14%)	6 6	13, 30, 65, 85	3 (0%)

All (123) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	94	ALA	5.9
1	B	216	PRO	5.8
1	B	96	TYR	5.4
1	B	205	VAL	5.2
1	B	97	LEU	5.2
1	B	408	VAL	4.8
1	B	104	GLN	4.5
1	B	242	TRP	4.5
1	A	207	GLY	4.3
1	B	91	THR	4.3
1	B	139	TYR	4.3
1	B	89	LEU	4.1
1	B	203	GLY	4.0
1	B	95	HIS	4.0
1	B	45	VAL	4.0
1	B	370	THR	3.9
1	B	312	ILE	3.9
1	B	316	LEU	3.8
1	A	488	GLY	3.7
1	B	85	VAL	3.7
1	B	99	ASP	3.5
1	B	138	LEU	3.5
1	B	90	THR	3.5
1	B	243	VAL	3.5

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Mol	Chain	Res	Type	RSRZ
1	B	127	MET	3.4
1	B	88	TRP	3.4
1	B	409	ARG	3.3
1	B	135	TYR	3.3
1	B	206	PRO	3.3
1	B	129	THR	3.2
1	B	241	PHE	3.2
1	B	130	ALA	3.2
1	B	218	LEU	3.2
1	B	50	ALA	3.1
1	B	200	SER	3.1
1	B	92	SER	3.1
1	B	313	ALA	3.1
1	B	373	LEU	3.1
1	B	103	THR	3.1
1	B	93	GLY	3.0
1	B	106	ALA	3.0
1	B	406	GLN	3.0
1	B	314	VAL	3.0
1	B	368	ARG	3.0
1	B	176	LEU	2.9
1	A	407	PRO	2.9
1	B	197	GLY	2.9
1	B	375	HIS	2.9
1	B	44	VAL	2.9
1	B	143	HIS	2.9
1	B	49	TYR	2.8
1	B	102	ALA	2.8
1	B	310	PHE	2.8
1	A	242	TRP	2.8
1	B	125	VAL	2.8
1	B	407	PRO	2.8
1	B	84	ASP	2.8
1	B	371	ASP	2.8
1	B	126	PRO	2.8
1	A	2	SER	2.7
1	B	109	ILE	2.7
1	B	46	GLN	2.7
1	B	51	LEU	2.7
1	B	304	ILE	2.7
1	B	311	ASP	2.6
1	B	119	LEU	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	83	PHE	2.6
1	B	81	GLU	2.6
1	B	207	GLY	2.6
1	B	140	GLY	2.5
1	B	128	MET	2.5
1	B	57	ALA	2.5
1	B	378	SER	2.4
1	B	98	LYS	2.4
1	B	411	ARG	2.4
1	B	245	PRO	2.4
1	B	405	GLN	2.4
1	B	309	LEU	2.4
1	B	100	ASN	2.4
1	B	434	VAL	2.4
1	B	101	ARG	2.3
1	A	96	TYR	2.3
1	A	241	PHE	2.3
1	B	132	TRP	2.3
1	B	202	ARG	2.3
1	B	196	LEU	2.3
1	B	42	ASN	2.3
1	B	136	ASN	2.3
1	B	105	ASP	2.3
1	B	141	THR	2.3
1	A	216	PRO	2.2
1	B	422	TYR	2.2
1	B	48	HIS	2.2
1	B	244	GLU	2.2
1	B	488	GLY	2.2
1	A	302	ALA	2.2
1	A	408	VAL	2.2
1	B	41	LEU	2.2
1	B	217	ASP	2.2
1	B	317	SER	2.2
1	B	107	PRO	2.2
1	B	315	LYS	2.2
1	A	261	ARG	2.2
1	B	86	GLN	2.2
1	B	134	THR	2.2
1	A	369	ASN	2.1
1	B	423	ASN	2.1
1	B	118	ILE	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	4	GLN	2.1
1	A	103	THR	2.1
1	B	82	SER	2.1
1	B	369	ASN	2.1
1	A	49	TYR	2.1
1	B	155	SER	2.1
1	A	316	LEU	2.1
1	B	198	LEU	2.1
1	B	110	ALA	2.1
1	B	379	ARG	2.1
1	A	372	LYS	2.0
1	B	124	ASN	2.0
1	B	3	GLU	2.0
1	A	448	MET	2.0
1	A	409	ARG	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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
1	SEB	A	179	16/17	0.85	0.16	14,30,44,47	16
1	SEB	B	179	6/17	0.85	0.11	37,39,44,51	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides 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.