



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 06:31 AM UTC

PDB ID : 3BSM / pdb_00003bsm
Title : Crystal structure of D-mannonate dehydratase from *Chromohalobacter salexigens*
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Deposited on : 2007-12-25
Resolution : 2.20 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
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

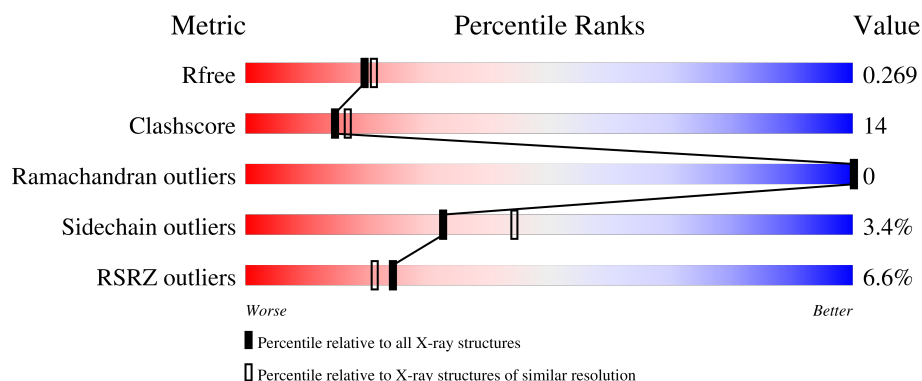
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	413	5% 63% 21% • 13%
1	B	413	5% 65% 19% • 13%
1	C	413	7% 63% 21% • 13%
1	D	413	7% 62% 22% • 13%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 11512 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 Mandelate racemase/muconate lactonizing enzyme.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	358	Total	C	N	O	S	0	0	0
			2825	1789	505	517	14			
1	B	358	Total	C	N	O	S	0	0	0
			2825	1789	505	517	14			
1	C	358	Total	C	N	O	S	0	0	0
			2825	1789	505	517	14			
1	D	358	Total	C	N	O	S	0	0	0
			2825	1789	505	517	14			

There are 44 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	expression tag	UNP Q1QT89
A	2	SER	-	expression tag	UNP Q1QT89
A	3	LEU	-	expression tag	UNP Q1QT89
A	406	GLU	-	expression tag	UNP Q1QT89
A	407	GLY	-	expression tag	UNP Q1QT89
A	408	HIS	-	expression tag	UNP Q1QT89
A	409	HIS	-	expression tag	UNP Q1QT89
A	410	HIS	-	expression tag	UNP Q1QT89
A	411	HIS	-	expression tag	UNP Q1QT89
A	412	HIS	-	expression tag	UNP Q1QT89
A	413	HIS	-	expression tag	UNP Q1QT89
B	1	MET	-	expression tag	UNP Q1QT89
B	2	SER	-	expression tag	UNP Q1QT89
B	3	LEU	-	expression tag	UNP Q1QT89
B	406	GLU	-	expression tag	UNP Q1QT89
B	407	GLY	-	expression tag	UNP Q1QT89
B	408	HIS	-	expression tag	UNP Q1QT89
B	409	HIS	-	expression tag	UNP Q1QT89
B	410	HIS	-	expression tag	UNP Q1QT89
B	411	HIS	-	expression tag	UNP Q1QT89
B	412	HIS	-	expression tag	UNP Q1QT89

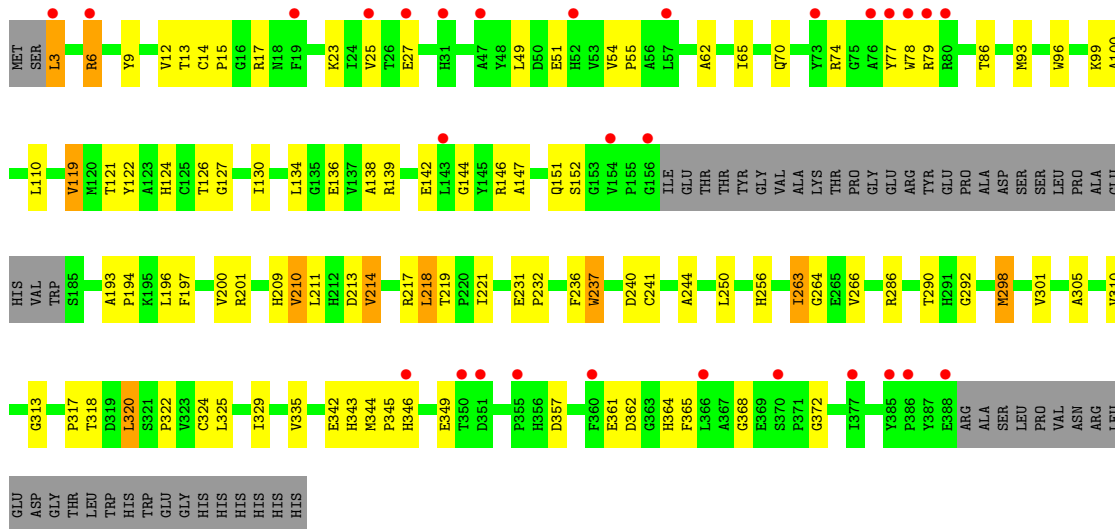
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Chain	Residue	Modelled	Actual	Comment	Reference
B	413	HIS	-	expression tag	UNP Q1QT89
C	1	MET	-	expression tag	UNP Q1QT89
C	2	SER	-	expression tag	UNP Q1QT89
C	3	LEU	-	expression tag	UNP Q1QT89
C	406	GLU	-	expression tag	UNP Q1QT89
C	407	GLY	-	expression tag	UNP Q1QT89
C	408	HIS	-	expression tag	UNP Q1QT89
C	409	HIS	-	expression tag	UNP Q1QT89
C	410	HIS	-	expression tag	UNP Q1QT89
C	411	HIS	-	expression tag	UNP Q1QT89
C	412	HIS	-	expression tag	UNP Q1QT89
C	413	HIS	-	expression tag	UNP Q1QT89
D	1	MET	-	expression tag	UNP Q1QT89
D	2	SER	-	expression tag	UNP Q1QT89
D	3	LEU	-	expression tag	UNP Q1QT89
D	406	GLU	-	expression tag	UNP Q1QT89
D	407	GLY	-	expression tag	UNP Q1QT89
D	408	HIS	-	expression tag	UNP Q1QT89
D	409	HIS	-	expression tag	UNP Q1QT89
D	410	HIS	-	expression tag	UNP Q1QT89
D	411	HIS	-	expression tag	UNP Q1QT89
D	412	HIS	-	expression tag	UNP Q1QT89
D	413	HIS	-	expression tag	UNP Q1QT89

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	67	Total O 67 67	0	0
2	B	60	Total O 60 60	0	0
2	C	46	Total O 46 46	0	0
2	D	39	Total O 39 39	0	0



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	110.12Å 163.07Å 84.55Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	24.93 – 2.20 24.93 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.5 (24.93-2.20) 99.4 (24.93-2.20)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.42 (at 2.00Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.243 , 0.268 0.243 , 0.269	Depositor DCC
R_{free} test set	5073 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	34.7	Xtriage
Anisotropy	0.600	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 46.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	11512	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.33% 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	0.45	1/2900 (0.0%)	0.92	5/3945 (0.1%)
1	B	0.45	1/2900 (0.0%)	0.92	4/3945 (0.1%)
1	C	0.44	1/2900 (0.0%)	0.93	4/3945 (0.1%)
1	D	0.43	1/2900 (0.0%)	0.92	5/3945 (0.1%)
All	All	0.44	4/11600 (0.0%)	0.92	18/15780 (0.1%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	298	MET	SD-CE	-6.70	1.62	1.79
1	A	298	MET	SD-CE	-6.12	1.64	1.79
1	C	298	MET	SD-CE	-5.51	1.65	1.79
1	B	298	MET	SD-CE	-5.46	1.65	1.79

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	219	THR	N-CA-C	-7.24	100.67	110.36
1	B	219	THR	N-CA-C	-7.06	100.90	110.36
1	D	219	THR	N-CA-C	-6.95	101.05	110.36
1	A	219	THR	N-CA-C	-6.82	101.23	110.36
1	D	236	PHE	N-CA-C	-6.67	104.09	111.36
1	C	236	PHE	N-CA-C	-6.59	104.09	111.28
1	A	236	PHE	N-CA-C	-6.49	104.21	111.28
1	B	236	PHE	N-CA-C	-6.38	104.33	111.28
1	B	244	ALA	N-CA-C	5.75	119.84	112.26
1	D	244	ALA	N-CA-C	5.71	119.79	112.26
1	A	218	LEU	N-CA-C	5.61	118.84	110.14
1	C	244	ALA	N-CA-C	5.61	119.67	112.26
1	A	244	ALA	N-CA-C	5.58	119.62	112.26
1	C	218	LEU	N-CA-C	5.25	118.28	110.14
1	B	218	LEU	N-CA-C	5.20	118.19	110.14
1	D	51	GLU	N-CA-C	5.04	118.69	112.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	51	GLU	N-CA-C	5.04	118.68	112.54
1	D	218	LEU	N-CA-C	5.01	117.90	110.14

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	2825	0	2733	83	1
1	B	2825	0	2733	75	0
1	C	2825	0	2733	80	0
1	D	2825	0	2733	82	1
2	A	67	0	0	6	1
2	B	60	0	0	2	0
2	C	46	0	0	2	0
2	D	39	0	0	1	1
All	All	11512	0	10932	308	2

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 14.

All (308) 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:292:GLY:HA3	1:C:298:MET:HE1	1.51	0.93
1:B:292:GLY:HA3	1:B:298:MET:HE1	1.53	0.91
1:D:292:GLY:HA3	1:D:298:MET:HE1	1.53	0.91
1:A:265:GLU:OE1	2:A:480:HOH:O	1.89	0.89
1:A:292:GLY:HA3	1:A:298:MET:HE1	1.55	0.89
1:D:93:MET:HE1	1:D:290:THR:HG22	1.56	0.88
1:C:151:GLN:HE22	1:C:213:ASP:H	1.22	0.88
1:D:151:GLN:HE22	1:D:213:ASP:H	1.22	0.87
1:C:93:MET:HE1	1:C:290:THR:HG22	1.55	0.87
1:B:93:MET:HE1	1:B:290:THR:HG22	1.56	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:151:GLN:HE22	1:B:213:ASP:H	1.23	0.86
1:A:151:GLN:HE22	1:A:213:ASP:H	1.20	0.85
1:A:93:MET:HE1	1:A:290:THR:HG22	1.59	0.84
1:A:99:LYS:HG2	1:A:110:LEU:HD21	1.65	0.79
1:B:99:LYS:HG2	1:B:110:LEU:HD21	1.64	0.77
1:D:99:LYS:HG2	1:D:110:LEU:HD21	1.65	0.77
1:C:99:LYS:HG2	1:C:110:LEU:HD21	1.64	0.76
1:C:119:VAL:HG22	1:C:365:PHE:HB2	1.67	0.76
1:B:119:VAL:HG22	1:B:365:PHE:HB2	1.68	0.76
1:D:119:VAL:HG22	1:D:365:PHE:HB2	1.67	0.75
1:A:119:VAL:HG22	1:A:365:PHE:HB2	1.69	0.74
1:A:266:VAL:HG13	1:C:79:ARG:HD3	1.72	0.72
1:D:197:PHE:O	1:D:200:VAL:HG22	1.91	0.71
1:A:197:PHE:O	1:A:200:VAL:HG22	1.91	0.70
1:B:290:THR:HG23	1:B:324:CYS:SG	2.30	0.70
1:D:298:MET:HA	1:D:301:VAL:HG22	1.74	0.70
1:A:317:PRO:HD2	1:A:320:LEU:HD22	1.74	0.70
1:C:317:PRO:HD2	1:C:320:LEU:HD22	1.74	0.69
1:B:317:PRO:HD2	1:B:320:LEU:HD22	1.74	0.69
1:B:266:VAL:HG13	1:D:79:ARG:HD3	1.74	0.69
1:C:197:PHE:O	1:C:200:VAL:HG22	1.92	0.69
1:D:263:ILE:HD13	1:D:264:GLY:H	1.57	0.69
1:D:317:PRO:HA	1:D:344:MET:HE2	1.76	0.68
1:B:197:PHE:O	1:B:200:VAL:HG22	1.93	0.68
1:C:290:THR:HG23	1:C:324:CYS:SG	2.35	0.67
1:D:318:THR:HG23	1:D:344:MET:HE3	1.74	0.67
1:D:317:PRO:HD2	1:D:320:LEU:HD22	1.75	0.67
1:B:263:ILE:HD13	1:B:264:GLY:H	1.58	0.67
1:C:263:ILE:HD13	1:C:264:GLY:H	1.59	0.66
1:B:318:THR:HG23	1:B:344:MET:HE3	1.77	0.66
1:B:317:PRO:HA	1:B:344:MET:HE2	1.78	0.66
1:A:263:ILE:HD13	1:A:264:GLY:H	1.61	0.65
1:A:290:THR:HG23	1:A:324:CYS:SG	2.36	0.65
1:A:298:MET:HA	1:A:301:VAL:HG22	1.77	0.65
1:D:100:ALA:HB3	1:D:372:GLY:HA2	1.77	0.65
1:D:290:THR:HG23	1:D:324:CYS:SG	2.37	0.65
1:A:138:ALA:O	1:A:142:GLU:HG2	1.97	0.65
1:C:298:MET:HA	1:C:301:VAL:HG22	1.77	0.65
1:C:317:PRO:HA	1:C:344:MET:HE2	1.78	0.65
1:C:138:ALA:O	1:C:142:GLU:HG2	1.97	0.64
1:C:100:ALA:HB3	1:C:372:GLY:HA2	1.79	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:100:ALA:HB3	1:B:372:GLY:HA2	1.77	0.64
1:D:138:ALA:O	1:D:142:GLU:HG2	1.97	0.64
1:A:100:ALA:HB3	1:A:372:GLY:HA2	1.78	0.64
1:A:317:PRO:HA	1:A:344:MET:HE2	1.80	0.63
1:B:138:ALA:O	1:B:142:GLU:HG2	1.97	0.63
1:C:271:HIS:HD2	2:C:441:HOH:O	1.80	0.63
1:D:74:ARG:HG3	1:D:78:TRP:CZ2	2.34	0.63
1:D:193:ALA:HB3	1:D:194:PRO:HD3	1.80	0.63
1:A:318:THR:HG23	1:A:344:MET:HE3	1.81	0.62
1:B:79:ARG:HD3	1:D:266:VAL:HG13	1.82	0.62
1:D:217:ARG:HG3	1:D:217:ARG:HH11	1.65	0.62
1:B:74:ARG:HG3	1:B:78:TRP:CZ2	2.35	0.62
1:A:79:ARG:HD3	1:C:266:VAL:HG13	1.81	0.61
1:A:217:ARG:HG3	1:A:217:ARG:HH11	1.64	0.61
1:B:298:MET:HA	1:B:301:VAL:HG22	1.80	0.61
1:A:193:ALA:HB3	1:A:194:PRO:HD3	1.83	0.61
1:A:74:ARG:HG3	1:A:78:TRP:CZ2	2.35	0.61
1:B:193:ALA:HB3	1:B:194:PRO:HD3	1.83	0.61
1:B:217:ARG:HG3	1:B:217:ARG:HH11	1.65	0.61
1:A:121:THR:CG2	1:A:122:TYR:N	2.64	0.61
1:C:74:ARG:HG3	1:C:78:TRP:CZ2	2.35	0.61
1:C:318:THR:HG23	1:C:344:MET:HE3	1.82	0.60
1:C:121:THR:CG2	1:C:122:TYR:N	2.64	0.60
1:A:241:CYS:SG	2:A:446:HOH:O	2.56	0.60
1:C:292:GLY:CA	1:C:298:MET:HE1	2.29	0.60
1:C:193:ALA:HB3	1:C:194:PRO:HD3	1.82	0.60
1:C:217:ARG:HG3	1:C:217:ARG:HH11	1.66	0.59
1:A:121:THR:HG22	1:A:122:TYR:N	2.17	0.59
1:B:292:GLY:CA	1:B:298:MET:HE1	2.31	0.59
1:D:292:GLY:CA	1:D:298:MET:HE1	2.30	0.59
1:D:121:THR:CG2	1:D:122:TYR:N	2.66	0.59
1:A:292:GLY:CA	1:A:298:MET:HE1	2.31	0.59
1:B:121:THR:CG2	1:B:122:TYR:N	2.66	0.58
1:C:121:THR:HG22	1:C:122:TYR:N	2.18	0.58
1:D:121:THR:HG22	1:D:122:TYR:N	2.19	0.58
1:D:256:HIS:HD2	2:D:417:HOH:O	1.87	0.58
1:A:325:LEU:O	1:A:329:ILE:HG13	2.04	0.57
1:B:13:THR:OG1	1:B:15:PRO:HD3	2.04	0.57
1:B:374:GLY:HA3	2:B:436:HOH:O	2.05	0.57
1:B:121:THR:HG22	1:B:122:TYR:N	2.20	0.56
1:C:325:LEU:O	1:C:329:ILE:HG13	2.06	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:3:LEU:HD13	1:C:3:LEU:N	2.21	0.56
1:A:70:GLN:O	1:A:74:ARG:HD3	2.06	0.56
1:D:325:LEU:O	1:D:329:ILE:HG13	2.05	0.56
1:A:266:VAL:HG21	1:C:77:TYR:HB3	1.88	0.55
1:B:292:GLY:HA3	1:B:298:MET:CE	2.32	0.55
1:D:74:ARG:HG3	1:D:78:TRP:CH2	2.41	0.55
1:B:74:ARG:HG3	1:B:78:TRP:CH2	2.42	0.55
1:A:61:ASP:OD2	2:A:441:HOH:O	2.18	0.55
1:D:70:GLN:O	1:D:74:ARG:HD3	2.06	0.55
1:A:74:ARG:HG3	1:A:78:TRP:CH2	2.42	0.55
1:C:19:PHE:O	2:C:456:HOH:O	2.18	0.55
1:A:318:THR:H	1:A:344:MET:HE2	1.72	0.54
1:C:74:ARG:HG3	1:C:78:TRP:CH2	2.42	0.54
1:C:292:GLY:HA3	1:C:298:MET:CE	2.32	0.54
1:D:119:VAL:CG2	1:D:365:PHE:HB2	2.37	0.54
1:C:9:TYR:OH	1:C:23:LYS:HD3	2.08	0.54
1:B:119:VAL:CG2	1:B:365:PHE:HB2	2.37	0.54
1:A:13:THR:OG1	1:A:15:PRO:HD3	2.07	0.54
1:C:70:GLN:O	1:C:74:ARG:HD3	2.08	0.54
1:C:13:THR:OG1	1:C:15:PRO:HD3	2.07	0.54
1:C:49:LEU:O	1:C:54:VAL:HG23	2.08	0.54
1:D:292:GLY:HA3	1:D:298:MET:CE	2.33	0.54
1:B:49:LEU:O	1:B:54:VAL:HG23	2.08	0.54
1:A:3:LEU:N	1:A:3:LEU:HD13	2.23	0.53
1:B:70:GLN:O	1:B:74:ARG:HD3	2.08	0.53
1:A:11:ILE:O	2:A:442:HOH:O	2.19	0.53
1:B:325:LEU:O	1:B:329:ILE:HG13	2.09	0.53
1:C:119:VAL:CG2	1:C:365:PHE:HB2	2.36	0.53
1:C:298:MET:O	1:C:301:VAL:HG22	2.09	0.53
1:B:139:ARG:O	1:B:142:GLU:HB2	2.09	0.53
1:B:298:MET:O	1:B:301:VAL:HG22	2.09	0.53
1:D:13:THR:OG1	1:D:15:PRO:HD3	2.09	0.53
1:B:318:THR:H	1:B:344:MET:HE2	1.74	0.52
1:D:263:ILE:CD1	1:D:264:GLY:H	2.23	0.52
1:C:286:ARG:HD2	1:C:313:GLY:O	2.09	0.52
1:C:298:MET:SD	1:C:301:VAL:HG21	2.50	0.52
1:C:318:THR:H	1:C:344:MET:HE2	1.74	0.52
1:D:318:THR:H	1:D:344:MET:HE2	1.74	0.52
1:A:286:ARG:HD2	1:A:313:GLY:O	2.10	0.52
1:A:298:MET:SD	1:A:301:VAL:HG21	2.50	0.52
1:B:3:LEU:HD13	1:B:3:LEU:N	2.24	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:9:TYR:OH	1:D:23:LYS:HD3	2.10	0.52
1:D:49:LEU:O	1:D:54:VAL:HG23	2.09	0.52
1:D:3:LEU:N	1:D:3:LEU:HD13	2.25	0.52
1:D:139:ARG:O	1:D:142:GLU:HB2	2.09	0.52
1:C:139:ARG:O	1:C:142:GLU:HB2	2.11	0.51
1:D:298:MET:SD	1:D:301:VAL:HG21	2.50	0.51
1:B:266:VAL:HG21	1:D:77:TYR:HB3	1.92	0.51
1:B:298:MET:SD	1:B:301:VAL:HG21	2.50	0.51
1:D:286:ARG:HD2	1:D:313:GLY:O	2.11	0.51
1:D:237:TRP:CD1	1:D:237:TRP:C	2.89	0.51
1:C:237:TRP:CD1	1:C:237:TRP:C	2.89	0.51
1:B:237:TRP:CD1	1:B:237:TRP:C	2.89	0.51
1:C:211:LEU:HD12	1:C:211:LEU:N	2.26	0.51
1:A:237:TRP:CD1	1:A:237:TRP:C	2.88	0.51
1:C:318:THR:N	1:C:344:MET:HE2	2.26	0.51
1:D:54:VAL:HB	1:D:55:PRO:HD3	1.93	0.51
1:D:211:LEU:N	1:D:211:LEU:HD12	2.26	0.51
1:A:147:ALA:CB	1:A:209:HIS:HB2	2.41	0.50
1:D:196:LEU:O	1:D:200:VAL:HG13	2.11	0.50
1:A:292:GLY:HA3	1:A:298:MET:CE	2.35	0.50
1:B:211:LEU:N	1:B:211:LEU:HD12	2.25	0.50
1:B:54:VAL:HB	1:B:55:PRO:HD3	1.93	0.50
1:A:263:ILE:CD1	1:A:264:GLY:H	2.25	0.50
1:C:54:VAL:HB	1:C:55:PRO:HD3	1.94	0.50
1:D:298:MET:O	1:D:301:VAL:HG22	2.12	0.50
1:C:196:LEU:O	1:C:200:VAL:HG13	2.12	0.50
1:D:151:GLN:NE2	1:D:213:ASP:H	2.02	0.50
1:D:318:THR:N	1:D:344:MET:HE2	2.27	0.50
1:B:9:TYR:OH	1:B:23:LYS:HD3	2.12	0.49
1:A:49:LEU:O	1:A:54:VAL:HG23	2.12	0.49
1:A:318:THR:N	1:A:344:MET:HE2	2.26	0.49
1:A:77:TYR:HB3	1:C:266:VAL:HG21	1.93	0.49
1:A:211:LEU:HD12	1:A:211:LEU:N	2.27	0.49
1:B:286:ARG:HD2	1:B:313:GLY:O	2.12	0.49
1:A:54:VAL:HB	1:A:55:PRO:HD3	1.94	0.49
1:B:196:LEU:O	1:B:200:VAL:HG13	2.13	0.49
1:A:196:LEU:O	1:A:200:VAL:HG13	2.11	0.49
1:A:119:VAL:CG2	1:A:365:PHE:HB2	2.40	0.49
1:B:214:VAL:HG23	1:B:214:VAL:O	2.11	0.49
1:B:263:ILE:CD1	1:B:264:GLY:H	2.26	0.49
1:D:214:VAL:HG23	1:D:214:VAL:O	2.11	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:214:VAL:O	1:A:214:VAL:HG23	2.12	0.48
1:A:298:MET:O	1:A:301:VAL:HG22	2.13	0.48
1:C:147:ALA:CB	1:C:209:HIS:HB2	2.43	0.48
1:C:263:ILE:CD1	1:C:264:GLY:H	2.25	0.48
1:B:318:THR:N	1:B:344:MET:HE2	2.28	0.48
1:A:139:ARG:O	1:A:142:GLU:HB2	2.13	0.48
1:B:147:ALA:CB	1:B:209:HIS:HB2	2.44	0.48
1:D:317:PRO:CA	1:D:344:MET:HE2	2.42	0.48
1:A:214:VAL:HG22	1:A:240:ASP:O	2.13	0.48
1:C:214:VAL:HG23	1:C:214:VAL:O	2.12	0.48
1:D:147:ALA:CB	1:D:209:HIS:HB2	2.43	0.48
1:C:362:ASP:O	1:C:364:HIS:HD2	1.97	0.48
1:D:6:ARG:HG3	1:D:25:VAL:O	2.14	0.48
1:A:362:ASP:O	1:A:364:HIS:HD2	1.95	0.48
1:B:214:VAL:HG22	1:B:240:ASP:O	2.14	0.48
1:B:362:ASP:O	1:B:364:HIS:HD2	1.96	0.48
1:C:136:GLU:OE2	1:C:139:ARG:HD3	2.14	0.48
1:C:119:VAL:O	1:C:119:VAL:HG23	2.13	0.47
1:A:305:ALA:HB1	1:A:310:VAL:HB	1.95	0.47
1:D:362:ASP:O	1:D:364:HIS:HD2	1.97	0.47
1:C:127:GLY:O	1:C:152:SER:HA	2.15	0.47
1:D:214:VAL:HG22	1:D:240:ASP:O	2.14	0.47
1:D:298:MET:CA	1:D:301:VAL:HG22	2.43	0.47
1:A:9:TYR:OH	1:A:23:LYS:HD3	2.14	0.47
1:A:136:GLU:OE2	1:A:139:ARG:HD3	2.15	0.47
1:A:147:ALA:HB2	1:A:209:HIS:HB2	1.96	0.47
1:C:6:ARG:HG3	1:C:25:VAL:O	2.14	0.47
1:D:136:GLU:OE2	1:D:139:ARG:HD3	2.15	0.47
1:B:119:VAL:O	1:B:119:VAL:HG23	2.14	0.47
1:D:201:ARG:NH2	1:D:210:VAL:HG13	2.30	0.47
1:D:119:VAL:HG23	1:D:119:VAL:O	2.13	0.46
1:D:357:ASP:CG	1:D:368:GLY:HA3	2.40	0.46
1:C:322:PRO:HG3	1:C:346:HIS:CD2	2.51	0.46
1:B:136:GLU:OE2	1:B:139:ARG:HD3	2.15	0.46
1:B:266:VAL:CG1	1:D:79:ARG:HD3	2.45	0.46
1:C:214:VAL:HG22	1:C:240:ASP:O	2.15	0.46
1:A:119:VAL:HG23	1:A:119:VAL:O	2.13	0.46
1:A:127:GLY:O	1:A:152:SER:HA	2.15	0.46
1:A:219:THR:OG1	1:A:222:GLU:HG3	2.16	0.46
1:B:77:TYR:HB3	1:D:266:VAL:HG21	1.97	0.46
1:A:266:VAL:CG1	1:C:79:ARG:HD3	2.43	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:322:PRO:HG3	1:D:346:HIS:CD2	2.51	0.46
1:A:201:ARG:NH2	1:A:210:VAL:HG13	2.31	0.46
1:B:201:ARG:NH2	1:B:210:VAL:HG13	2.31	0.46
1:C:317:PRO:CA	1:C:344:MET:HE2	2.44	0.46
1:D:147:ALA:HB2	1:D:209:HIS:HB2	1.98	0.46
1:A:322:PRO:HG3	1:A:346:HIS:CD2	2.51	0.46
1:B:127:GLY:O	1:B:152:SER:HA	2.15	0.46
1:D:292:GLY:O	1:D:298:MET:HE1	2.16	0.46
1:A:6:ARG:HG3	1:A:25:VAL:O	2.16	0.45
1:D:349:GLU:CD	1:D:349:GLU:H	2.25	0.45
1:C:219:THR:OG1	1:C:222:GLU:HG3	2.16	0.45
1:D:305:ALA:HB1	1:D:310:VAL:HB	1.98	0.45
1:A:357:ASP:CG	1:A:368:GLY:HA3	2.42	0.45
1:D:127:GLY:O	1:D:152:SER:HA	2.16	0.45
1:B:322:PRO:HG3	1:B:346:HIS:CD2	2.52	0.45
1:B:317:PRO:CA	1:B:344:MET:HE2	2.44	0.45
1:C:357:ASP:CG	1:C:368:GLY:HA3	2.42	0.45
1:A:298:MET:CA	1:A:301:VAL:HG22	2.46	0.44
1:A:346:HIS:HD2	2:A:456:HOH:O	1.99	0.44
1:A:292:GLY:O	1:A:298:MET:HE1	2.18	0.44
1:A:317:PRO:CA	1:A:344:MET:HE2	2.45	0.44
1:B:357:ASP:CG	1:B:368:GLY:HA3	2.43	0.44
1:C:147:ALA:HB2	1:C:209:HIS:HB2	1.99	0.44
1:C:298:MET:CA	1:C:301:VAL:HG22	2.46	0.44
1:C:119:VAL:CG2	1:C:119:VAL:O	2.66	0.44
1:C:305:ALA:HB1	1:C:310:VAL:HB	1.99	0.44
1:C:130:ILE:O	1:C:134:LEU:HG	2.18	0.44
1:B:6:ARG:HG3	1:B:25:VAL:O	2.17	0.44
1:C:201:ARG:NH2	1:C:210:VAL:HG13	2.33	0.43
1:B:124:HIS:HE1	1:B:342:GLU:OE2	2.01	0.43
1:C:144:GLY:O	1:C:146:ARG:NH1	2.51	0.43
1:A:349:GLU:CD	1:A:349:GLU:H	2.27	0.43
1:D:86:THR:HA	1:D:290:THR:O	2.19	0.43
1:D:119:VAL:CG2	1:D:119:VAL:O	2.66	0.43
1:C:214:VAL:CG2	1:C:241:CYS:HA	2.49	0.43
1:D:130:ILE:O	1:D:134:LEU:HG	2.19	0.43
1:B:349:GLU:CD	1:B:349:GLU:H	2.27	0.43
1:B:130:ILE:O	1:B:134:LEU:HG	2.19	0.43
1:B:147:ALA:HB2	1:B:209:HIS:HB2	2.00	0.43
1:A:119:VAL:CG2	1:A:119:VAL:O	2.66	0.42
1:B:305:ALA:HB1	1:B:310:VAL:HB	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:144:GLY:O	1:D:146:ARG:NH1	2.52	0.42
1:B:3:LEU:HA	1:B:27:GLU:OE2	2.20	0.42
1:B:119:VAL:CG2	1:B:119:VAL:O	2.67	0.42
1:C:12:VAL:O	1:C:12:VAL:HG13	2.19	0.42
1:B:374:GLY:CA	2:B:436:HOH:O	2.67	0.42
1:C:124:HIS:HE1	1:C:342:GLU:OE2	2.03	0.42
1:A:86:THR:HA	1:A:290:THR:O	2.20	0.42
1:C:86:THR:HA	1:C:290:THR:O	2.20	0.42
1:C:349:GLU:CD	1:C:349:GLU:H	2.27	0.42
1:A:214:VAL:CG2	1:A:241:CYS:HA	2.50	0.42
1:C:292:GLY:O	1:C:298:MET:HE1	2.20	0.42
1:D:121:THR:CG2	1:D:343:HIS:HB3	2.50	0.42
1:B:86:THR:HA	1:B:290:THR:O	2.19	0.42
1:D:214:VAL:CG2	1:D:241:CYS:HA	2.50	0.42
1:C:65:ILE:HG21	1:C:96:TRP:NE1	2.35	0.42
1:A:79:ARG:HD3	1:C:266:VAL:HG22	2.02	0.41
1:A:130:ILE:O	1:A:134:LEU:HG	2.20	0.41
1:B:214:VAL:CG2	1:B:241:CYS:HA	2.50	0.41
1:D:126:THR:HG22	1:D:151:GLN:HB2	2.02	0.41
1:D:12:VAL:HG13	1:D:12:VAL:O	2.20	0.41
1:D:3:LEU:HA	1:D:27:GLU:OE2	2.20	0.41
1:A:124:HIS:HE1	1:A:342:GLU:OE2	2.04	0.41
1:C:126:THR:HG22	1:C:151:GLN:HB2	2.02	0.41
1:C:231:GLU:N	1:C:232:PRO:CD	2.84	0.41
1:D:231:GLU:N	1:D:232:PRO:CD	2.84	0.41
1:B:79:ARG:HD3	1:D:266:VAL:CG1	2.49	0.41
1:D:3:LEU:HD23	1:D:62:ALA:HB3	2.02	0.41
1:D:221:ILE:HD12	1:D:221:ILE:HA	1.96	0.41
1:A:12:VAL:HG13	1:A:12:VAL:O	2.21	0.41
1:A:49:LEU:HA	1:A:53:VAL:HB	2.03	0.41
1:B:200:VAL:HG23	1:B:201:ARG:N	2.36	0.41
1:C:3:LEU:HA	1:C:27:GLU:OE2	2.21	0.41
1:A:3:LEU:HA	1:A:27:GLU:OE2	2.20	0.41
1:A:144:GLY:O	1:A:146:ARG:NH1	2.53	0.41
1:A:221:ILE:HD12	1:A:221:ILE:HA	1.95	0.41
1:B:144:GLY:O	1:B:146:ARG:NH1	2.53	0.41
1:B:292:GLY:O	1:B:298:MET:HE1	2.21	0.41
1:A:65:ILE:HG21	1:A:96:TRP:NE1	2.35	0.40
1:B:65:ILE:HG21	1:B:96:TRP:NE1	2.36	0.40
1:B:219:THR:OG1	1:B:222:GLU:HG3	2.21	0.40
1:C:3:LEU:HD23	1:C:62:ALA:HB3	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:14:CYS:HA	1:D:17:ARG:O	2.21	0.40
1:D:65:ILE:HG21	1:D:96:TRP:NE1	2.35	0.40
1:D:343:HIS:NE2	1:D:345:PRO:HG3	2.37	0.40
1:A:300:ARG:NH1	2:A:450:HOH:O	2.41	0.40
1:C:335:VAL:O	1:C:335:VAL:HG13	2.20	0.40
1:D:124:HIS:HE1	1:D:342:GLU:OE2	2.04	0.40
1:A:14:CYS:HA	1:A:17:ARG:O	2.21	0.40
1:A:217:ARG:HG3	1:A:217:ARG:NH1	2.35	0.40
1:B:14:CYS:O	1:B:14:CYS:SG	2.80	0.40
1:C:14:CYS:HA	1:C:17:ARG:O	2.22	0.40
1:D:335:VAL:O	1:D:335:VAL:HG13	2.21	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:385:TYR:OH	1:D:361:GLU:OE2[3_556]	2.05	0.15
2:A:478:HOH:O	2:D:447:HOH:O[2_655]	2.13	0.07

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	354/413 (86%)	343 (97%)	11 (3%)	0	100	100
1	B	354/413 (86%)	341 (96%)	13 (4%)	0	100	100
1	C	354/413 (86%)	342 (97%)	12 (3%)	0	100	100
1	D	354/413 (86%)	339 (96%)	15 (4%)	0	100	100
All	All	1416/1652 (86%)	1365 (96%)	51 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	292/339 (86%)	282 (97%)	10 (3%)	32	44
1	B	292/339 (86%)	282 (97%)	10 (3%)	32	44
1	C	292/339 (86%)	282 (97%)	10 (3%)	32	44
1	D	292/339 (86%)	282 (97%)	10 (3%)	32	44
All	All	1168/1356 (86%)	1128 (97%)	40 (3%)	32	44

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	LEU
1	A	6	ARG
1	A	119	VAL
1	A	210	VAL
1	A	214	VAL
1	A	218	LEU
1	A	237	TRP
1	A	250	LEU
1	A	263	ILE
1	A	320	LEU
1	B	3	LEU
1	B	6	ARG
1	B	119	VAL
1	B	210	VAL
1	B	214	VAL
1	B	218	LEU
1	B	237	TRP
1	B	250	LEU
1	B	263	ILE
1	B	320	LEU
1	C	3	LEU
1	C	6	ARG
1	C	119	VAL
1	C	210	VAL

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Mol	Chain	Res	Type
1	C	214	VAL
1	C	218	LEU
1	C	237	TRP
1	C	250	LEU
1	C	263	ILE
1	C	320	LEU
1	D	3	LEU
1	D	6	ARG
1	D	119	VAL
1	D	210	VAL
1	D	214	VAL
1	D	218	LEU
1	D	237	TRP
1	D	250	LEU
1	D	263	ILE
1	D	320	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (45) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	124	HIS
1	A	128	GLN
1	A	140	HIS
1	A	151	GLN
1	A	234	HIS
1	A	268	ASN
1	A	291	HIS
1	A	341	GLN
1	A	343	HIS
1	A	346	HIS
1	A	364	HIS
1	B	124	HIS
1	B	128	GLN
1	B	140	HIS
1	B	151	GLN
1	B	268	ASN
1	B	278	GLN
1	B	291	HIS
1	B	341	GLN
1	B	343	HIS
1	B	346	HIS
1	B	364	HIS

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Mol	Chain	Res	Type
1	C	124	HIS
1	C	128	GLN
1	C	140	HIS
1	C	151	GLN
1	C	234	HIS
1	C	268	ASN
1	C	271	HIS
1	C	291	HIS
1	C	341	GLN
1	C	343	HIS
1	C	346	HIS
1	C	364	HIS
1	D	124	HIS
1	D	128	GLN
1	D	140	HIS
1	D	151	GLN
1	D	234	HIS
1	D	268	ASN
1	D	291	HIS
1	D	341	GLN
1	D	343	HIS
1	D	346	HIS
1	D	364	HIS

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 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	358/413 (86%)	0.42	19 (5%) 32 29	25, 39, 62, 75	0
1	B	358/413 (86%)	0.44	19 (5%) 32 29	25, 41, 61, 76	0
1	C	358/413 (86%)	0.73	27 (7%) 20 17	26, 47, 64, 76	0
1	D	358/413 (86%)	0.73	29 (8%) 18 15	26, 48, 65, 76	0
All	All	1432/1652 (86%)	0.58	94 (6%) 24 21	25, 44, 63, 76	0

All (94) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	78	TRP	6.3
1	C	78	TRP	6.0
1	D	78	TRP	5.4
1	B	78	TRP	4.5
1	B	77	TYR	4.2
1	B	388	GLU	3.6
1	C	76	ALA	3.6
1	B	3	LEU	3.5
1	B	155	PRO	3.4
1	D	3	LEU	3.4
1	C	156	GLY	3.3
1	C	3	LEU	3.3
1	A	77	TYR	3.3
1	A	385	TYR	3.3
1	D	366	LEU	3.3
1	A	76	ALA	3.2
1	B	76	ALA	3.2
1	C	388	GLU	3.2
1	B	6	ARG	3.1
1	C	77	TYR	3.1
1	C	27	GLU	3.1

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Mol	Chain	Res	Type	RSRZ
1	B	27	GLU	3.0
1	B	80	ARG	3.0
1	D	27	GLU	2.9
1	A	388	GLU	2.9
1	B	30	THR	2.9
1	C	6	ARG	2.9
1	A	41	ARG	2.8
1	A	79	ARG	2.8
1	D	77	TYR	2.8
1	C	44	ALA	2.8
1	B	48	TYR	2.8
1	C	385	TYR	2.8
1	C	154	VAL	2.8
1	A	80	ARG	2.7
1	A	27	GLU	2.7
1	D	79	ARG	2.7
1	C	80	ARG	2.6
1	B	126	THR	2.6
1	C	29	GLY	2.6
1	B	154	VAL	2.6
1	D	350	THR	2.6
1	A	3	LEU	2.6
1	D	156	GLY	2.6
1	D	385	TYR	2.6
1	D	76	ALA	2.5
1	D	377	ILE	2.5
1	C	126	THR	2.4
1	C	31	HIS	2.4
1	A	29	GLY	2.4
1	C	30	THR	2.4
1	C	51	GLU	2.4
1	A	383	ALA	2.4
1	D	370	SER	2.4
1	D	143	LEU	2.4
1	C	64	ARG	2.3
1	C	19	PHE	2.3
1	A	143	LEU	2.3
1	B	16	GLY	2.3
1	C	5	ILE	2.3
1	B	12	VAL	2.3
1	D	25	VAL	2.3
1	A	154	VAL	2.3

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Mol	Chain	Res	Type	RSRZ
1	D	80	ARG	2.2
1	C	48	TYR	2.2
1	D	388	GLU	2.2
1	A	141	VAL	2.2
1	A	155	PRO	2.2
1	D	386	PRO	2.2
1	C	128	GLN	2.2
1	B	156	GLY	2.2
1	D	73	TYR	2.2
1	B	381	LEU	2.2
1	D	154	VAL	2.2
1	D	19	PHE	2.2
1	A	142	GLU	2.2
1	C	266	VAL	2.1
1	B	128	GLN	2.1
1	C	58	ILE	2.1
1	C	10	THR	2.1
1	B	385	TYR	2.1
1	D	31	HIS	2.1
1	D	360	PHE	2.1
1	A	81	GLY	2.1
1	D	6	ARG	2.1
1	A	126	THR	2.1
1	D	47	ALA	2.1
1	D	351	ASP	2.1
1	C	62	ALA	2.0
1	D	52	HIS	2.0
1	D	57	LEU	2.0
1	C	54	VAL	2.0
1	D	346	HIS	2.0
1	D	355	PRO	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.