



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 12:20 PM UTC

PDB ID : 2E1V / pdb_00002e1v
Title : Crystal structure of Dendranthema morifolium DmAT, seleno-methionine derivative
Authors : Unno, H.; Ichimaida, F.; Kusunoki, M.; Nakayama, T.
Deposited on : 2006-10-28
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

<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

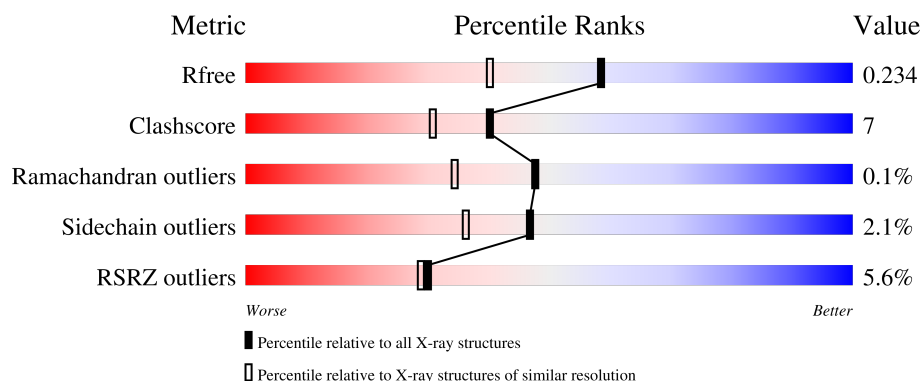
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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}	180053	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (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 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	454	5% 88% 8% ..
1	B	454	6% 85% 11% ..

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7878 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 acyl transferase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	440	Total	C	N	O	S	Se	0	0	0
			3466	2229	573	646	12	6			
1	B	440	Total	C	N	O	S	Se	0	0	0
			3466	2229	573	646	12	6			

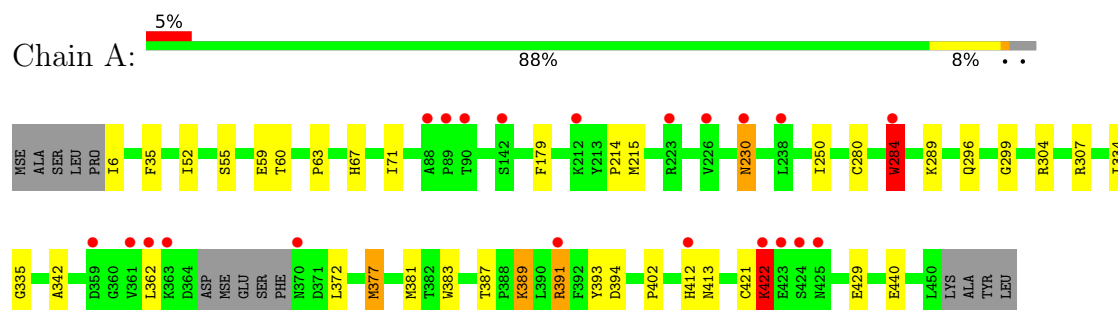
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	477	Total	O	0	0
			477	477		
2	B	469	Total	O	0	0
			469	469		

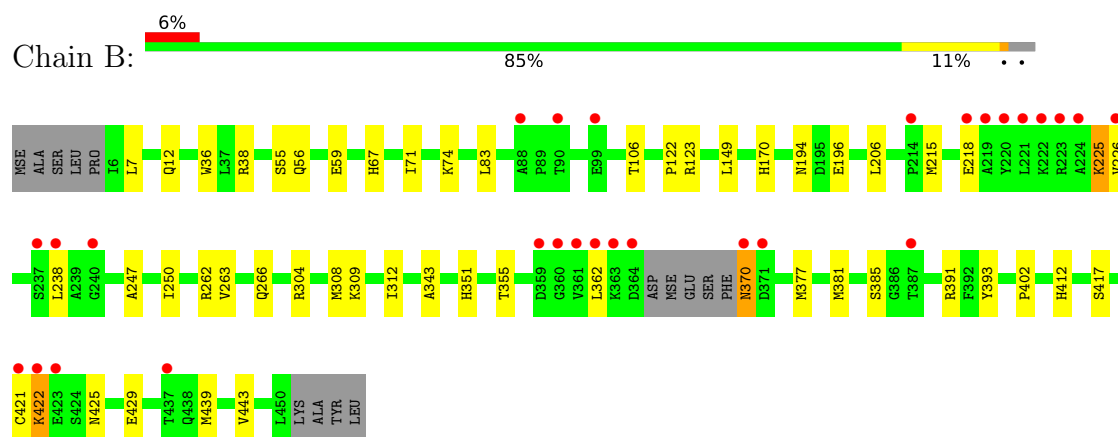
3 Residue-property plots [i](#)

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: acyl transferase



- Molecule 1: acyl transferase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	52.78Å 122.91Å 70.33Å 90.00° 94.38° 90.00°	Depositor
Resolution (Å)	30.46 – 1.80 30.46 – 1.80	Depositor EDS
% Data completeness (in resolution range)	98.5 (30.46-1.80) 98.5 (30.46-1.80)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.64 (at 1.80Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.196 , 0.234 0.196 , 0.234	Depositor DCC
R_{free} test set	4079 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å ²)	23.7	Xtriage
Anisotropy	0.063	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 45.4	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	0.96	EDS
Total number of atoms	7878	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 5.03% 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	0.55	0/3543	0.75	2/4803 (0.0%)
1	B	0.52	0/3543	0.74	0/4803
All	All	0.53	0/7086	0.75	2/9606 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	284	TRP	N-CA-CB	-5.63	101.84	110.01
1	A	422	LYS	N-CA-C	5.59	121.90	114.12

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3466	0	3457	53	0
1	B	3466	0	3457	41	0
2	A	477	0	0	8	0
2	B	469	0	0	9	0
All	All	7878	0	6914	91	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 7.

All (91) 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:B:55:SER:O	1:B:59:GLU:HG2	1.46	1.16
1:A:393:TYR:HB3	1:A:422:LYS:HG2	1.41	1.02
1:A:393:TYR:HB3	1:A:422:LYS:CG	1.93	0.98
1:A:289:LYS:HE2	1:A:335:GLY:H	1.25	0.96
1:B:12:GLN:OE1	1:B:106:THR:HG22	1.72	0.90
1:A:393:TYR:CB	1:A:422:LYS:HG2	2.04	0.87
1:A:284:TRP:HZ3	1:A:296:GLN:OE1	1.60	0.85
1:B:262:ARG:HD2	1:B:266:GLN:HE21	1.42	0.83
1:B:439:MSE:HE3	1:B:443:VAL:CG2	2.09	0.82
1:A:284:TRP:CZ2	2:A:638:HOH:O	2.32	0.80
1:A:289:LYS:CE	1:A:335:GLY:H	1.99	0.74
1:B:421:CYS:SG	1:B:429:GLU:HG3	2.29	0.72
1:B:439:MSE:SE	2:B:851:HOH:O	2.58	0.71
1:A:284:TRP:HZ2	2:A:638:HOH:O	1.70	0.71
1:A:381:MSE:SE	1:B:381:MSE:SE	3.09	0.69
1:B:71:ILE:O	1:B:74:LYS:HG2	1.93	0.69
1:B:391:ARG:NH1	1:B:421:CYS:O	2.18	0.68
1:B:106:THR:HG23	2:B:459:HOH:O	1.93	0.68
1:A:289:LYS:HE2	1:A:335:GLY:N	2.06	0.67
1:A:280:CYS:O	1:A:284:TRP:HB2	1.95	0.66
1:B:247:ALA:HB2	1:B:439:MSE:HE2	1.79	0.65
1:A:179:PHE:HZ	1:A:391:ARG:HD2	1.63	0.63
1:B:122:PRO:HG3	1:B:439:MSE:HE2	1.80	0.63
1:B:38:ARG:NH2	1:B:225:LYS:O	2.25	0.62
1:A:230:ASN:H	1:A:230:ASN:HD22	1.46	0.61
1:A:284:TRP:NE1	2:A:638:HOH:O	2.34	0.59
1:B:247:ALA:CB	1:B:439:MSE:HE1	2.33	0.59
1:A:284:TRP:CE2	2:A:638:HOH:O	2.56	0.59
1:B:351:HIS:O	1:B:355:THR:HG23	2.02	0.58
1:B:238:LEU:HB2	2:B:856:HOH:O	2.03	0.58
1:A:393:TYR:HB3	1:A:422:LYS:CB	2.33	0.57
1:B:412:HIS:HE1	2:B:831:HOH:O	1.87	0.57
1:A:284:TRP:HH2	1:A:296:GLN:HB3	1.69	0.57
1:A:284:TRP:CZ3	1:A:296:GLN:OE1	2.49	0.57
1:B:247:ALA:CB	1:B:439:MSE:CE	2.82	0.57
1:A:284:TRP:CH2	1:A:296:GLN:HB3	2.41	0.56
1:A:391:ARG:HG3	2:A:555:HOH:O	2.06	0.56
1:B:250:ILE:CD1	1:B:429:GLU:HG2	2.37	0.54
1:A:412:HIS:ND1	1:A:413:ASN:ND2	2.55	0.54
1:A:55:SER:O	1:A:59:GLU:HG2	2.07	0.53
1:A:230:ASN:ND2	2:A:791:HOH:O	2.42	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:250:ILE:CD1	1:A:429:GLU:HG2	2.39	0.53
1:A:284:TRP:HH2	1:A:296:GLN:CB	2.21	0.53
1:A:421:CYS:SG	1:A:429:GLU:HG3	2.49	0.52
1:A:179:PHE:CZ	1:A:391:ARG:HD2	2.43	0.52
1:B:262:ARG:HD2	1:B:266:GLN:NE2	2.18	0.52
1:B:56:GLN:HG2	2:B:909:HOH:O	2.09	0.52
1:B:122:PRO:HG3	1:B:439:MSE:CE	2.39	0.52
1:A:289:LYS:HG2	1:A:334:ILE:HG23	1.92	0.51
1:A:440:GLU:HB3	2:A:868:HOH:O	2.11	0.51
1:B:215:MSE:SE	2:B:538:HOH:O	2.79	0.51
1:B:304:ARG:HH22	1:B:362:LEU:HA	1.76	0.51
1:B:247:ALA:HB2	1:B:439:MSE:CE	2.40	0.51
1:A:391:ARG:HD3	1:A:394:ASP:HB2	1.92	0.50
1:A:393:TYR:HB3	1:A:422:LYS:HB2	1.92	0.50
1:B:247:ALA:HB3	1:B:439:MSE:HE1	1.93	0.50
1:A:393:TYR:HB2	1:A:422:LYS:HG2	1.93	0.49
1:A:304:ARG:HH22	1:A:362:LEU:HA	1.77	0.49
1:A:389:LYS:HE3	1:A:389:LYS:HA	1.94	0.49
1:B:391:ARG:HD3	1:B:425:ASN:ND2	2.28	0.49
1:B:391:ARG:HH22	1:B:422:LYS:HA	1.78	0.48
1:A:214:PRO:C	1:A:215:MSE:HE2	2.39	0.47
1:A:393:TYR:O	1:A:402:PRO:HD2	2.15	0.47
1:A:284:TRP:CH2	1:A:296:GLN:CB	2.98	0.47
1:A:391:ARG:HB2	1:A:422:LYS:NZ	2.30	0.46
1:B:194:ASN:OD1	1:B:196:GLU:HB3	2.16	0.46
1:A:383:TRP:HB2	1:B:377:MSE:HG3	1.98	0.46
1:B:439:MSE:HE3	1:B:443:VAL:HG23	1.95	0.45
1:A:372:LEU:HB3	1:A:377:MSE:CE	2.47	0.45
1:B:308:MSE:HG3	1:B:312:ILE:HG12	1.98	0.45
1:A:230:ASN:HD22	1:A:230:ASN:N	2.13	0.45
1:B:263:VAL:HG13	2:B:855:HOH:O	2.16	0.45
1:A:307:ARG:HD2	2:A:543:HOH:O	2.17	0.45
1:A:377:MSE:HA	1:A:377:MSE:HE2	1.99	0.45
1:A:421:CYS:O	1:A:422:LYS:HE3	2.17	0.44
1:B:309:LYS:HB2	2:B:906:HOH:O	2.17	0.43
1:A:391:ARG:HB2	1:A:422:LYS:HZ2	1.83	0.43
1:B:83:LEU:HB2	1:B:149:LEU:HD21	1.99	0.43
1:B:393:TYR:O	1:B:402:PRO:HD2	2.18	0.43
1:A:372:LEU:HB3	1:A:377:MSE:HE3	2.00	0.43
1:B:36:TRP:HB3	1:B:170:HIS:HB3	2.02	0.42
1:A:299:GLY:HA3	1:A:383:TRP:CZ3	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:284:TRP:CD1	1:A:342:ALA:HB1	2.55	0.41
1:A:372:LEU:CB	1:A:377:MSE:HE3	2.50	0.41
1:A:35:PHE:CZ	1:B:370:ASN:HB2	2.56	0.41
1:B:67:HIS:CE1	1:B:71:ILE:HD11	2.56	0.41
1:A:391:ARG:CB	1:A:422:LYS:NZ	2.84	0.41
1:A:60:THR:C	1:A:63:PRO:HD2	2.46	0.41
1:A:67:HIS:CE1	1:A:71:ILE:HD11	2.57	0.40
1:B:385:SER:O	1:B:417:SER:HA	2.22	0.40
1:B:343:ALA:HB3	2:B:855:HOH:O	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	436/454 (96%)	429 (98%)	7 (2%)	0	100	100
1	B	436/454 (96%)	428 (98%)	7 (2%)	1 (0%)	43	31
All	All	872/908 (96%)	857 (98%)	14 (2%)	1 (0%)	48	34

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	225	LYS

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	388/392 (99%)	379 (98%)	9 (2%)	44	33
1	B	388/392 (99%)	381 (98%)	7 (2%)	51	43
All	All	776/784 (99%)	760 (98%)	16 (2%)	47	36

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

Mol	Chain	Res	Type
1	A	6	ILE
1	A	52	ILE
1	A	230	ASN
1	A	284	TRP
1	A	377	MSE
1	A	387	THR
1	A	389	LYS
1	A	391	ARG
1	A	422	LYS
1	B	7	LEU
1	B	123	ARG
1	B	206	LEU
1	B	218	GLU
1	B	226	VAL
1	B	370	ASN
1	B	422	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	67	HIS
1	A	230	ASN
1	A	236	GLN
1	A	266	GLN
1	A	438	GLN
1	A	444	HIS
1	B	67	HIS
1	B	266	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides 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	434/454 (95%)	0.25	21 (4%) 35 34	14, 23, 40, 60	0
1	B	434/454 (95%)	0.28	28 (6%) 25 23	14, 23, 40, 64	0
All	All	868/908 (95%)	0.27	49 (5%) 30 29	14, 23, 40, 64	0

All (49) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	224	ALA	6.9
1	A	370	ASN	5.7
1	B	361	VAL	4.9
1	B	90	THR	4.6
1	A	422	LYS	4.4
1	A	89	PRO	4.3
1	B	359	ASP	4.3
1	A	412	HIS	4.3
1	B	423	GLU	4.2
1	B	370	ASN	4.0
1	A	90	THR	3.8
1	B	226	VAL	3.6
1	A	391	ARG	3.4
1	A	423	GLU	3.3
1	B	362	LEU	3.3
1	B	238	LEU	3.2
1	A	212	LYS	3.1
1	A	226	VAL	3.1
1	B	221	LEU	3.1
1	B	223	ARG	3.0
1	B	220	TYR	3.0
1	A	88	ALA	2.9
1	A	284	TRP	2.9
1	B	218	GLU	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	387	THR	2.8
1	A	359	ASP	2.8
1	B	222	LYS	2.8
1	B	422	LYS	2.8
1	B	437	THR	2.7
1	A	230	ASN	2.6
1	B	371	ASP	2.5
1	A	361	VAL	2.5
1	B	240	GLY	2.4
1	B	360	GLY	2.4
1	A	362	LEU	2.3
1	B	88	ALA	2.3
1	A	142	SER	2.3
1	A	363	LYS	2.3
1	B	363	LYS	2.3
1	A	424	SER	2.2
1	B	364	ASP	2.2
1	B	237	SER	2.2
1	A	425	ASN	2.2
1	B	99	GLU	2.1
1	A	223	ARG	2.1
1	A	238	LEU	2.1
1	B	214	PRO	2.1
1	B	219	ALA	2.1
1	B	421	CYS	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.