Full wwPDB X-ray Structure Validation Report (i)

Jun 12, 2024 – 03:47 PM EDT

PDB ID	:	3B3G
Title	:	The 2.4 A crystal structure of the apo catalytic domain of coactivator-associated arginine methyl transferase I(CARM1,140-480).
Authors	:	Troffer-Charlier, N.; Cura, V.; Hassenboehler, P.; Moras, D.; Cavarelli, J.
Deposited on	:	2007-10-22
Resolution	:	2.40 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 (i) 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 (1)) were used in the production of this report:

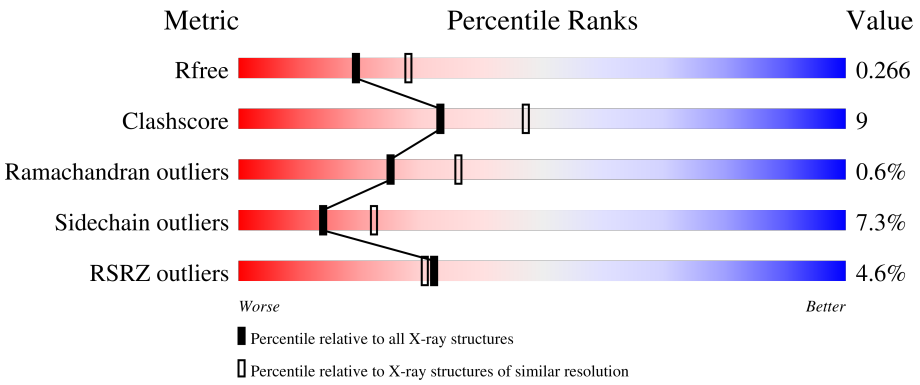
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.20.1
EDS	:	2.36.2
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

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 <=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	341	4%, 74%, 19%, •• 5%
1	B	341	4%, 73%, 19%, • 5%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 5233 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 Histone-arginine methyltransferase CARM1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	324	2588	1671	430	473	14	0	0	0
1	B	324	2588	1671	430	473	14	0	0	0

- Molecule 2 is water.

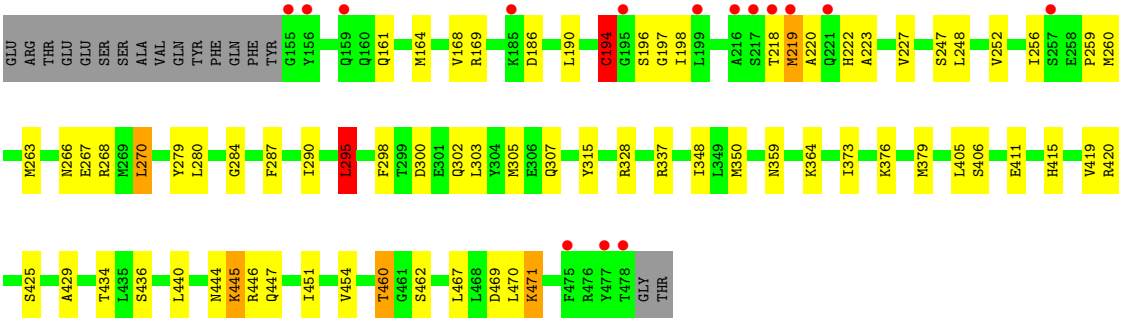
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	34	34	34	0	0
2	B	23	23	23	0	0

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.

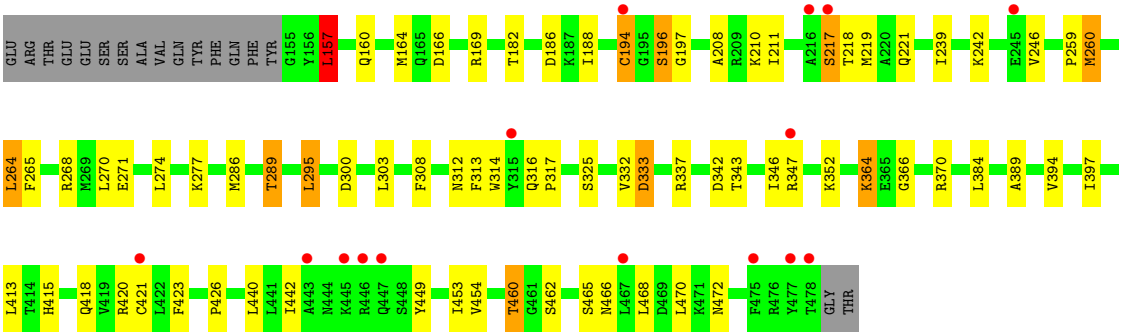
• Molecule 1: Histone-arginine methyltransferase CARM1

Chain A:



• Molecule 1: Histone-arginine methyltransferase CARM1

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α, β, γ	73.82Å 98.12Å 207.13Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.49 – 2.40 29.49 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.0 (29.49-2.40) 99.0 (29.49-2.40)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	0.04	Depositor
$< I/\sigma(I) >$ [1]	7.68 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.205 , 0.265 0.205 , 0.266	Depositor DCC
R_{free} test set	1511 reflections (5.11%)	wwPDB-VP
Wilson B-factor (Å^2)	44.3	Xtriage
Anisotropy	0.207	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 46.6	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	5233	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 4.03% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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	0.60	0/2653	0.69	1/3596 (0.0%)
1	B	0.57	0/2653	0.69	2/3596 (0.1%)
All	All	0.59	0/5306	0.69	3/7192 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	295	LEU	CA-CB-CG	6.31	129.82	115.30
1	B	157	LEU	CB-CG-CD1	5.15	119.76	111.00
1	B	157	LEU	CA-CB-CG	5.04	126.89	115.30

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	2588	0	2557	44	0
1	B	2588	0	2557	45	0
2	A	34	0	0	0	0
2	B	23	0	0	1	0
All	All	5233	0	5114	88	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 9.

All (88) 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:445:LYS:H	1:A:445:LYS:HE2	1.43	0.83
1:B:466:ASN:HB3	2:B:42:HOH:O	1.81	0.80
1:A:194:CYS:HB2	1:A:197:GLY:HA2	1.66	0.76
1:B:194:CYS:HB3	1:B:197:GLY:HA2	1.67	0.76
1:A:446:ARG:NH1	1:A:469:ASP:OD2	2.20	0.74
1:B:289:THR:CG2	1:B:397:ILE:H	2.00	0.73
1:A:169:ARG:HD2	1:A:415:HIS:HE1	1.51	0.73
1:A:263:MET:HE2	1:A:419:VAL:HG11	1.70	0.73
1:A:376:LYS:HE2	1:A:436:SER:HB3	1.69	0.72
1:A:445:LYS:H	1:A:445:LYS:CE	2.03	0.72
1:A:300:ASP:OD2	1:A:420:ARG:NH2	2.24	0.70
1:B:160:GLN:NE2	1:B:219:MET:SD	2.64	0.69
1:B:364:LYS:HE2	1:B:366:GLY:H	1.57	0.69
1:A:460:THR:HG22	1:A:462:SER:H	1.60	0.67
1:A:169:ARG:HD2	1:A:415:HIS:CE1	2.29	0.66
1:B:169:ARG:HH11	1:B:415:HIS:CE1	2.13	0.66
1:B:460:THR:HG22	1:B:462:SER:H	1.61	0.65
1:B:364:LYS:CE	1:B:366:GLY:H	2.10	0.64
1:B:342:ASP:OD2	1:B:420:ARG:NH1	2.31	0.63
1:A:337:ARG:HG2	1:A:467:LEU:O	2.01	0.60
1:A:164:MET:HG3	1:A:196:SER:HB2	1.84	0.59
1:A:446:ARG:O	1:A:471:LYS:HE3	2.04	0.57
1:A:266:ASN:O	1:A:267:GLU:HB2	2.05	0.55
1:A:247:SER:HA	1:A:279:TYR:OH	2.06	0.55
1:B:268:ARG:HD2	1:B:271:GLU:OE2	2.06	0.55
1:A:280:LEU:HD11	1:A:284:GLY:HA3	1.87	0.55
1:A:379:MET:HG3	1:A:429:ALA:HB1	1.90	0.54
1:B:188:ILE:HD12	1:B:210:LYS:HB3	1.89	0.54
1:B:421:CYS:SG	1:B:468:LEU:HD22	2.48	0.54
1:B:166:ASP:OD2	1:B:415:HIS:ND1	2.40	0.53
1:A:259:PRO:HD2	1:A:260:MET:SD	2.48	0.53
1:A:266:ASN:HB2	1:A:470:LEU:HD21	1.90	0.53
1:B:260:MET:HB3	1:B:264:LEU:HD23	1.90	0.53
1:A:350:MET:HA	1:A:350:MET:HE2	1.91	0.52
1:A:290:ILE:HG22	1:A:359:ASN:HA	1.93	0.51
1:A:223:ALA:O	1:A:227:VAL:HG23	2.11	0.51
1:B:289:THR:HG23	1:B:397:ILE:H	1.73	0.50
1:A:263:MET:HE3	1:A:451:ILE:CD1	2.41	0.50
1:B:242:LYS:HB2	1:B:242:LYS:HZ3	1.76	0.50
1:A:218:THR:C	1:A:220:ALA:H	2.15	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:333:ASP:O	1:B:337:ARG:HG3	2.12	0.49
1:B:217:SER:C	1:B:219:MET:H	2.16	0.49
1:B:314:TRP:O	1:B:325:SER:HA	2.13	0.49
1:A:376:LYS:HE2	1:A:436:SER:CB	2.42	0.48
1:B:208:ALA:HB3	1:B:211:ILE:HD11	1.96	0.48
1:A:259:PRO:HG2	1:A:270:LEU:HD13	1.96	0.48
1:B:169:ARG:NH1	1:B:415:HIS:NE2	2.62	0.48
1:B:164:MET:CG	1:B:196:SER:HB3	2.45	0.47
1:B:423:PHE:HD1	1:B:466:ASN:ND2	2.13	0.47
1:A:168:VAL:HG12	1:A:405:LEU:HD11	1.94	0.47
1:B:295:LEU:HA	1:B:389:ALA:O	2.15	0.47
1:A:219:MET:HE2	1:A:222:HIS:HD2	1.79	0.47
1:B:384:LEU:HD23	1:B:426:PRO:HB2	1.97	0.46
1:B:217:SER:HB2	1:B:219:MET:H	1.80	0.46
1:B:300:ASP:OD2	1:B:420:ARG:NH2	2.47	0.46
1:B:370:ARG:HG3	1:B:442:ILE:CD1	2.46	0.46
1:A:190:LEU:HD13	1:A:248:LEU:HD11	1.97	0.46
1:B:217:SER:C	1:B:219:MET:N	2.69	0.45
1:A:406:SER:HB3	1:A:411:GLU:HG3	1.99	0.45
1:A:259:PRO:CG	1:A:270:LEU:HD13	2.46	0.45
1:B:277:LYS:HD2	1:B:286:MET:SD	2.56	0.45
1:B:415:HIS:HD1	1:B:415:HIS:H	1.64	0.44
1:A:256:ILE:HG22	1:A:287:PHE:HB2	1.98	0.44
1:B:259:PRO:HD2	1:B:260:MET:SD	2.58	0.44
1:A:196:SER:OG	1:A:198:ILE:HG22	2.17	0.44
1:A:194:CYS:CB	1:A:197:GLY:HA2	2.39	0.44
1:A:161:GLN:HG3	1:B:313:PHE:CD2	2.54	0.43
1:B:164:MET:HG3	1:B:196:SER:HB3	2.00	0.43
1:A:263:MET:HE3	1:A:451:ILE:HD13	2.00	0.43
1:B:270:LEU:O	1:B:274:LEU:HG	2.18	0.43
1:A:168:VAL:HG12	1:A:405:LEU:CD1	2.49	0.42
1:B:454:VAL:HG22	1:B:465:SER:HB2	2.01	0.42
1:B:423:PHE:CD1	1:B:466:ASN:ND2	2.87	0.42
1:A:190:LEU:HB2	1:A:252:VAL:HG11	2.01	0.42
1:A:268:ARG:HD3	1:A:447:GLN:HA	2.02	0.42
1:B:316:GLN:HA	1:B:317:PRO:HD3	1.90	0.41
1:B:352:LYS:HB2	1:B:352:LYS:HE2	1.96	0.41
1:A:295:LEU:HD11	1:A:373:ILE:HG21	2.01	0.41
1:A:298:PHE:HB2	1:A:348:ILE:HD12	2.03	0.41
1:B:242:LYS:O	1:B:246:VAL:HG13	2.20	0.41
1:A:444:ASN:HB2	1:A:445:LYS:HE2	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:157:LEU:HA	1:B:160:GLN:HG3	2.03	0.41
1:A:445:LYS:HE2	1:A:445:LYS:N	2.23	0.41
1:B:343:THR:HB	1:B:413:LEU:HD21	2.02	0.41
1:B:453:ILE:O	1:B:465:SER:HA	2.21	0.41
1:B:286:MET:O	1:B:289:THR:HB	2.20	0.40
1:A:315:TYR:HB2	1:A:328:ARG:HG3	2.03	0.40
1:B:265:PHE:HB3	1:B:449:TYR:CE1	2.57	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	322/341 (94%)	306 (95%)	14 (4%)	2 (1%)	25	36
1	B	322/341 (94%)	310 (96%)	10 (3%)	2 (1%)	25	36
All	All	644/682 (94%)	616 (96%)	24 (4%)	4 (1%)	25	36

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	194	CYS
1	A	219	MET
1	B	194	CYS
1	B	218	THR

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	281/296 (95%)	265 (94%)	16 (6%)	20 33
1	B	281/296 (95%)	256 (91%)	25 (9%)	9 14
All	All	562/592 (95%)	521 (93%)	41 (7%)	14 22

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	186	ASP
1	A	194	CYS
1	A	270	LEU
1	A	295	LEU
1	A	302	GLN
1	A	303	LEU
1	A	305	MET
1	A	307	GLN
1	A	364	LYS
1	A	425	SER
1	A	434	THR
1	A	440	LEU
1	A	445	LYS
1	A	454	VAL
1	A	460	THR
1	A	471	LYS
1	B	157	LEU
1	B	182	THR
1	B	186	ASP
1	B	196	SER
1	B	217	SER
1	B	221	GLN
1	B	239	ILE
1	B	260	MET
1	B	264	LEU
1	B	289	THR
1	B	295	LEU
1	B	303	LEU
1	B	308	PHE
1	B	312	ASN
1	B	332	VAL
1	B	333	ASP
1	B	346	ILE

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Mol	Chain	Res	Type
1	B	347	ARG
1	B	364	LYS
1	B	394	VAL
1	B	418	GLN
1	B	440	LEU
1	B	460	THR
1	B	470	LEU
1	B	472	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	159	GLN
1	A	162	ASN
1	A	179	GLN
1	A	222	HIS
1	A	381	HIS
1	A	447	GLN
1	B	160	GLN
1	B	179	GLN
1	B	221	GLN
1	B	266	ASN
1	B	312	ASN
1	B	472	ASN

5.3.3 RNA (i)

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 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 (i)

6.1 Protein, DNA and RNA chains (i)

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	324/341 (95%)	0.09	15 (4%)	32	31	21, 41, 63, 74	0
1	B	324/341 (95%)	0.07	15 (4%)	32	31	29, 45, 66, 79	0
All	All	648/682 (95%)	0.08	30 (4%)	32	31	21, 43, 64, 79	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	478	THR	7.1
1	B	477	TYR	6.1
1	A	478	THR	4.9
1	A	477	TYR	4.6
1	A	155	GLY	4.5
1	A	195	GLY	4.4
1	B	315	TYR	3.8
1	A	219	MET	3.7
1	A	218	THR	3.6
1	B	445	LYS	3.6
1	A	217	SER	3.2
1	A	156	TYR	3.2
1	B	194	CYS	3.1
1	A	199	LEU	3.0
1	B	217	SER	2.9
1	B	216	ALA	2.8
1	A	221	GLN	2.8
1	A	159	GLN	2.7
1	B	446	ARG	2.7
1	B	347	ARG	2.6
1	B	245	GLU	2.4
1	B	421	CYS	2.3
1	A	257	SER	2.2
1	A	475	PHE	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	467	LEU	2.1
1	A	216	ALA	2.1
1	B	443	ALA	2.0
1	B	447	GLN	2.0
1	B	475	PHE	2.0
1	A	185	LYS	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.