



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 02:01 AM UTC

PDB ID : 7QDF / pdb_00007qdf
Title : Hexameric HIV-1 (M-group) CA R120 mutant
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Deposited on : 2021-11-26
Resolution : 2.30 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

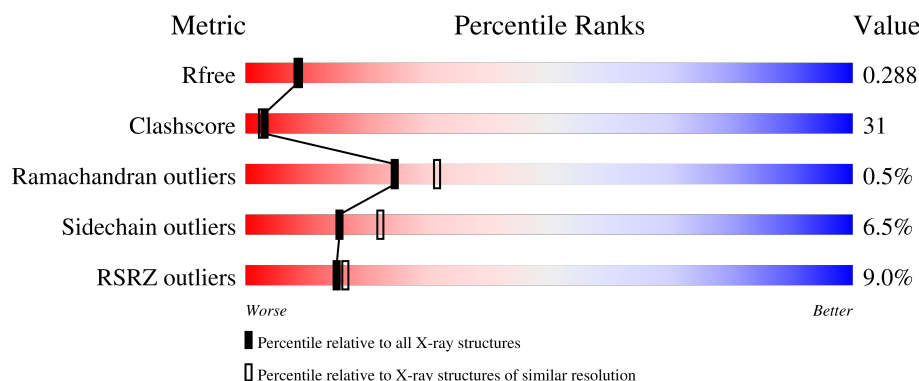
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	AAA	232	9% 55% 37% . .

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	IOD	AAA	302	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 1772 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 Gag polyprotein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	AAA	222	Total	C	N	O	S	0	3	0
			1747	1100	306	328	13			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
AAA	120	ARG	-	insertion	UNP P12493

- Molecule 2 is IODIDE ION (CCD ID: IOD) (formula: I).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	AAA	2	Total	I	0	0
			2	2		

- Molecule 3 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	AAA	5	Total	Cl	0	0
			5	5		

- Molecule 4 is BETA-MERCAPTOETHANOL (CCD ID: BME) (formula: C₂H₆OS).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	AAA	1	Total	C	O	S	0	0
			4	2	1	1		

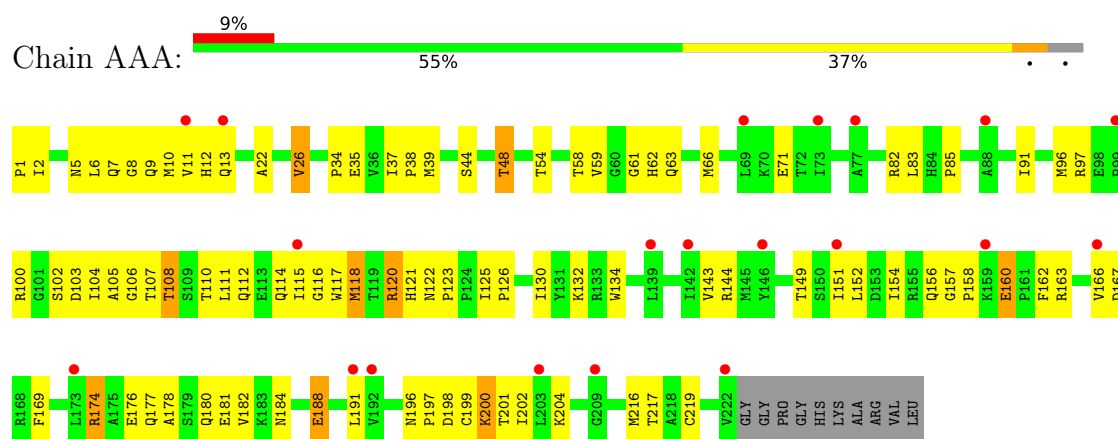
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	AAA	14	Total	O	0	0
			14	14		

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: Gag polypeptide



4 Data and refinement statistics

Property	Value	Source
Space group	P 6	Depositor
Cell constants a, b, c, α , β , γ	92.36Å 92.36Å 57.44Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	46.66 – 2.30 46.66 – 2.30	Depositor EDS
% Data completeness (in resolution range)	98.4 (46.66-2.30) 98.4 (46.66-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.42 (at 2.32Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.242 , 0.278 0.243 , 0.288	Depositor DCC
R_{free} test set	618 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	69.0	Xtriage
Anisotropy	0.171	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 71.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.062 for h,-h-k,-l	Xtriage
Reported twinning fraction	0.918 for H, K, L 0.082 for -K, -H, -L	Depositor
Outliers	1 of 12277 reflections (0.008%)	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	1772	wwPDB-VP
Average B, all atoms (Å ²)	93.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.64% 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: BME, CL, IOD

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	AAA	0.71	0/1785	1.48	19/2426 (0.8%)

There are no bond length outliers.

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AAA	59	VAL	CA-C-N	-11.50	113.93	122.33
1	AAA	59	VAL	C-N-CA	-11.50	113.93	122.33
1	AAA	174	ARG	NE-CZ-NH2	-7.57	112.39	119.20
1	AAA	200	LYS	CA-C-N	6.88	129.73	120.65
1	AAA	200	LYS	C-N-CA	6.88	129.73	120.65
1	AAA	71	GLU	CB-CG-CD	6.73	124.04	112.60
1	AAA	149	THR	CA-C-N	6.03	130.04	121.42
1	AAA	149	THR	C-N-CA	6.03	130.04	121.42
1	AAA	108	THR	CA-CB-OG1	-5.78	100.92	109.60
1	AAA	180	GLN	CB-CA-C	-5.55	101.58	110.79
1	AAA	48	THR	CA-CB-OG1	-5.45	101.43	109.60
1	AAA	201	THR	CA-CB-OG1	-5.41	101.48	109.60
1	AAA	191	LEU	CA-C-N	-5.32	112.99	121.02
1	AAA	191	LEU	C-N-CA	-5.32	112.99	121.02
1	AAA	184	ASN	CB-CA-C	-5.30	102.32	110.81
1	AAA	176	GLU	CA-C-N	5.22	130.27	121.14
1	AAA	176	GLU	C-N-CA	5.22	130.27	121.14
1	AAA	105	ALA	N-CA-C	-5.08	107.05	113.20
1	AAA	174	ARG	NH1-CZ-NH2	5.05	125.86	119.30

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	AAA	1747	0	1741	109	1
2	AAA	2	0	0	3	0
3	AAA	5	0	0	2	0
4	AAA	4	0	6	1	0
5	AAA	14	0	0	3	0
All	All	1772	0	1747	110	1

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 31.

All (110) 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:AAA:96:MET:HE2	1:AAA:117:TRP:CD1	1.55	1.41
1:AAA:48:THR:HG22	1:AAA:118:MET:HE1	1.28	1.15
1:AAA:110:THR:O	1:AAA:114:GLN:HG3	1.49	1.12
1:AAA:35:GLU:O	1:AAA:38:PRO:HD2	1.56	1.04
1:AAA:111:LEU:O	1:AAA:115:ILE:HG13	1.56	1.04
1:AAA:48:THR:HG22	1:AAA:118:MET:CE	1.88	1.02
1:AAA:96:MET:CE	1:AAA:117:TRP:CD1	2.42	1.01
3:AAA:306:CL:CL	5:AAA:410:HOH:O	2.17	0.98
1:AAA:35:GLU:C	1:AAA:38:PRO:HD2	1.92	0.94
1:AAA:96:MET:HE2	1:AAA:117:TRP:HD1	1.15	0.92
1:AAA:174:ARG:NH1	2:AAA:301:IOD:I	2.73	0.91
1:AAA:96:MET:SD	1:AAA:120:ARG:HD3	2.11	0.91
1:AAA:144:ARG:HD3	1:AAA:177:GLN:CA	2.02	0.89
1:AAA:196:ASN:OD1	1:AAA:197:PRO:HD3	1.71	0.88
1:AAA:48:THR:CG2	1:AAA:118:MET:HE1	2.04	0.88
1:AAA:196:ASN:OD1	1:AAA:197:PRO:CD	2.23	0.87
1:AAA:44:SER:OG	1:AAA:132:LYS:HE2	1.76	0.86
1:AAA:144:ARG:HD3	1:AAA:177:GLN:HA	1.58	0.84
1:AAA:158:PRO:HA	1:AAA:196:ASN:ND2	1.93	0.84
1:AAA:12:HIS:CE1	1:AAA:111:LEU:HD21	2.14	0.82
1:AAA:196:ASN:CG	1:AAA:197:PRO:HD2	2.04	0.82
1:AAA:37:ILE:HD11	1:AAA:143:VAL:HG21	1.64	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AAA:103:ASP:OD1	1:AAA:108:THR:HB	1.82	0.79
1:AAA:6:LEU:C	1:AAA:6:LEU:HD12	2.09	0.78
1:AAA:126:PRO:O	1:AAA:130:ILE:HG13	1.85	0.77
1:AAA:6:LEU:HD12	1:AAA:6:LEU:O	1.85	0.76
1:AAA:61:GLY:HA2	1:AAA:63[B]:GLN:HE21	1.49	0.76
1:AAA:144:ARG:CD	1:AAA:177:GLN:HB3	2.16	0.75
1:AAA:163:ARG:HG3	1:AAA:216:MET:CE	2.18	0.74
1:AAA:158:PRO:HA	1:AAA:196:ASN:HD22	1.52	0.72
1:AAA:196:ASN:CG	1:AAA:197:PRO:CD	2.63	0.71
1:AAA:200:LYS:O	1:AAA:204:LYS:HG3	1.92	0.70
1:AAA:34:PRO:O	1:AAA:38:PRO:HD3	1.91	0.70
1:AAA:37:ILE:HD11	1:AAA:143:VAL:CG2	2.22	0.70
1:AAA:82:ARG:NH2	1:AAA:83:LEU:HD11	2.07	0.69
1:AAA:174:ARG:HG3	5:AAA:402:HOH:O	1.91	0.69
1:AAA:96:MET:HE1	1:AAA:116:GLY:C	2.18	0.68
1:AAA:144:ARG:HD2	1:AAA:177:GLN:HB3	1.76	0.67
1:AAA:61:GLY:HA2	1:AAA:63[B]:GLN:NE2	2.09	0.67
1:AAA:111:LEU:HG	1:AAA:115:ILE:HD11	1.77	0.66
1:AAA:82:ARG:HG2	1:AAA:82:ARG:HH11	1.61	0.66
1:AAA:110:THR:O	1:AAA:114:GLN:CG	2.37	0.65
1:AAA:151:ILE:O	1:AAA:151:ILE:HG13	1.96	0.64
1:AAA:196:ASN:OD1	1:AAA:197:PRO:HD2	1.96	0.64
1:AAA:82:ARG:NH2	1:AAA:83:LEU:CD1	2.61	0.64
1:AAA:103:ASP:OD1	1:AAA:108:THR:CB	2.47	0.63
1:AAA:12:HIS:NE2	1:AAA:111:LEU:HD11	2.14	0.62
1:AAA:97:ARG:HH22	1:AAA:103:ASP:CG	2.08	0.62
1:AAA:116:GLY:O	1:AAA:120:ARG:N	2.31	0.62
1:AAA:48:THR:CG2	1:AAA:118:MET:CE	2.72	0.61
1:AAA:96:MET:HE1	1:AAA:117:TRP:N	2.16	0.60
1:AAA:114:GLN:O	1:AAA:118:MET:HG3	2.03	0.59
1:AAA:158:PRO:HD2	2:AAA:302:IOD:I	2.72	0.59
1:AAA:156:GLN:HG3	1:AAA:160:GLU:HG2	1.85	0.58
1:AAA:144:ARG:HD3	1:AAA:177:GLN:CB	2.32	0.58
1:AAA:144:ARG:HD3	1:AAA:177:GLN:HB3	1.86	0.58
1:AAA:48:THR:HG22	1:AAA:118:MET:HE3	1.82	0.58
1:AAA:111:LEU:C	1:AAA:115:ILE:HG13	2.29	0.58
1:AAA:96:MET:SD	1:AAA:120:ARG:CD	2.90	0.56
1:AAA:120:ARG:HG3	1:AAA:123:PRO:O	2.05	0.56
1:AAA:37:ILE:N	1:AAA:38:PRO:CD	2.68	0.56
1:AAA:48:THR:CB	1:AAA:118:MET:HE1	2.35	0.55
1:AAA:54:THR:O	1:AAA:58:THR:HG23	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AAA:82:ARG:HH22	1:AAA:83:LEU:HD11	1.71	0.53
1:AAA:199:CYS:SG	1:AAA:219:CYS:SG	3.06	0.53
1:AAA:120:ARG:NH1	1:AAA:125:ILE:HD11	2.24	0.53
1:AAA:122:ASN:N	1:AAA:123:PRO:CD	2.71	0.53
1:AAA:96:MET:HE1	1:AAA:117:TRP:HA	1.90	0.52
1:AAA:122:ASN:H	1:AAA:123:PRO:CD	2.21	0.52
1:AAA:11:VAL:HG12	1:AAA:13:GLN:HG3	1.92	0.52
1:AAA:108:THR:HG22	1:AAA:108:THR:O	2.09	0.52
1:AAA:122:ASN:H	1:AAA:123:PRO:HD3	1.75	0.52
1:AAA:103:ASP:OD1	1:AAA:108:THR:CG2	2.59	0.51
1:AAA:62:HIS:O	1:AAA:66:MET:HG2	2.10	0.50
1:AAA:5:ASN:HB2	1:AAA:8:GLY:H	1.76	0.50
1:AAA:163:ARG:HB2	3:AAA:305:CL:CL	2.48	0.50
1:AAA:91:ILE:HD12	1:AAA:91:ILE:N	2.26	0.50
1:AAA:96:MET:CG	1:AAA:120:ARG:HH21	2.24	0.50
1:AAA:144:ARG:CD	1:AAA:177:GLN:HA	2.35	0.49
1:AAA:9:GLN:O	1:AAA:9:GLN:HG3	2.13	0.48
1:AAA:110:THR:OG1	1:AAA:112[B]:GLN:HG3	2.15	0.47
1:AAA:96:MET:HE1	1:AAA:117:TRP:CA	2.44	0.47
1:AAA:103:ASP:OD1	1:AAA:108:THR:HG22	2.15	0.47
1:AAA:163:ARG:HG3	1:AAA:216:MET:HE2	1.93	0.47
1:AAA:63[B]:GLN:NE2	5:AAA:401:HOH:O	2.28	0.47
1:AAA:134:TRP:CD1	1:AAA:134:TRP:N	2.82	0.46
1:AAA:10:MET:HB2	1:AAA:10:MET:HE2	1.61	0.46
1:AAA:157:GLY:HA3	2:AAA:302:IOD:I	2.86	0.46
1:AAA:102:SER:HB2	1:AAA:108:THR:OG1	2.16	0.46
1:AAA:111:LEU:O	1:AAA:115:ILE:CG1	2.44	0.46
1:AAA:154:ILE:HG21	1:AAA:169:PHE:HA	1.98	0.45
1:AAA:37:ILE:HG22	4:AAA:308:BME:H21	1.98	0.45
1:AAA:12:HIS:CD2	1:AAA:111:LEU:HD11	2.52	0.45
1:AAA:162:PHE:O	1:AAA:166:VAL:HG23	2.16	0.45
1:AAA:163:ARG:HB3	1:AAA:163:ARG:NH1	2.32	0.45
1:AAA:96:MET:CE	1:AAA:117:TRP:HA	2.46	0.45
1:AAA:6:LEU:O	1:AAA:6:LEU:CD1	2.60	0.44
1:AAA:100:ARG:CZ	1:AAA:100:ARG:HB3	2.48	0.44
1:AAA:37:ILE:CD1	1:AAA:143:VAL:HG21	2.41	0.43
1:AAA:83:LEU:O	1:AAA:85:PRO:HD3	2.18	0.43
1:AAA:178:ALA:HB1	1:AAA:182:VAL:HB	2.00	0.43
1:AAA:1:PRO:N	1:AAA:13:GLN:O	2.33	0.43
1:AAA:26:VAL:HG11	1:AAA:39:MET:HG2	2.01	0.43
1:AAA:104:ILE:C	1:AAA:106:GLY:H	2.27	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AAA:120:ARG:HB3	1:AAA:121:HIS:H	1.44	0.42
1:AAA:12:HIS:NE2	1:AAA:111:LEU:HD21	2.34	0.42
1:AAA:22:ALA:O	1:AAA:26:VAL:HG13	2.20	0.41
1:AAA:144:ARG:HD3	1:AAA:177:GLN:N	2.35	0.41
1:AAA:34:PRO:O	1:AAA:38:PRO:CD	2.67	0.41
1:AAA:188:GLU:H	1:AAA:188:GLU:HG3	1.66	0.40

All (1) 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:AAA:63[B]:GLN:OE1	1:AAA:167:ASP:O[6_555]	1.81	0.39

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	AAA	223/232 (96%)	210 (94%)	12 (5%)	1 (0%)	30 38

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	AAA	7	GLN

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	AAA	189/195 (97%)	177 (94%)	12 (6%)	16	23

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	AAA	2	ILE
1	AAA	26	VAL
1	AAA	107	THR
1	AAA	118	MET
1	AAA	120	ARG
1	AAA	152	LEU
1	AAA	160	GLU
1	AAA	181	GLU
1	AAA	188	GLU
1	AAA	198	ASP
1	AAA	202	ILE
1	AAA	217	THR

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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 8 ligands modelled in this entry, 7 are monoatomic - leaving 1 for Mogul analysis.

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	$\# Z > 2$
4	BME	AAA	308	-	3,3,3	0.27	0	2,2,2	0.26	0

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
4	BME	AAA	308	-	-	0/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	AAA	308	BME	1	0

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	AAA	222/232 (95%)	0.90	20 (9%) 15 16	29, 89, 142, 181	3 (1%)

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	AAA	203	LEU	7.4
1	AAA	13	GLN	6.8
1	AAA	11	VAL	4.2
1	AAA	142	ILE	3.6
1	AAA	151	ILE	3.4
1	AAA	222	VAL	3.4
1	AAA	146	TYR	3.2
1	AAA	209	GLY	3.1
1	AAA	115	ILE	3.0
1	AAA	192	VAL	3.0
1	AAA	73	ILE	2.7
1	AAA	99	PRO	2.4
1	AAA	166	VAL	2.4
1	AAA	69	LEU	2.3
1	AAA	88	ALA	2.2
1	AAA	173	LEU	2.1
1	AAA	191	LEU	2.1
1	AAA	139	LEU	2.1
1	AAA	159	LYS	2.0
1	AAA	77	ALA	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](#)

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($\text{\AA}^2$)	Q<0.9
3	CL	AAA	304	1/1	0.91	0.17	90,90,90,90	1
3	CL	AAA	305	1/1	0.91	0.15	85,85,85,85	1
4	BME	AAA	308	4/4	0.93	0.12	51,52,70,84	0
3	CL	AAA	306	1/1	0.94	0.11	55,55,55,55	1
3	CL	AAA	303	1/1	0.95	0.10	55,55,55,55	1
2	IOD	AAA	302	1/1	0.96	0.08	96,96,96,96	1
3	CL	AAA	307	1/1	0.98	0.10	48,48,48,48	1
2	IOD	AAA	301	1/1	0.99	0.04	67,67,67,67	1

6.5 Other polymers [i](#)

There are no such residues in this entry.