



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 11:41 AM UTC

PDB ID : 7Q8Q / pdb_00007q8q
Title : Peptide RLSAKP in complex with human cathepsin V C25S mutant
Authors : Loboda, J.; Sosnowski, P.; Tusar, L.; Vidmar, R.; Vizovisek, M.; Horvat, J.;
Kosec, G.; Impens, F.; Demol, H.; Turk, B.; Gevaert, K.; Turk, D.
Deposited on : 2021-11-11
Resolution : 2.13 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

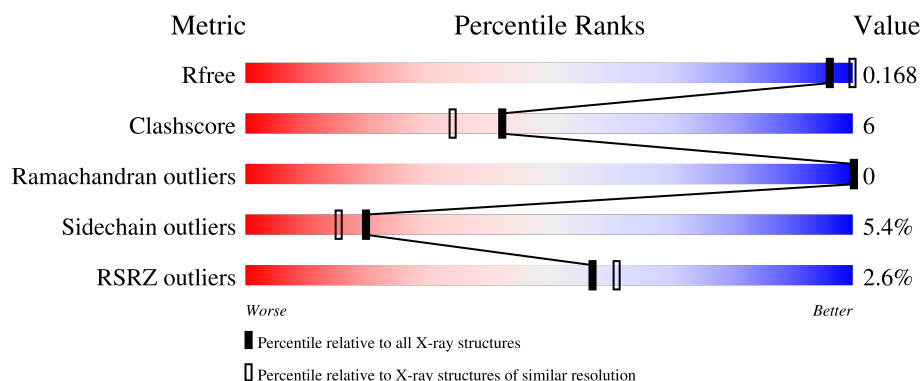
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.13 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	3689 (2.16-2.12)
Clashscore	190562	3812 (2.16-2.12)
Ramachandran outliers	187476	3773 (2.16-2.12)
Sidechain outliers	187428	3772 (2.16-2.12)
RSRZ outliers	180081	3691 (2.16-2.12)

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	AA	221	79% 17% .
1	BA	221	3% 78% 19% .
2	PA	6	33% 50% 33% 17%
2	PB	6	83% 17%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 5154 atoms, of which 1377 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 Cathepsin L2.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	AA	221	Total	C	H	N	O	S	395	0	0
			2080	1065	390	290	325	10			
1	BA	221	Total	C	H	N	O	S	393	0	0
			2081	1065	391	290	325	10			

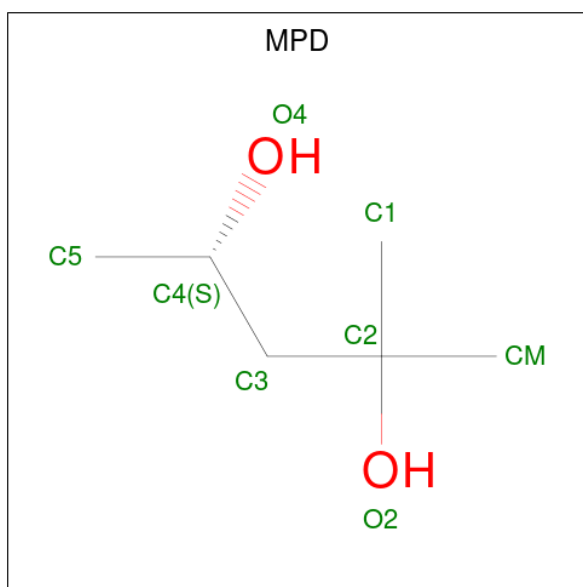
There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
AA	25	SER	CYS	engineered mutation	UNP O60911
AA	108	GLN	ASN	engineered mutation	UNP O60911
AA	179	GLN	ASN	engineered mutation	UNP O60911
BA	246	SER	CYS	engineered mutation	UNP O60911
BA	329	GLN	ASN	engineered mutation	UNP O60911
BA	400	GLN	ASN	engineered mutation	UNP O60911

- Molecule 2 is a protein called RLSAKP Peptide.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	PA	5	Total	C	H	N	O	14	0	0
			53	24	14	9	6			
2	PB	6	Total	C	H	N	O	14	0	0
			61	29	14	10	8			

- Molecule 3 is (4S)-2-METHYL-2,4-PENTANEDIOL (CCD ID: MPD) (formula: C₆H₁₄O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	AA	1	Total	C	O	0	0
			8	6	2		
3	BA	1	Total	C	O	0	0
			8	6	2		
3	PA	1	Total	C	O	0	0
			8	6	2		

- Molecule 4 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	AA	1	Total	Cl	0	0
			1	1		
4	BA	2	Total	Cl	0	0
			2	2		

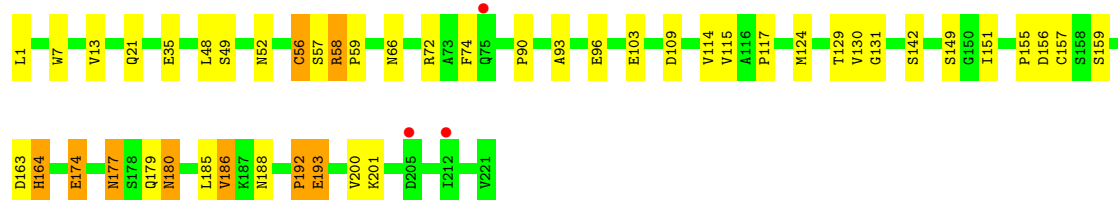
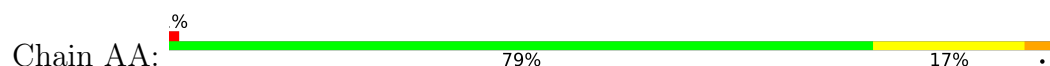
- Molecule 5 is water.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	AA	137	Total	H	O	274	0
			411	274	137		
5	BA	139	Total	H	O	278	0
			417	278	139		
5	PA	6	Total	H	O	12	0
			18	12	6		
5	PB	2	Total	H	O	4	0
			6	4	2		

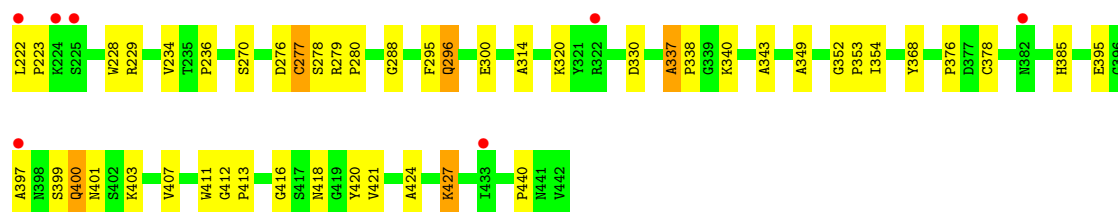
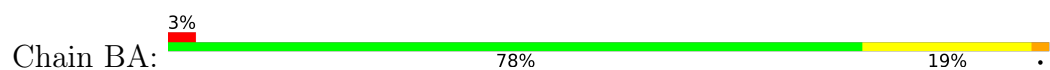
3 Residue-property plots [i](#)

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.

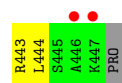
- Molecule 1: Cathepsin L2



- Molecule 1: Cathepsin L2



- Molecule 2: RLSAKP Peptide



- Molecule 2: RLSAKP Peptide



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	92.75Å 92.75Å 128.09Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.61 – 2.13 43.61 – 2.13	Depositor EDS
% Data completeness (in resolution range)	100.0 (43.61-2.13) 99.4 (43.61-2.13)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.73 (at 2.14Å)	Xtriage
Refinement program	MAIN	Depositor
R, R_{free}	0.211 , 0.239 (Not available) , 0.168	Depositor DCC
R_{free} test set	1577 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	36.3	Xtriage
Anisotropy	0.021	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 56.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5154	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.47% 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: MPD, CL

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	AA	1.22	8/1732 (0.5%)	1.41	22/2340 (0.9%)
1	BA	0.86	1/1732 (0.1%)	1.33	23/2340 (1.0%)
2	PA	0.77	0/38	1.55	0/48
2	PB	0.68	0/47	1.31	0/60
All	All	1.05	9/3549 (0.3%)	1.37	45/4788 (0.9%)

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	AA	164	HIS	C-N	-7.46	1.28	1.33
1	AA	124	MET	SD-CE	-6.17	1.64	1.79
1	AA	21	GLN	CA-C	-6.14	1.46	1.53
1	AA	142	SER	CA-C	-6.03	1.44	1.52
1	AA	13	VAL	CA-C	-5.59	1.45	1.52
1	AA	186	VAL	CA-C	-5.32	1.46	1.52
1	BA	296	GLN	CA-C	-5.13	1.46	1.52
1	AA	159	SER	CA-C	-5.07	1.46	1.52
1	AA	164	HIS	CD2-NE2	-5.02	1.32	1.37

All (45) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AA	130	VAL	N-CA-C	9.92	119.94	111.90
1	BA	276	ASP	N-CA-C	8.32	121.48	111.82
1	BA	234	VAL	N-CA-C	8.22	119.62	108.11
1	AA	56	CYS	N-CA-C	7.67	123.57	113.30
1	AA	49	SER	N-CA-C	7.64	120.51	108.67
1	AA	13	VAL	N-CA-C	7.11	118.35	108.12
1	BA	277	CYS	N-CA-C	7.07	122.77	113.30
1	AA	109	ASP	CA-CB-CG	6.84	119.44	112.60

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	BA	400	GLN	CA-C-N	-6.80	113.06	122.93
1	BA	400	GLN	C-N-CA	-6.80	113.06	122.93
1	BA	314	ALA	N-CA-C	6.69	120.81	112.24
1	AA	90	PRO	N-CA-C	6.62	121.90	111.38
1	AA	186	VAL	N-CA-C	6.42	117.11	108.11
1	AA	93	ALA	N-CA-C	6.39	120.66	112.86
1	AA	103	GLU	N-CA-C	-6.32	104.83	112.54
1	BA	330	ASP	CA-CB-CG	6.12	118.72	112.60
1	BA	270	SER	N-CA-C	6.06	118.65	108.23
1	BA	397	ALA	CA-C-N	-5.96	115.09	122.84
1	BA	397	ALA	C-N-CA	-5.96	115.09	122.84
1	AA	74	PHE	N-CA-C	-5.84	104.91	111.28
1	AA	130	VAL	N-CA-CB	-5.64	102.40	111.82
1	BA	407	VAL	N-CA-C	5.58	115.92	108.11
1	AA	13	VAL	N-CA-CB	-5.57	103.39	111.52
1	AA	149	SER	N-CA-C	5.53	116.37	108.74
1	BA	278	SER	N-CA-C	5.49	118.98	112.72
1	AA	1	LEU	CA-C-N	-5.40	114.68	120.03
1	AA	1	LEU	C-N-CA	-5.40	114.68	120.03
1	AA	57	SER	N-CA-C	5.40	118.67	112.57
1	AA	129	THR	N-CA-C	5.39	120.01	113.38
1	BA	401	ASN	CA-CB-CG	5.38	117.98	112.60
1	BA	338	PRO	N-CA-C	5.34	119.26	111.03
1	BA	276	ASP	CA-CB-CG	5.30	117.90	112.60
1	BA	337	ALA	CA-C-N	-5.30	114.10	119.83
1	BA	337	ALA	C-N-CA	-5.30	114.10	119.83
1	BA	338	PRO	CB-CA-C	5.22	118.21	111.64
1	AA	185	LEU	N-CA-C	-5.20	100.69	108.96
1	AA	201	LYS	N-CA-C	-5.20	100.77	109.24
1	BA	427	LYS	N-CA-C	-5.20	102.63	110.06
1	BA	385	HIS	N-CA-C	5.17	116.45	107.93
1	BA	354	ILE	N-CA-C	5.12	115.85	108.48
1	AA	72	ARG	N-CA-C	-5.09	105.73	111.28
1	AA	58	ARG	N-CA-C	-5.08	106.56	113.57
1	BA	349	ALA	CA-C-N	-5.01	114.38	122.24
1	BA	349	ALA	C-N-CA	-5.01	114.38	122.24
1	AA	117	PRO	N-CA-C	5.00	118.73	111.03

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	AA	1690	390	1622	18	1
1	BA	1690	391	1621	20	1
2	PA	39	14	46	2	0
2	PB	47	14	53	2	0
3	AA	8	0	14	2	0
3	BA	8	0	14	1	0
3	PA	8	0	14	2	0
4	AA	1	0	0	0	0
4	BA	2	0	0	1	0
5	AA	137	274	0	4	0
5	BA	139	278	0	1	0
5	PA	6	12	0	0	0
5	PB	2	4	0	0	0
All	All	3777	1377	3384	42	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 6.

All (42) 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:AA:151:ILE:HD11	1:AA:174:GLU:HG3	1.49	0.94
1:BA:295:PHE:CE1	1:BA:440:PRO:HG3	2.04	0.93
1:AA:66:ASN:ND2	5:AA:401:HOH:O	2.07	0.85
1:AA:177:ASN:HB3	1:AA:180:ASN:HD22	1.45	0.81
1:AA:177:ASN:HB3	1:AA:180:ASN:ND2	2.04	0.72
1:AA:180:ASN:O	5:AA:402:HOH:O	2.09	0.69
1:AA:177:ASN:HD22	1:AA:179:GLN:H	1.42	0.68
1:AA:192:PRO:HD2	1:AA:193:GLU:OE1	1.96	0.64
1:AA:7:TRP:CE2	1:AA:131:GLY:HA2	2.36	0.60
1:BA:295:PHE:CD1	1:BA:440:PRO:HG3	2.38	0.59
1:BA:228:TRP:CE2	1:BA:352:GLY:HA2	2.39	0.57
3:PA:501:MPD:O4	3:PA:501:MPD:HM2	2.04	0.57
3:AA:301:MPD:H53	3:PA:501:MPD:O2	2.05	0.56
1:AA:151:ILE:CD1	1:AA:174:GLU:HG3	2.30	0.56

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AA:164:HIS:HD1	2:PA:443:ARG:N	2.03	0.56
1:AA:188:ASN:HB3	5:AA:437:HOH:O	2.09	0.53
1:BA:413:PRO:O	1:BA:418:ASN:HA	2.09	0.53
1:BA:295:PHE:CE1	1:BA:440:PRO:CG	2.88	0.52
1:AA:186:VAL:HB	1:AA:200:VAL:HG13	1.93	0.51
1:BA:279:ARG:N	1:BA:280:PRO:CD	2.75	0.50
1:BA:295:PHE:CZ	1:BA:440:PRO:HG3	2.46	0.49
4:BA:501:CL:CL	2:PB:448:ARG:N	2.84	0.48
1:AA:180:ASN:ND2	5:AA:402:HOH:O	2.47	0.48
1:AA:58:ARG:N	1:AA:59:PRO:CD	2.77	0.47
1:BA:340:LYS:HB2	1:BA:343:ALA:HB3	1.97	0.46
1:BA:337:ALA:CB	1:BA:340:LYS:HG2	2.44	0.46
1:BA:400:GLN:O	1:BA:403:LYS:HE2	2.15	0.46
1:AA:151:ILE:HD11	1:AA:174:GLU:CG	2.34	0.43
1:AA:163:ASP:CG	2:PA:443:ARG:HD2	2.44	0.43
1:BA:368:TYR:CZ	1:BA:421:VAL:HG23	2.53	0.42
1:AA:35:GLU:HG3	1:AA:48:LEU:HG	2.01	0.42
1:BA:368:TYR:O	1:BA:416:GLY:HA2	2.20	0.42
1:BA:412:GLY:HA3	5:BA:635:HOH:O	2.19	0.42
1:BA:229:ARG:HD3	1:BA:420:TYR:CE2	2.55	0.42
1:BA:229:ARG:HD3	1:BA:420:TYR:CZ	2.55	0.42
1:BA:340:LYS:HB2	1:BA:343:ALA:CB	2.50	0.42
1:BA:222:LEU:HA	1:BA:223:PRO:HD3	1.93	0.41
1:BA:411:TRP:HZ2	2:PB:448:ARG:HG3	1.85	0.41
3:AA:301:MPD:O4	3:AA:301:MPD:HM1	2.20	0.41
1:BA:424:ALA:HB1	1:BA:427:LYS:HG3	2.03	0.41
1:BA:288:GLY:HA2	3:BA:502:MPD:H51	2.03	0.40
1:AA:114:VAL:HG12	1:AA:115:VAL:N	2.37	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:AA:96:GLU:OE1	1:BA:320:LYS:HZ1[3_554]	1.38	0.22

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	AA	219/221 (99%)	213 (97%)	6 (3%)	0	100	100
1	BA	219/221 (99%)	210 (96%)	9 (4%)	0	100	100
2	PA	3/6 (50%)	3 (100%)	0	0	100	100
2	PB	4/6 (67%)	4 (100%)	0	0	100	100
All	All	445/454 (98%)	430 (97%)	15 (3%)	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	AA	179/179 (100%)	169 (94%)	10 (6%)	19	14
1	BA	179/179 (100%)	170 (95%)	9 (5%)	22	17
2	PA	4/5 (80%)	3 (75%)	1 (25%)	0	0
2	PB	5/5 (100%)	5 (100%)	0	100	100
All	All	367/368 (100%)	347 (95%)	20 (5%)	20	15

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	AA	52	ASN
1	AA	56	CYS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	AA	155	PRO
1	AA	156	ASP
1	AA	157	CYS
1	AA	174	GLU
1	AA	177	ASN
1	AA	180	ASN
1	AA	192	PRO
1	AA	193	GLU
1	BA	236	PRO
1	BA	277	CYS
1	BA	296	GLN
1	BA	300	GLU
1	BA	353	PRO
1	BA	376	PRO
1	BA	378	CYS
1	BA	395	GLU
1	BA	399	SER
2	PA	444	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	AA	177	ASN
1	AA	180	ASN
1	AA	220	ASN
1	BA	296	GLN
1	BA	428	ASN
1	BA	441	ASN

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 6 ligands modelled in this entry, 3 are monoatomic - leaving 3 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	$\# Z > 2$
3	MPD	AA	301	-	7,7,7	0.39	0	9,10,10	0.73	0
3	MPD	PA	501	-	7,7,7	0.60	0	9,10,10	0.46	0
3	MPD	BA	502	-	7,7,7	0.97	0	9,10,10	0.68	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
3	MPD	AA	301	-	-	0/5/5/5	-
3	MPD	PA	501	-	-	3/5/5/5	-
3	MPD	BA	502	-	-	2/5/5/5	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	PA	501	MPD	O2-C2-C3-C4
3	PA	501	MPD	C2-C3-C4-O4
3	BA	502	MPD	C1-C2-C3-C4
3	BA	502	MPD	CM-C2-C3-C4
3	PA	501	MPD	C1-C2-C3-C4

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	AA	301	MPD	2	0
3	PA	501	MPD	2	0
3	BA	502	MPD	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	AA	221/221 (100%)	-0.03	3 (1%) 73 77	16, 28, 48, 72	2 (0%)
1	BA	221/221 (100%)	0.16	7 (3%) 50 54	18, 33, 56, 100	1 (0%)
2	PA	5/6 (83%)	2.17	2 (40%) 1 1	32, 37, 85, 106	0
2	PB	6/6 (100%)	0.60	0 100 100	32, 34, 42, 65	0
All	All	453/454 (99%)	0.09	12 (2%) 57 61	16, 30, 53, 106	3 (0%)

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	PA	446	ALA	5.8
2	PA	447	LYS	3.4
1	BA	433	ILE	3.1
1	AA	205	ASP	2.9
1	BA	397	ALA	2.9
1	BA	382	ASN	2.7
1	BA	222	LEU	2.6
1	BA	224	LYS	2.3
1	BA	322	ARG	2.3
1	AA	75	GLN	2.1
1	AA	212	ILE	2.1
1	BA	225	SER	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	MPD	PA	501	8/8	0.87	0.20	80,87,92,94	0
3	MPD	BA	502	8/8	0.89	0.21	68,75,78,89	0
3	MPD	AA	301	8/8	0.90	0.16	60,63,64,66	0
4	CL	BA	503	1/1	0.94	0.09	56,56,56,56	0
4	CL	BA	501	1/1	0.97	0.05	38,38,38,38	0
4	CL	AA	302	1/1	0.99	0.04	31,31,31,31	0

6.5 Other polymers [i](#)

There are no such residues in this entry.