



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 01:10 PM UTC

PDB ID : 6NNV / pdb_00006nnv
Title : PD-L1 IgV domain complex with macro-cyclic peptide
Authors : Zhao, B.; Perry, E.
Deposited on : 2019-01-15
Resolution : 1.92 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

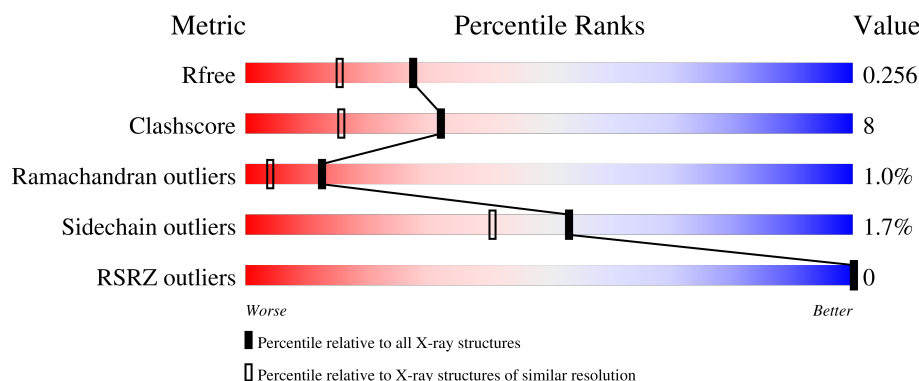
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.92 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	1188 (1.92-1.92)
Clashscore	190562	1209 (1.92-1.92)
Ramachandran outliers	187476	1195 (1.92-1.92)
Sidechain outliers	187428	1195 (1.92-1.92)
RSRZ outliers	180081	1188 (1.92-1.92)

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	127	
1	B	127	
1	C	127	
1	D	127	
2	I	16	

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Mol	Chain	Length	Quality of chain
2	J	16	
2	K	16	
2	L	16	

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	MEA	K	8	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4561 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 Programmed cell death 1 ligand 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	119	Total	C	N	O	S	0	0	0
			925	594	150	176	5			
1	B	121	Total	C	N	O	S	0	0	0
			935	599	151	180	5			
1	C	123	Total	C	N	O	S	0	0	0
			959	614	159	181	5			
1	D	121	Total	C	N	O	S	0	0	0
			928	594	152	177	5			

There are 44 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	76	THR	VAL	engineered mutation	UNP Q9NZQ7
A	135	ALA	-	expression tag	UNP Q9NZQ7
A	136	ALA	-	expression tag	UNP Q9NZQ7
A	137	ALA	-	expression tag	UNP Q9NZQ7
A	138	LEU	-	expression tag	UNP Q9NZQ7
A	139	HIS	-	expression tag	UNP Q9NZQ7
A	140	GLU	-	expression tag	UNP Q9NZQ7
A	141	HIS	-	expression tag	UNP Q9NZQ7
A	142	HIS	-	expression tag	UNP Q9NZQ7
A	143	HIS	-	expression tag	UNP Q9NZQ7
A	144	HIS	-	expression tag	UNP Q9NZQ7
B	76	THR	VAL	engineered mutation	UNP Q9NZQ7
B	135	ALA	-	expression tag	UNP Q9NZQ7
B	136	ALA	-	expression tag	UNP Q9NZQ7
B	137	ALA	-	expression tag	UNP Q9NZQ7
B	138	LEU	-	expression tag	UNP Q9NZQ7
B	139	HIS	-	expression tag	UNP Q9NZQ7
B	140	GLU	-	expression tag	UNP Q9NZQ7
B	141	HIS	-	expression tag	UNP Q9NZQ7
B	142	HIS	-	expression tag	UNP Q9NZQ7
B	143	HIS	-	expression tag	UNP Q9NZQ7

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Chain	Residue	Modelled	Actual	Comment	Reference
B	144	HIS	-	expression tag	UNP Q9NZQ7
C	76	THR	VAL	engineered mutation	UNP Q9NZQ7
C	135	ALA	-	expression tag	UNP Q9NZQ7
C	136	ALA	-	expression tag	UNP Q9NZQ7
C	137	ALA	-	expression tag	UNP Q9NZQ7
C	138	LEU	-	expression tag	UNP Q9NZQ7
C	139	HIS	-	expression tag	UNP Q9NZQ7
C	140	GLU	-	expression tag	UNP Q9NZQ7
C	141	HIS	-	expression tag	UNP Q9NZQ7
C	142	HIS	-	expression tag	UNP Q9NZQ7
C	143	HIS	-	expression tag	UNP Q9NZQ7
C	144	HIS	-	expression tag	UNP Q9NZQ7
D	76	THR	VAL	engineered mutation	UNP Q9NZQ7
D	135	ALA	-	expression tag	UNP Q9NZQ7
D	136	ALA	-	expression tag	UNP Q9NZQ7
D	137	ALA	-	expression tag	UNP Q9NZQ7
D	138	LEU	-	expression tag	UNP Q9NZQ7
D	139	HIS	-	expression tag	UNP Q9NZQ7
D	140	GLU	-	expression tag	UNP Q9NZQ7
D	141	HIS	-	expression tag	UNP Q9NZQ7
D	142	HIS	-	expression tag	UNP Q9NZQ7
D	143	HIS	-	expression tag	UNP Q9NZQ7
D	144	HIS	-	expression tag	UNP Q9NZQ7

- Molecule 2 is a protein called macrocyclic peptide.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	I	16	Total 134	C 97	N 17	O 19	S 1	0	0	1
2	J	16	Total 134	C 97	N 17	O 19	S 1	0	0	1
2	K	16	Total 134	C 97	N 17	O 19	S 1	0	0	1
2	L	16	Total 134	C 97	N 17	O 19	S 1	0	0	1

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	58	Total	O	0	0
			58	58		
3	B	62	Total	O	0	0
			62	62		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	C	63	Total 63	O 63	0	0
3	D	72	Total 72	O 72	0	0
3	I	7	Total 7	O 7	0	0
3	J	7	Total 7	O 7	0	0
3	K	5	Total 5	O 5	0	0
3	L	4	Total 4	O 4	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: Programmed cell death 1 ligand 1

Chain A: 



- Molecule 1: Programmed cell death 1 ligand 1

Chain B: 



- Molecule 1: Programmed cell death 1 ligand 1

Chain C: 



- Molecule 1: Programmed cell death 1 ligand 1

Chain D: 

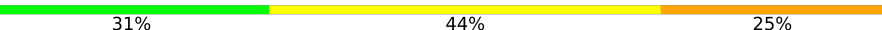


- Molecule 2: macrocyclic peptide

Chain I: 



- Molecule 2: macrocyclic peptide

Chain J: 



● Molecule 2: macrocyclic peptide



● Molecule 2: macrocyclic peptide



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	56.84Å 61.68Å 80.89Å 90.00° 90.17° 90.00°	Depositor
Resolution (Å)	29.10 – 1.92 29.10 – 1.92	Depositor EDS
% Data completeness (in resolution range)	97.0 (29.10-1.92) 96.3 (29.10-1.92)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.56 (at 1.92Å)	Xtriage
Refinement program	PHENIX (1.14_3260: ???)	Depositor
R, R_{free}	0.215 , 0.253 0.219 , 0.256	Depositor DCC
R_{free} test set	2016 reflections (3.85%)	wwPDB-VP
Wilson B-factor (Å ²)	17.6	Xtriage
Anisotropy	0.855	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 23.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	0.139 for h,-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	4561	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 39.78 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 3.0147e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 9KK, ACE, NH2, SAR, MEA

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	0.17	0/942	0.45	2/1276 (0.2%)
1	B	0.15	0/951	0.40	0/1291
1	C	0.19	0/976	0.48	0/1322
1	D	0.27	0/944	0.60	1/1283 (0.1%)
2	I	0.98	1/89 (1.1%)	1.46	1/116 (0.9%)
2	J	0.98	1/89 (1.1%)	1.42	1/116 (0.9%)
2	K	0.99	1/89 (1.1%)	1.37	1/116 (0.9%)
2	L	0.98	1/89 (1.1%)	1.53	1/116 (0.9%)
All	All	0.35	4/4169 (0.1%)	0.63	7/5636 (0.1%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	J	6	ASP	CG-OD2	5.16	1.35	1.25
2	I	6	ASP	CG-OD2	5.14	1.35	1.25
2	K	6	ASP	CG-OD2	5.04	1.34	1.25
2	L	6	ASP	CG-OD2	5.03	1.34	1.25

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	I	2	PHE	CA-CB-CG	-6.91	106.89	113.80
2	L	2	PHE	CA-CB-CG	-6.70	107.10	113.80
2	J	2	PHE	CA-CB-CG	-6.21	107.59	113.80
2	K	2	PHE	CA-CB-CG	-6.07	107.73	113.80
1	A	60	GLU	N-CA-C	-5.37	99.01	108.08
1	D	60	GLU	CB-CA-C	-5.11	110.70	116.63
1	A	60	GLU	CB-CA-C	-5.08	110.73	116.63

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	925	0	908	11	0
1	B	935	0	902	8	0
1	C	959	0	936	16	0
1	D	928	0	890	12	0
2	I	134	0	113	7	0
2	J	134	0	113	6	0
2	K	134	0	112	12	0
2	L	134	0	113	7	0
3	A	58	0	0	1	0
3	B	62	0	0	2	0
3	C	63	0	0	2	0
3	D	72	0	0	0	0
3	I	7	0	0	1	0
3	J	7	0	0	0	0
3	K	5	0	0	0	0
3	L	4	0	0	0	0
All	All	4561	0	4087	67	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 8.

All (67) 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:115:MET:SD	3:B:253:HOH:O	2.40	0.79
1:C:36:MET:HE2	1:C:101:ILE:HD12	1.68	0.75
1:B:115:MET:HE1	2:I:12:TYR:HB3	1.69	0.73
1:D:115:MET:HE1	2:L:12:TYR:HB3	1.75	0.68
1:C:115:MET:HE3	2:K:8:MEA:HB1	1.79	0.64
1:A:115:MET:HE1	2:J:12:TYR:HB3	1.81	0.63
1:B:21:VAL:HG23	1:B:122:ASP:HB3	1.80	0.62
2:L:13:LEU:O	2:L:15:LYS:N	2.33	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:I:13:LEU:O	2:I:15:LYS:N	2.32	0.59
2:I:16:NH2:N	3:I:103:HOH:O	2.36	0.58
1:C:115:MET:CE	2:K:8:MEA:HB1	2.35	0.56
1:C:54:ILE:HB	1:C:117:SER:HB3	1.88	0.56
1:D:59:MET:SD	1:D:60:GLU:HG3	2.47	0.55
1:D:36:MET:O	1:D:100:GLN:HA	2.07	0.55
1:C:115:MET:CE	2:K:8:MEA:CB	2.86	0.54
1:D:115:MET:HE2	2:L:8:MEA:CB	2.38	0.53
1:A:124:LYS:NZ	3:A:207:HOH:O	2.42	0.52
1:C:72:GLU:OE1	1:C:74:LEU:HD21	2.09	0.52
2:K:2:PHE:HE1	2:K:8:MEA:HC2	1.74	0.52
1:A:36:MET:HE1	1:A:128:VAL:HG11	1.91	0.51
1:C:64:ILE:O	1:C:65:ILE:HG13	2.10	0.51
1:A:85:ALA:HB2	1:A:101:ILE:HG12	1.94	0.50
1:C:115:MET:HE2	2:K:8:MEA:CB	2.42	0.49
1:A:65:ILE:HG21	1:A:87:LEU:HB2	1.94	0.49
2:L:9:TYR:HA	2:L:10:SAR:HN1	1.61	0.48
1:C:43:PRO:O	3:C:201:HOH:O	2.20	0.47
1:A:67:PHE:CD2	1:A:87:LEU:HD23	2.49	0.47
1:B:115:MET:HG3	1:B:123:TYR:HB2	1.96	0.47
2:K:9:TYR:HA	2:K:10:SAR:HN1	1.62	0.47
1:D:115:MET:HE2	2:L:8:MEA:HB2	1.97	0.46
1:B:115:MET:HG3	1:B:123:TYR:CB	2.45	0.46
1:B:66:GLN:OE1	1:B:76:THR:HG21	2.15	0.46
1:D:78:HIS:HB3	1:D:81:TYR:HD2	1.81	0.45
1:A:59:MET:SD	1:A:60:GLU:HG3	2.56	0.45
1:D:36:MET:HE2	1:D:104:VAL:HG22	1.98	0.45
2:I:2:PHE:HA	2:I:3:MEA:HC1	1.71	0.45
2:K:4:9KK:O	2:K:5:SAR:C	2.64	0.45
2:L:8:MEA:HC3	2:L:12:TYR:HE1	1.81	0.45
1:D:82:ARG:O	1:D:83:GLN:HB2	2.16	0.44
2:I:8:MEA:HC3	2:I:12:TYR:HE1	1.81	0.44
1:A:109:ALA:HB2	1:A:130:VAL:HG23	1.99	0.44
1:B:122:ASP:OD1	3:B:201:HOH:O	2.21	0.44
1:C:21:VAL:HG23	1:C:122:ASP:HB3	1.99	0.44
1:C:82:ARG:O	1:C:83:GLN:HB2	2.17	0.44
1:C:115:MET:HE1	2:K:12:TYR:HB3	2.00	0.44
1:D:78:HIS:HB3	1:D:81:TYR:CD2	2.52	0.44
1:A:59:MET:O	1:A:60:GLU:HB2	2.17	0.44
2:L:2:PHE:N	2:L:14:CYS:HB3	2.33	0.44
1:C:86:ARG:NH2	3:C:208:HOH:O	2.50	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:J:9:TYR:HA	2:J:10:SAR:HN1	1.61	0.43
1:D:65:ILE:HG21	1:D:87:LEU:HB2	1.99	0.43
1:A:21:VAL:HG23	1:A:122:ASP:HB3	2.01	0.43
1:C:115:MET:HE2	2:K:8:MEA:CG	2.48	0.43
1:A:38:ILE:HD11	1:A:99:LEU:HD23	2.01	0.43
1:C:36:MET:HE1	1:C:112:TYR:CZ	2.53	0.43
2:J:8:MEA:HD1	2:J:11:TRP:CZ3	2.55	0.42
1:D:57:TRP:CH2	1:D:114:CYS:HB3	2.55	0.42
1:C:115:MET:HE3	2:K:11:TRP:HB2	2.02	0.41
1:D:56:TYR:CZ	1:D:58:GLU:HB2	2.56	0.41
2:J:8:MEA:HC3	2:J:12:TYR:HE1	1.84	0.41
2:J:15:LYS:HZ3	2:J:15:LYS:HG2	1.77	0.41
1:B:78:HIS:ND1	1:B:80:SER:OG	2.50	0.41
2:I:2:PHE:N	2:I:14:CYS:HB3	2.36	0.41
2:K:2:PHE:CE1	2:K:8:MEA:HC2	2.55	0.41
2:I:9:TYR:HA	2:I:10:SAR:HN1	1.71	0.40
2:J:2:PHE:HA	2:J:3:MEA:HC1	1.75	0.40
2:K:8:MEA:HC3	2:K:12:TYR:HE1	1.86	0.40

There are no symmetry-related clashes.

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	117/127 (92%)	111 (95%)	6 (5%)	0	100	100
1	B	117/127 (92%)	112 (96%)	5 (4%)	0	100	100
1	C	119/127 (94%)	114 (96%)	5 (4%)	0	100	100
1	D	117/127 (92%)	112 (96%)	5 (4%)	0	100	100
2	I	9/16 (56%)	6 (67%)	1 (11%)	2 (22%)	0	0
2	J	9/16 (56%)	8 (89%)	0	1 (11%)	0	0

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	K	9/16 (56%)	7 (78%)	2 (22%)	0	100	100
2	L	9/16 (56%)	7 (78%)	0	2 (22%)	0	0
All	All	506/572 (88%)	477 (94%)	24 (5%)	5 (1%)	12	4

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	L	7	VAL
2	L	14	CYS
2	I	14	CYS
2	I	7	VAL
2	J	7	VAL

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	97/109 (89%)	97 (100%)	0	100	100
1	B	96/109 (88%)	95 (99%)	1 (1%)	68	64
1	C	99/109 (91%)	99 (100%)	0	100	100
1	D	94/109 (86%)	94 (100%)	0	100	100
2	I	9/9 (100%)	7 (78%)	2 (22%)	1	0
2	J	9/9 (100%)	8 (89%)	1 (11%)	6	0
2	K	9/9 (100%)	8 (89%)	1 (11%)	6	0
2	L	9/9 (100%)	7 (78%)	2 (22%)	1	0
All	All	422/472 (89%)	415 (98%)	7 (2%)	53	41

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	45	GLU
2	I	6	ASP

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Mol	Chain	Res	Type
2	I	15	LYS
2	J	15	LYS
2	K	15	LYS
2	L	13	LEU
2	L	15	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	77	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

20 non-standard protein/DNA/RNA residues are modelled in this entry.

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	$\# Z > 2$
2	SAR	L	5	2	3,4,5	0.90	0	1,3,5	1.09	0
2	MEA	K	8	2	11,12,13	0.96	0	13,14,16	1.37	1 (7%)
2	SAR	J	10	2	3,4,5	0.84	0	1,3,5	1.91	0
2	9KK	J	4	2	7,8,9	1.04	1 (14%)	5,8,10	1.28	1 (20%)
2	MEA	J	8	2	11,12,13	0.99	0	13,14,16	1.67	1 (7%)
2	MEA	K	3	2	11,12,13	1.55	2 (18%)	13,14,16	3.31	5 (38%)
2	SAR	L	10	2	3,4,5	0.80	0	1,3,5	1.73	0
2	SAR	K	10	2	3,4,5	0.86	0	1,3,5	1.69	0
2	SAR	K	5	2	3,4,5	0.76	0	1,3,5	1.71	0
2	9KK	K	4	2	7,8,9	1.74	1 (14%)	5,8,10	1.79	1 (20%)
2	SAR	I	10	2	3,4,5	0.80	0	1,3,5	1.80	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	9KK	L	4	2	7,8,9	1.06	1 (14%)	5,8,10	1.47	1 (20%)
2	SAR	J	5	2	3,4,5	0.87	0	1,3,5	1.49	0
2	MEA	J	3	2	11,12,13	1.56	1 (9%)	13,14,16	2.40	4 (30%)
2	MEA	L	3	2	11,12,13	1.60	1 (9%)	13,14,16	2.44	5 (38%)
2	SAR	I	5	2	3,4,5	0.88	0	1,3,5	1.53	0
2	9KK	I	4	2	7,8,9	1.10	1 (14%)	5,8,10	1.34	1 (20%)
2	MEA	L	8	2	11,12,13	0.91	0	13,14,16	0.96	1 (7%)
2	MEA	I	3	2	11,12,13	1.55	1 (9%)	13,14,16	2.44	5 (38%)
2	MEA	I	8	2	11,12,13	0.94	0	13,14,16	1.46	1 (7%)

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
2	SAR	L	5	2	-	1/1/2/3	-
2	MEA	K	8	2	-	1/5/8/10	0/1/1/1
2	SAR	J	10	2	-	1/1/2/3	-
2	9KK	J	4	2	-	2/5/8/10	-
2	MEA	J	8	2	-	0/5/8/10	0/1/1/1
2	MEA	K	3	2	-	0/5/8/10	0/1/1/1
2	SAR	L	10	2	-	1/1/2/3	-
2	SAR	K	10	2	-	1/1/2/3	-
2	SAR	K	5	2	-	1/1/2/3	-
2	9KK	K	4	2	-	4/5/8/10	-
2	SAR	I	10	2	-	1/1/2/3	-
2	9KK	L	4	2	-	3/5/8/10	-
2	SAR	J	5	2	-	1/1/2/3	-
2	MEA	J	3	2	-	0/5/8/10	0/1/1/1
2	MEA	L	3	2	-	0/5/8/10	0/1/1/1
2	SAR	I	5	2	-	1/1/2/3	-
2	9KK	I	4	2	-	2/5/8/10	-
2	MEA	L	8	2	-	0/5/8/10	0/1/1/1
2	MEA	I	3	2	-	1/5/8/10	0/1/1/1
2	MEA	I	8	2	-	0/5/8/10	0/1/1/1

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	K	4	9KK	CA-N	3.96	1.54	1.47
2	L	3	MEA	CA-N	3.89	1.54	1.47
2	I	3	MEA	CA-N	3.64	1.53	1.47
2	J	3	MEA	CA-N	3.61	1.53	1.47
2	K	3	MEA	CA-N	3.17	1.52	1.47
2	K	3	MEA	CB-CA	2.22	1.59	1.54
2	I	4	9KK	CA-N	2.19	1.51	1.47
2	L	4	9KK	CA-N	2.13	1.51	1.47
2	J	4	9KK	CA-N	2.09	1.51	1.47

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	K	3	MEA	CB-CA-C	6.94	125.60	111.81
2	K	3	MEA	O-C-CA	-6.26	108.67	124.77
2	J	8	MEA	CG-CB-CA	5.31	121.08	113.51
2	K	3	MEA	CB-CA-N	-5.16	102.88	110.48
2	L	3	MEA	O-C-CA	-4.87	112.25	124.77
2	J	3	MEA	CB-CA-C	4.85	121.45	111.81
2	I	3	MEA	O-C-CA	-4.84	112.33	124.77
2	I	3	MEA	CB-CA-C	4.68	121.11	111.81
2	J	3	MEA	O-C-CA	-4.61	112.92	124.77
2	K	8	MEA	CG-CB-CA	4.26	119.58	113.51
2	L	3	MEA	CB-CA-C	4.26	120.28	111.81
2	K	3	MEA	C1-N-CA	4.18	126.19	113.70
2	I	8	MEA	CG-CB-CA	4.11	119.36	113.51
2	L	3	MEA	CG-CB-CA	-4.02	107.80	113.51
2	I	3	MEA	CG-CB-CA	-3.60	108.39	113.51
2	J	3	MEA	C1-N-CA	3.53	124.25	113.70
2	J	3	MEA	CG-CB-CA	-3.48	108.56	113.51
2	I	3	MEA	C1-N-CA	3.48	124.10	113.70
2	L	3	MEA	C1-N-CA	3.46	124.04	113.70
2	K	4	9KK	O-C-CA	-3.31	116.27	124.77
2	L	4	9KK	O-C-CA	-2.99	117.08	124.77
2	I	4	9KK	O-C-CA	-2.80	117.58	124.77
2	J	4	9KK	O-C-CA	-2.68	117.87	124.77
2	L	8	MEA	CG-CB-CA	2.45	117.00	113.51
2	I	3	MEA	CB-CA-N	-2.33	107.05	110.48
2	L	3	MEA	CB-CA-N	-2.24	107.18	110.48
2	K	3	MEA	CB-CG-CD2	2.08	124.78	120.90

There are no chirality outliers.

All (21) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	I	3	MEA	O-C-CA-CB
2	K	4	9KK	C-CA-CB-CG
2	I	5	SAR	C-CA-N-CN
2	J	5	SAR	C-CA-N-CN
2	K	5	SAR	C-CA-N-CN
2	L	5	SAR	C-CA-N-CN
2	K	8	MEA	O-C-CA-CB
2	I	10	SAR	C-CA-N-CN
2	J	10	SAR	C-CA-N-CN
2	K	10	SAR	C-CA-N-CN
2	L	10	SAR	C-CA-N-CN
2	K	4	9KK	N-CA-CB-CG
2	I	4	9KK	N-CA-CB-CG
2	L	4	9KK	CE-CD-CG-CB
2	K	4	9KK	CA-CB-CG-CD
2	I	4	9KK	CA-CB-CG-CD
2	J	4	9KK	CA-CB-CG-CD
2	L	4	9KK	CA-CB-CG-CD
2	L	4	9KK	N-CA-CB-CG
2	J	4	9KK	CE-CD-CG-CB
2	K	4	9KK	CE-CD-CG-CB

There are no ring outliers.

12 monomers are involved in 21 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	K	8	MEA	8	0
2	J	10	SAR	1	0
2	J	8	MEA	2	0
2	L	10	SAR	1	0
2	K	10	SAR	1	0
2	K	5	SAR	1	0
2	K	4	9KK	1	0
2	I	10	SAR	1	0
2	J	3	MEA	1	0
2	L	8	MEA	3	0
2	I	3	MEA	1	0
2	I	8	MEA	1	0

5.5 Carbohydrates

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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	D	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	D	44:VAL	C	47:GLU	N	6.24

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	119/127 (93%)	-1.03	0 100 100	15, 23, 38, 52	0
1	B	121/127 (95%)	-1.05	0 100 100	12, 22, 33, 48	0
1	C	123/127 (96%)	-1.03	0 100 100	16, 22, 39, 46	0
1	D	121/127 (95%)	-1.05	0 100 100	16, 23, 41, 46	0
2	I	9/16 (56%)	-0.99	0 100 100	11, 17, 26, 34	0
2	J	9/16 (56%)	-1.00	0 100 100	11, 21, 30, 41	0
2	K	9/16 (56%)	-1.05	0 100 100	14, 19, 26, 31	0
2	L	9/16 (56%)	-0.99	0 100 100	15, 21, 30, 41	0
All	All	520/572 (90%)	-1.04	0 100 100	11, 23, 39, 52	0

There are no RSRZ outliers to report.

6.2 Non-standard residues in protein, DNA, RNA chains [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(Å ²)	Q<0.9
2	MEA	L	3	12/13	0.98	0.04	23,33,37,37	0
2	SAR	I	5	5/6	0.98	0.04	18,19,20,21	0
2	SAR	J	5	5/6	0.98	0.04	21,23,24,24	0
2	SAR	K	5	5/6	0.98	0.04	12,16,17,19	0
2	SAR	L	5	5/6	0.98	0.06	23,23,28,29	0
2	MEA	L	8	12/13	0.98	0.04	7,14,21,22	0
2	SAR	L	10	5/6	0.98	0.05	16,16,17,19	0
2	9KK	L	4	9/10	0.99	0.05	22,26,32,32	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	MEA	J	3	12/13	0.99	0.04	24,27,28,28	0
2	MEA	K	3	12/13	0.99	0.04	20,24,31,33	0
2	MEA	I	3	12/13	0.99	0.03	23,26,32,34	0
2	9KK	I	4	9/10	0.99	0.05	23,28,34,36	0
2	MEA	I	8	12/13	0.99	0.03	10,16,22,23	0
2	MEA	J	8	12/13	0.99	0.03	14,16,21,22	0
2	MEA	K	8	12/13	0.99	0.03	7,13,17,21	0
2	9KK	J	4	9/10	0.99	0.04	27,29,33,34	0
2	SAR	I	10	5/6	0.99	0.04	14,15,16,17	0
2	SAR	J	10	5/6	0.99	0.03	15,15,19,19	0
2	SAR	K	10	5/6	0.99	0.04	19,20,20,21	0
2	9KK	K	4	9/10	0.99	0.06	21,27,33,37	0

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.