



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:38 AM UTC

PDB ID : 1RGB / pdb_00001rgb
Title : Phospholipase A2 from Vipera ammodytes meridionalis
Authors : Georgieva, D.N.
Deposited on : 2003-11-12
Resolution : 3.30 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)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

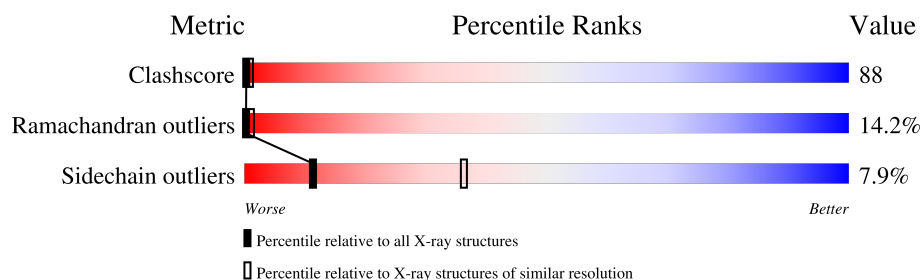
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.30 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))
Clashscore	190562	1209 (3.32-3.28)
Ramachandran outliers	187476	1188 (3.32-3.28)
Sidechain outliers	187428	1187 (3.32-3.28)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	122	
1	B	122	
1	K	122	
1	L	122	

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	ELD	B	134	-	-	X	-
2	ELD	L	134	-	-	X	-

2 Entry composition [i](#)

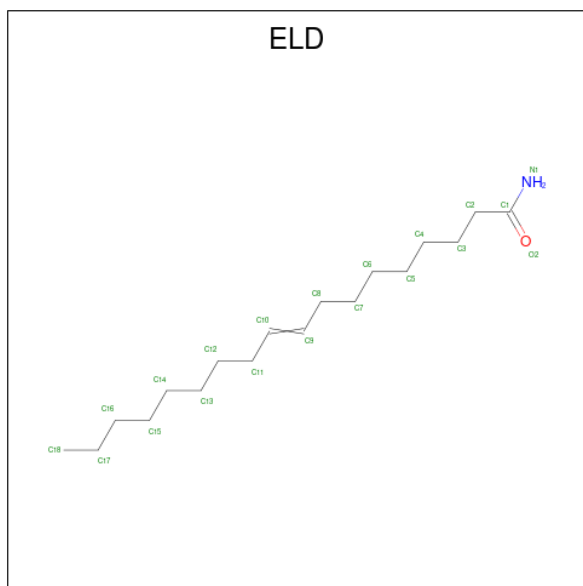
There are 2 unique types of molecules in this entry. The entry contains 3892 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Phospholipase A2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	122	Total	C	N	O	S	84	0	0
			963	598	176	174	15			
1	B	122	Total	C	N	O	S	65	0	0
			963	598	176	174	15			
1	K	122	Total	C	N	O	S	87	0	0
			963	598	176	174	15			
1	L	122	Total	C	N	O	S	81	0	0
			963	598	176	174	15			

- Molecule 2 is (9E)-OCTADEC-9-ENAMIDE (CCD ID: ELD) (formula: C₁₈H₃₅NO).



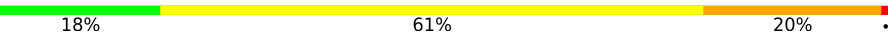
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	B	1	Total	C	N	O	0	0
			20	18	1	1		
2	L	1	Total	C	N	O	0	0
			20	18	1	1		

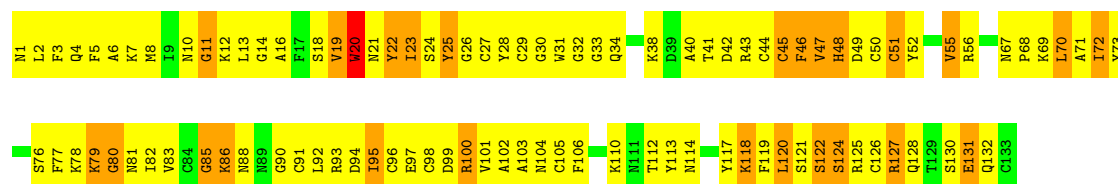
3 Residue-property plots

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. 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. 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.

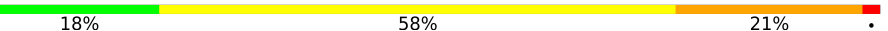
Note EDS was not executed.

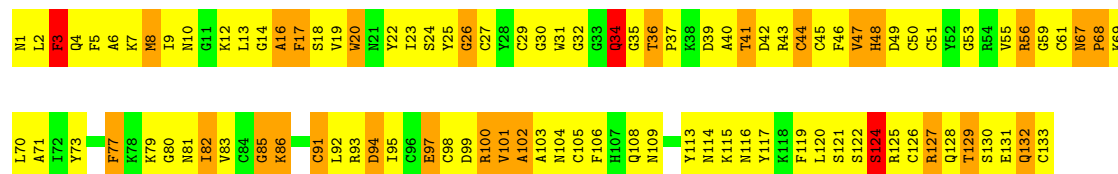
• Molecule 1: Phospholipase A2

Chain A: 

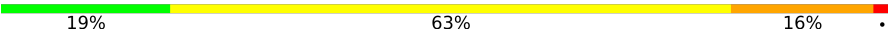


• Molecule 1: Phospholipase A2

Chain B: 

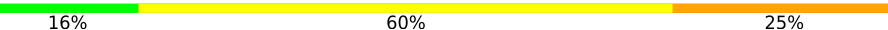


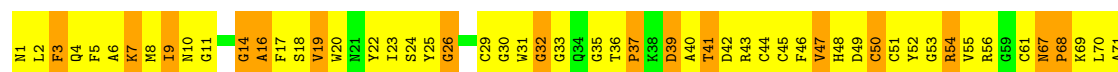
• Molecule 1: Phospholipase A2

Chain K: 



• Molecule 1: Phospholipase A2

Chain L: 



I72	Y73	S74	Y75	S76	F77		G80	N81	I82	V83	G84	G85	K86	N88	N89	G90	C91	L92	R93	D94	I95	C96	E97	G98	D99	R100	V101	A102	A103	N104	C105	F106	H107	Q108	N109	K110	N111	T112	Y113	N114	K115	N116	Y117	K118	F119	L120	S121	S122	S124	R125	C126	R127	Q128	T129	S130		C133
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4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	46.57 Å 82.67 Å 119.47 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.53 – 3.30	Depositor
% Data completeness (in resolution range)	94.5 (19.53-3.30)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.234 , 0.286	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3892	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ELD

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	0.61	0/985	1.11	10/1320 (0.8%)
1	B	0.63	0/985	1.18	14/1320 (1.1%)
1	K	0.58	0/985	1.05	10/1320 (0.8%)
1	L	0.56	0/985	1.07	9/1320 (0.7%)
All	All	0.60	0/3940	1.10	43/5280 (0.8%)

There are no bond length outliers.

All (43) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	122	SER	N-CA-C	-8.13	102.26	113.56
1	A	23	ILE	N-CA-C	-7.91	102.55	110.62
1	A	85	GLY	N-CA-C	7.79	122.20	112.14
1	B	44	CYS	N-CA-C	-7.75	101.81	111.11
1	A	120	LEU	N-CA-C	-7.19	98.37	109.52
1	A	48	HIS	N-CA-C	-6.87	102.64	111.02
1	A	22	TYR	N-CA-C	-6.86	102.35	113.19
1	B	3	PHE	N-CA-C	-6.81	103.94	111.36
1	L	80	GLY	N-CA-C	6.74	123.01	114.25
1	B	93	ARG	N-CA-C	-6.66	103.12	111.11
1	L	50	CYS	N-CA-C	-6.62	103.17	111.11
1	L	98	CYS	N-CA-C	-6.57	103.00	111.02
1	B	132	GLN	N-CA-C	6.42	119.08	110.35
1	K	48	HIS	N-CA-C	-6.36	104.04	110.97
1	B	91	CYS	N-CA-C	-6.27	104.50	111.71
1	K	21	ASN	N-CA-C	-6.26	104.75	112.38
1	B	82	ILE	N-CA-C	6.22	117.77	109.37
1	B	41	THR	N-CA-C	-6.19	104.44	111.07
1	B	73	TYR	N-CA-C	6.19	117.97	109.18
1	K	36	THR	CA-C-N	6.16	127.54	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	K	36	THR	C-N-CA	6.16	127.54	119.84
1	B	34	GLN	N-CA-C	6.13	116.53	107.88
1	L	92	LEU	N-CA-C	5.92	117.42	110.97
1	K	41	THR	N-CA-C	-5.88	104.79	111.14
1	L	67	ASN	N-CA-C	5.87	118.04	109.71
1	K	44	CYS	N-CA-C	-5.86	103.87	111.02
1	K	127	ARG	N-CA-C	-5.85	105.90	113.16
1	L	70	LEU	N-CA-C	5.79	120.63	113.50
1	A	70	LEU	N-CA-C	5.68	119.91	112.92
1	L	103	ALA	N-CA-C	-5.54	105.24	111.28
1	B	8	MET	N-CA-C	-5.50	105.38	111.71
1	L	126	CYS	N-CA-C	5.48	117.02	110.44
1	A	124	SER	N-CA-C	-5.48	106.95	113.97
1	A	51	CYS	N-CA-C	-5.41	105.08	110.97
1	K	40	ALA	N-CA-C	5.38	119.01	112.23
1	B	56	ARG	N-CA-C	5.36	117.47	109.59
1	L	102	ALA	N-CA-C	-5.34	105.10	111.03
1	K	35	GLY	N-CA-C	5.33	117.75	111.35
1	B	67	ASN	CA-C-N	5.14	126.27	119.84
1	B	67	ASN	C-N-CA	5.14	126.27	119.84
1	A	118	LYS	N-CA-C	-5.06	100.91	108.96
1	B	59	GLY	N-CA-C	5.03	125.11	113.18
1	K	61	CYS	N-CA-C	5.01	116.66	108.55

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	963	0	904	159	0
1	B	963	0	904	161	0
1	K	963	0	904	146	0
1	L	963	0	904	168	0
2	B	20	0	35	12	0
2	L	20	0	35	16	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	3892	0	3686	597	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 88.

All (597) 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:L:114:ASN:HD22	1:L:116:ASN:HB3	1.14	1.11
1:B:124:SER:HB3	1:B:127:ARG:HH21	1.20	1.06
1:A:25:TYR:O	1:A:29:CYS:HB2	1.52	1.05
1:L:20:TRP:HA	1:L:23:ILE:HG12	1.34	1.04
1:B:82:ILE:HG21	1:B:100:ARG:HG3	1.41	1.01
1:K:2:LEU:HD12	1:L:31:TRP:HA	1.45	0.99
1:K:69:LYS:HG2	1:L:49:ASP:OD1	1.66	0.95
1:L:41:THR:O	1:L:44:CYS:HB2	1.67	0.95
1:A:11:GLY:HA3	1:A:77:PHE:CZ	2.02	0.95
1:B:25:TYR:HB3	1:B:29:CYS:HB2	1.52	0.92
1:L:101:VAL:HA	1:L:104:ASN:HB2	1.53	0.91
1:K:1:ASN:HB2	1:L:33:GLY:HA3	1.49	0.91
1:K:43:ARG:O	1:K:47:VAL:HG23	1.68	0.90
1:A:69:LYS:HG3	1:A:70:LEU:HD12	1.54	0.90
1:L:114:ASN:ND2	1:L:116:ASN:HB3	1.88	0.88
1:L:5:PHE:CE1	1:L:99:ASP:HA	2.09	0.87
1:L:25:TYR:O	1:L:29:CYS:HB2	1.75	0.86
1:L:45:CYS:SG	2:L:134:ELD:H151	2.16	0.85
1:K:35:GLY:H	1:K:130:SER:HA	1.41	0.85
1:K:1:ASN:HD21	1:K:4:GLN:HG3	1.39	0.85
1:L:128:GLN:CD	1:L:128:GLN:H	1.82	0.85
1:K:118:LYS:HG3	1:K:119:PHE:CD2	2.13	0.84
1:A:24:SER:HA	1:A:30:GLY:HA3	1.59	0.83
1:K:1:ASN:CB	1:L:33:GLY:HA3	2.08	0.83
1:A:23:ILE:HG12	1:B:23:ILE:HD11	1.62	0.81
1:B:25:TYR:HB3	1:B:29:CYS:CB	2.11	0.81
1:L:5:PHE:HD1	1:L:99:ASP:CG	1.88	0.81
1:A:131:GLU:HG3	1:A:132:GLN:H	1.47	0.80
1:A:11:GLY:HA3	1:A:77:PHE:HZ	1.43	0.80
1:A:122:SER:HA	1:B:3:PHE:CZ	2.17	0.79
1:B:8:MET:SD	1:B:99:ASP:HB3	2.21	0.79
1:L:19:VAL:HG13	1:L:20:TRP:H	1.47	0.79
1:A:100:ARG:O	1:A:103:ALA:HB3	1.83	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:124:SER:HB3	1:B:127:ARG:NH2	1.96	0.78
1:K:25:TYR:O	1:K:29:CYS:HB2	1.82	0.78
1:K:100:ARG:NH2	1:K:101:VAL:HG22	1.98	0.78
1:B:117:TYR:C	1:B:120:LEU:HD21	2.09	0.77
1:A:101:VAL:HA	1:A:104:ASN:HD22	1.50	0.77
1:A:41:THR:O	1:A:44:CYS:HB2	1.85	0.76
1:L:48:HIS:O	1:L:51:CYS:HB3	1.85	0.76
1:A:120:LEU:HB3	1:A:125:ARG:HB2	1.67	0.76
1:B:82:ILE:CG2	1:B:100:ARG:HG3	2.16	0.76
1:B:19:VAL:HG13	1:B:20:TRP:CE3	2.20	0.76
1:A:48:HIS:HB3	2:B:134:ELD:HN11	1.51	0.75
1:L:20:TRP:HA	1:L:23:ILE:CG1	2.15	0.75
1:A:122:SER:HA	1:B:3:PHE:CE1	2.21	0.75
1:L:36:THR:HG23	1:L:37:PRO:HD2	1.67	0.75
1:A:90:GLY:O	1:A:93:ARG:HB3	1.86	0.74
1:A:1:ASN:HB3	1:A:71:ALA:O	1.88	0.74
1:B:45:CYS:SG	2:B:134:ELD:H172	2.28	0.74
1:L:5:PHE:HE1	1:L:99:ASP:HA	1.51	0.74
1:B:23:ILE:O	1:B:30:GLY:HA3	1.87	0.74
1:L:94:ASP:O	1:L:97:GLU:HB2	1.88	0.74
1:K:49:ASP:OD1	1:L:69:LYS:HG2	1.88	0.74
1:K:29:CYS:SG	1:K:42:ASP:HA	2.28	0.74
1:K:3:PHE:CE1	1:L:32:GLY:HA2	2.23	0.73
1:B:103:ALA:O	1:B:106:PHE:HB2	1.88	0.73
1:L:18:SER:HB3	1:L:22:TYR:HE2	1.53	0.73
1:L:20:TRP:CA	1:L:23:ILE:HG12	2.17	0.73
1:B:94:ASP:O	1:B:97:GLU:HB2	1.87	0.73
1:K:91:CYS:C	1:K:93:ARG:H	1.95	0.73
1:A:25:TYR:HD2	1:A:118:LYS:HG2	1.52	0.73
1:B:102:ALA:HB1	2:B:134:ELD:H183	1.70	0.73
1:B:132:GLN:HG2	1:B:133:CYS:N	2.04	0.73
1:K:45:CYS:O	1:K:48:HIS:HB3	1.88	0.72
1:L:91:CYS:SG	1:L:92:LEU:N	2.63	0.72
1:L:118:LYS:HB3	1:L:119:PHE:CD2	2.24	0.72
1:B:47:VAL:O	1:B:50:CYS:N	2.22	0.72
1:A:49:ASP:OD1	1:B:69:LYS:HG2	1.90	0.72
1:K:72:ILE:O	1:K:92:LEU:HD11	1.89	0.72
1:B:45:CYS:HA	2:B:134:ELD:H172	1.72	0.71
1:A:43:ARG:O	1:A:47:VAL:HG23	1.90	0.71
1:A:8:MET:SD	1:A:82:ILE:HD13	2.31	0.71
1:A:28:TYR:HE2	1:A:33:GLY:C	1.99	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:26:GLY:HA3	1:L:42:ASP:CG	2.16	0.70
1:K:71:ALA:C	1:K:72:ILE:HG13	2.16	0.70
1:L:88:ASN:ND2	1:L:93:ARG:HB2	2.06	0.70
1:B:16:ALA:O	1:B:17:PHE:HB2	1.91	0.70
1:K:51:CYS:O	1:K:55:VAL:HG12	1.92	0.70
1:K:59:GLY:HA3	1:K:91:CYS:HB3	1.73	0.70
1:A:71:ALA:O	1:A:72:ILE:HG13	1.92	0.70
1:A:1:ASN:CG	1:A:4:GLN:HG3	2.17	0.69
1:A:26:GLY:HA3	1:A:42:ASP:OD1	1.92	0.69
1:B:41:THR:O	1:B:44:CYS:HB2	1.92	0.69
1:A:47:VAL:O	1:A:51:CYS:N	2.21	0.69
1:K:35:GLY:N	1:K:130:SER:HA	2.08	0.69
1:K:114:ASN:C	1:K:116:ASN:H	1.97	0.69
1:L:44:CYS:HB3	2:L:134:ELD:H181	1.75	0.69
1:K:27:CYS:O	1:K:33:GLY:HA2	1.92	0.68
1:L:54:ARG:HA	1:L:54:ARG:CZ	2.23	0.68
1:K:67:ASN:O	1:K:71:ALA:N	2.27	0.68
1:K:71:ALA:O	1:K:72:ILE:HG13	1.93	0.68
1:K:124:SER:O	1:K:127:ARG:HG2	1.93	0.68
1:A:102:ALA:O	1:A:105:CYS:HB3	1.93	0.68
1:K:91:CYS:O	1:K:93:ARG:N	2.25	0.68
1:A:28:TYR:HE2	1:A:34:GLN:N	1.92	0.68
1:B:29:CYS:SG	1:B:42:ASP:HA	2.34	0.68
1:B:122:SER:O	1:B:125:ARG:N	2.27	0.67
1:K:67:ASN:CG	1:K:70:LEU:HD22	2.19	0.67
1:B:5:PHE:CE1	1:B:99:ASP:HA	2.30	0.67
1:A:3:PHE:CE2	1:B:32:GLY:HA2	2.29	0.67
1:B:24:SER:HB3	1:B:119:PHE:CD1	2.29	0.67
1:L:19:VAL:HG13	1:L:20:TRP:N	2.08	0.67
1:L:47:VAL:O	1:L:48:HIS:C	2.38	0.67
1:A:126:CYS:O	1:A:128:GLN:N	2.29	0.66
1:A:82:ILE:HG22	1:A:100:ARG:HG3	1.76	0.66
1:B:48:HIS:CG	2:B:134:ELD:H161	2.29	0.66
1:B:26:GLY:HA3	1:B:42:ASP:CG	2.20	0.66
1:B:47:VAL:HG12	1:B:48:HIS:N	2.10	0.66
1:K:51:CYS:HA	1:K:54:ARG:HE	1.61	0.65
1:K:103:ALA:O	1:K:106:PHE:HB2	1.96	0.65
1:K:69:LYS:HG3	1:K:70:LEU:HD13	1.78	0.65
1:A:19:VAL:HG13	1:A:20:TRP:H	1.60	0.65
1:L:91:CYS:SG	1:L:92:LEU:HG	2.37	0.65
1:K:11:GLY:HA3	1:K:77:PHE:CZ	2.32	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:122:SER:C	1:B:125:ARG:H	2.04	0.65
1:K:1:ASN:ND2	1:K:4:GLN:HG3	2.11	0.65
1:L:101:VAL:CA	1:L:104:ASN:HB2	2.26	0.64
1:A:126:CYS:O	1:A:127:ARG:C	2.40	0.64
1:A:88:ASN:CG	1:A:93:ARG:HA	2.22	0.64
1:K:70:LEU:HD11	1:L:49:ASP:HB3	1.79	0.64
1:K:102:ALA:O	1:K:105:CYS:HB3	1.97	0.64
1:B:121:SER:O	1:B:125:ARG:HB2	1.98	0.64
1:K:3:PHE:HE1	1:L:32:GLY:HA2	1.63	0.64
1:B:25:TYR:O	1:B:26:GLY:O	2.16	0.64
1:L:67:ASN:OD1	1:L:67:ASN:C	2.40	0.63
1:K:10:ASN:HD21	1:K:19:VAL:CG1	2.12	0.63
1:A:38:LYS:HB3	1:A:117:TYR:CE1	2.34	0.63
1:B:5:PHE:HB2	1:B:99:ASP:OD2	1.99	0.63
1:L:29:CYS:SG	1:L:42:ASP:HA	2.39	0.63
1:L:77:PHE:HD2	1:L:81:ASN:O	1.82	0.63
1:A:16:ALA:C	1:A:18:SER:H	2.07	0.63
1:A:82:ILE:CG2	1:A:100:ARG:HG3	2.29	0.63
1:A:25:TYR:HA	1:A:118:LYS:HA	1.81	0.62
1:L:10:ASN:HB3	1:L:16:ALA:HA	1.81	0.62
1:L:46:PHE:CE1	1:L:50:CYS:SG	2.93	0.62
1:K:78:LYS:C	1:K:80:GLY:H	2.07	0.62
1:K:10:ASN:HA	1:K:14:GLY:O	1.99	0.62
1:L:88:ASN:ND2	1:L:90:GLY:H	1.97	0.62
1:K:37:PRO:O	1:K:39:ASP:N	2.32	0.62
1:L:67:ASN:O	1:L:71:ALA:N	2.33	0.62
1:B:45:CYS:O	1:B:46:PHE:C	2.43	0.62
1:K:24:SER:O	1:K:25:TYR:HB2	1.98	0.62
1:A:3:PHE:CD2	1:B:32:GLY:HA2	2.35	0.62
1:A:28:TYR:HD1	1:A:46:PHE:HB2	1.65	0.62
1:B:105:CYS:O	1:B:108:GLN:HB2	2.00	0.62
1:B:122:SER:O	1:B:126:CYS:N	2.32	0.61
1:L:56:ARG:HE	1:L:56:ARG:HA	1.65	0.61
1:L:98:CYS:O	1:L:102:ALA:CB	2.48	0.61
1:K:52:TYR:O	1:K:55:VAL:HG13	2.00	0.61
1:A:100:ARG:NH2	1:A:101:VAL:HG22	2.15	0.61
1:A:19:VAL:O	1:A:21:ASN:N	2.34	0.60
1:B:67:ASN:O	1:B:71:ALA:HB2	2.00	0.60
1:B:67:ASN:ND2	1:B:70:LEU:HD12	2.17	0.60
1:K:97:GLU:O	1:K:101:VAL:HG23	2.01	0.60
1:A:5:PHE:CD1	1:A:99:ASP:HB3	2.35	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:61:CYS:SG	1:L:95:ILE:HD12	2.42	0.60
1:L:92:LEU:O	1:L:93:ARG:C	2.43	0.60
1:B:41:THR:HG22	1:B:109:ASN:HB2	1.82	0.60
1:L:36:THR:CG2	1:L:37:PRO:HD2	2.31	0.60
1:L:25:TYR:C	1:L:29:CYS:HB2	2.26	0.60
1:A:70:LEU:HD13	1:B:49:ASP:HB3	1.83	0.59
1:K:74:SER:O	1:K:75:TYR:HB3	2.02	0.59
1:A:71:ALA:C	1:A:72:ILE:HG13	2.27	0.59
1:B:39:ASP:O	1:B:43:ARG:N	2.35	0.59
1:B:16:ALA:O	1:B:17:PHE:CB	2.50	0.59
1:L:5:PHE:CD1	1:L:99:ASP:HA	2.36	0.59
1:L:6:ALA:O	1:L:9:ILE:N	2.35	0.59
1:A:2:LEU:O	1:A:5:PHE:HB3	2.03	0.59
1:B:122:SER:C	1:B:125:ARG:N	2.60	0.59
1:L:16:ALA:O	1:L:17:PHE:HB2	2.02	0.59
1:B:1:ASN:N	1:B:4:GLN:OE1	2.35	0.59
1:K:17:PHE:CD2	1:K:20:TRP:HD1	2.21	0.59
1:L:51:CYS:O	1:L:55:VAL:HG23	2.01	0.59
1:L:126:CYS:HA	1:L:128:GLN:OE1	2.03	0.59
1:K:69:LYS:HG2	1:L:49:ASP:CG	2.29	0.58
1:L:128:GLN:H	1:L:128:GLN:NE2	2.00	0.58
1:L:46:PHE:O	1:L:47:VAL:C	2.45	0.58
1:K:2:LEU:CD1	1:L:31:TRP:HA	2.27	0.58
1:B:101:VAL:O	1:B:105:CYS:N	2.33	0.58
1:B:18:SER:O	1:B:22:TYR:HD2	1.87	0.58
1:B:129:THR:HG22	1:B:130:SER:O	2.03	0.58
1:A:118:LYS:O	1:A:119:PHE:HB2	2.04	0.58
1:B:91:CYS:SG	1:B:92:LEU:HD12	2.44	0.58
1:K:77:PHE:CE2	1:K:82:ILE:HG23	2.38	0.58
1:L:121:SER:O	1:L:122:SER:C	2.44	0.58
1:L:1:ASN:O	1:L:5:PHE:N	2.30	0.58
1:L:18:SER:O	1:L:19:VAL:C	2.46	0.58
1:B:45:CYS:HA	2:B:134:ELD:C17	2.34	0.57
1:A:1:ASN:OD1	1:A:3:PHE:HB2	2.04	0.57
1:A:94:ASP:O	1:A:97:GLU:N	2.37	0.57
1:L:114:ASN:HD22	1:L:116:ASN:CB	2.04	0.57
1:B:67:ASN:C	1:B:67:ASN:OD1	2.46	0.57
1:L:56:ARG:HA	1:L:56:ARG:NE	2.20	0.57
1:B:8:MET:SD	1:B:99:ASP:CB	2.91	0.57
1:B:8:MET:SD	1:B:100:ARG:N	2.78	0.57
1:K:114:ASN:O	1:K:116:ASN:N	2.37	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:110:LYS:O	1:K:113:TYR:HB2	2.04	0.57
1:A:46:PHE:O	1:A:49:ASP:HB2	2.05	0.57
1:B:114:ASN:O	1:B:117:TYR:HB2	2.04	0.56
1:K:1:ASN:C	1:K:3:PHE:N	2.60	0.56
1:B:23:ILE:HG23	1:B:24:SER:N	2.20	0.56
1:A:23:ILE:HG12	1:B:23:ILE:CD1	2.35	0.56
1:A:45:CYS:O	1:A:46:PHE:C	2.47	0.56
1:L:122:SER:O	1:L:124:SER:C	2.48	0.56
1:A:49:ASP:HB3	1:B:70:LEU:HD11	1.87	0.56
1:K:54:ARG:O	1:K:55:VAL:C	2.49	0.56
1:L:98:CYS:O	1:L:102:ALA:HB3	2.06	0.56
1:B:69:LYS:NZ	2:B:134:ELD:H22	2.20	0.55
1:K:111:ASN:C	1:K:113:TYR:H	2.14	0.55
1:B:18:SER:O	1:B:22:TYR:CD2	2.59	0.55
1:K:111:ASN:O	1:K:113:TYR:N	2.36	0.55
1:A:45:CYS:O	1:A:48:HIS:N	2.38	0.55
1:L:100:ARG:O	1:L:104:ASN:N	2.27	0.55
1:A:8:MET:HE3	1:A:99:ASP:HB3	1.88	0.55
1:B:102:ALA:O	1:B:103:ALA:C	2.48	0.55
1:A:49:ASP:CG	2:B:134:ELD:HN12	2.15	0.55
1:K:6:ALA:O	1:K:9:ILE:N	2.39	0.55
1:K:114:ASN:C	1:K:116:ASN:N	2.64	0.55
1:K:29:CYS:HA	2:L:134:ELD:O2	2.06	0.55
1:K:68:PRO:HG3	1:K:95:ILE:HD11	1.87	0.55
1:K:130:SER:O	1:K:131:GLU:HB3	2.06	0.55
1:A:4:GLN:HE22	1:A:73:TYR:N	2.05	0.55
1:B:46:PHE:O	1:B:47:VAL:C	2.50	0.55
1:K:3:PHE:CE2	1:L:122:SER:HA	2.42	0.55
1:L:128:GLN:CD	1:L:128:GLN:N	2.59	0.55
1:K:1:ASN:HD22	1:K:1:ASN:N	2.04	0.54
1:L:45:CYS:O	1:L:48:HIS:HB3	2.07	0.54
1:A:101:VAL:O	1:A:102:ALA:C	2.50	0.54
1:B:132:GLN:HG2	1:B:133:CYS:O	2.07	0.54
1:A:47:VAL:O	1:A:50:CYS:N	2.41	0.54
1:B:34:GLN:HG2	1:B:126:CYS:HB2	1.90	0.54
1:K:16:ALA:C	1:K:18:SER:H	2.15	0.54
1:B:117:TYR:CA	1:B:120:LEU:HD21	2.38	0.54
1:L:20:TRP:HE3	1:L:23:ILE:HG13	1.73	0.54
1:A:8:MET:SD	1:A:100:ARG:HA	2.48	0.54
1:L:6:ALA:O	1:L:7:LYS:C	2.49	0.54
1:K:105:CYS:O	1:K:109:ASN:OD1	2.25	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:103:ALA:HA	1:L:106:PHE:CD2	2.43	0.54
1:A:23:ILE:CG2	1:A:24:SER:N	2.71	0.53
1:A:10:ASN:HA	1:A:14:GLY:O	2.08	0.53
1:A:25:TYR:CD1	1:A:25:TYR:C	2.86	0.53
1:A:28:TYR:CE2	1:A:33:GLY:C	2.83	0.53
1:B:67:ASN:OD1	1:B:70:LEU:N	2.37	0.53
1:B:106:PHE:CZ	2:B:134:ELD:H171	2.44	0.53
1:B:115:LYS:C	1:B:117:TYR:H	2.16	0.53
1:L:39:ASP:OD2	1:L:39:ASP:C	2.52	0.53
1:A:70:LEU:CD1	1:B:49:ASP:HB3	2.39	0.53
1:A:94:ASP:O	1:A:95:ILE:C	2.50	0.53
1:B:99:ASP:O	1:B:103:ALA:N	2.41	0.53
1:A:97:GLU:O	1:A:100:ARG:HB3	2.08	0.53
1:K:1:ASN:O	1:K:3:PHE:N	2.42	0.53
1:K:97:GLU:HA	1:K:100:ARG:NH1	2.24	0.53
1:L:82:ILE:CG2	1:L:100:ARG:HG3	2.38	0.53
1:A:1:ASN:O	1:A:2:LEU:C	2.49	0.53
1:L:48:HIS:O	1:L:51:CYS:CB	2.55	0.53
1:A:92:LEU:O	1:A:93:ARG:C	2.51	0.53
1:L:96:CYS:O	1:L:100:ARG:N	2.35	0.53
1:A:48:HIS:HB3	2:B:134:ELD:N1	2.24	0.52
1:B:24:SER:HB3	1:B:119:PHE:HD1	1.73	0.52
1:K:94:ASP:O	1:K:97:GLU:HB2	2.09	0.52
1:A:101:VAL:O	1:A:105:CYS:N	2.28	0.52
1:A:19:VAL:C	1:A:21:ASN:N	2.67	0.52
1:A:85:GLY:O	1:A:86:LYS:C	2.51	0.52
1:B:27:CYS:HB3	1:B:34:GLN:O	2.08	0.52
1:K:95:ILE:O	1:K:96:CYS:C	2.51	0.52
1:L:88:ASN:HD22	1:L:93:ARG:HB2	1.74	0.52
1:L:94:ASP:O	1:L:97:GLU:N	2.43	0.52
1:L:96:CYS:O	1:L:99:ASP:HB2	2.10	0.52
1:L:97:GLU:O	1:L:98:CYS:C	2.51	0.52
1:K:3:PHE:CD1	1:L:32:GLY:HA2	2.44	0.52
1:L:6:ALA:C	1:L:8:MET:N	2.66	0.52
1:B:25:TYR:CD1	1:B:26:GLY:N	2.71	0.52
1:L:48:HIS:O	1:L:52:TYR:N	2.41	0.52
1:A:23:ILE:HD12	1:B:31:TRP:NE1	2.25	0.52
1:A:47:VAL:HA	1:A:50:CYS:HB2	1.92	0.52
1:L:20:TRP:O	1:L:23:ILE:HG12	2.10	0.52
1:A:22:TYR:OH	1:A:106:PHE:HB3	2.09	0.52
1:K:45:CYS:SG	1:K:106:PHE:HZ	2.33	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:32:GLY:HA2	1:B:3:PHE:CG	2.46	0.51
1:B:22:TYR:CD2	1:B:22:TYR:N	2.78	0.51
1:B:100:ARG:O	1:B:101:VAL:C	2.53	0.51
1:A:3:PHE:O	1:A:4:GLN:C	2.53	0.51
1:A:38:LYS:HB3	1:A:117:TYR:HE1	1.75	0.51
1:B:98:CYS:O	1:B:102:ALA:N	2.38	0.51
1:K:73:TYR:HA	1:K:92:LEU:HD11	1.91	0.51
1:B:9:ILE:HG22	1:B:10:ASN:N	2.24	0.51
1:K:122:SER:HA	1:L:3:PHE:CZ	2.45	0.51
1:A:13:LEU:O	1:A:18:SER:HB3	2.09	0.51
1:A:47:VAL:O	1:A:48:HIS:C	2.54	0.51
1:A:77:PHE:O	1:A:79:LYS:N	2.43	0.51
1:L:3:PHE:O	1:L:7:LYS:HG3	2.11	0.51
1:K:45:CYS:SG	1:K:106:PHE:CZ	3.04	0.51
1:K:91:CYS:C	1:K:93:ARG:N	2.64	0.51
1:A:55:VAL:CG1	1:A:56:ARG:N	2.74	0.50
1:B:6:ALA:O	1:B:9:ILE:HB	2.11	0.50
1:K:1:ASN:O	1:K:4:GLN:N	2.44	0.50
1:K:20:TRP:C	1:K:22:TYR:N	2.67	0.50
1:K:46:PHE:O	1:K:49:ASP:HB2	2.11	0.50
1:L:4:GLN:NE2	1:L:75:TYR:HE1	2.09	0.50
1:L:10:ASN:OD1	1:L:19:VAL:HG12	2.11	0.50
1:A:19:VAL:C	1:A:21:ASN:H	2.18	0.50
1:K:78:LYS:O	1:K:80:GLY:N	2.43	0.50
1:A:10:ASN:O	1:A:14:GLY:O	2.29	0.50
1:B:9:ILE:O	1:B:10:ASN:C	2.54	0.50
1:L:94:ASP:O	1:L:95:ILE:C	2.54	0.50
1:A:25:TYR:HD2	1:A:118:LYS:CG	2.22	0.50
1:B:47:VAL:O	1:B:48:HIS:C	2.53	0.50
1:B:34:GLN:HG3	1:B:35:GLY:N	2.26	0.50
1:K:35:GLY:CA	1:K:130:SER:HA	2.41	0.50
1:L:48:HIS:NE2	1:L:52:TYR:CZ	2.80	0.50
1:A:67:ASN:OD1	1:A:69:LYS:HB3	2.11	0.50
1:A:1:ASN:ND2	1:A:4:GLN:HG3	2.27	0.50
1:A:67:ASN:O	1:A:71:ALA:N	2.45	0.50
1:B:117:TYR:HA	1:B:120:LEU:HD21	1.94	0.50
1:A:8:MET:HE1	1:A:100:ARG:N	2.27	0.50
1:B:5:PHE:O	1:B:9:ILE:HG13	2.11	0.50
1:B:99:ASP:O	1:B:100:ARG:C	2.53	0.50
1:K:1:ASN:ND2	1:K:4:GLN:H	2.10	0.50
1:K:2:LEU:HB3	1:L:31:TRP:CE3	2.47	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:24:SER:OG	1:L:119:PHE:HA	2.12	0.49
1:B:4:GLN:O	1:B:5:PHE:C	2.55	0.49
1:B:25:TYR:CD1	1:B:42:ASP:HB2	2.46	0.49
1:A:5:PHE:O	1:A:8:MET:HB2	2.13	0.49
1:B:101:VAL:O	1:B:102:ALA:C	2.55	0.49
1:L:5:PHE:CE1	1:L:48:HIS:CD2	3.00	0.49
1:A:95:ILE:O	1:A:98:CYS:HB2	2.12	0.49
1:K:1:ASN:ND2	1:K:1:ASN:N	2.60	0.49
1:K:18:SER:O	1:K:21:ASN:N	2.40	0.49
1:A:97:GLU:CD	1:A:100:ARG:HE	2.19	0.49
1:L:30:GLY:H	2:L:134:ELD:H122	1.78	0.49
1:B:82:ILE:HD12	1:B:104:ASN:ND2	2.28	0.49
1:L:111:ASN:C	1:L:113:TYR:H	2.21	0.49
1:A:23:ILE:O	1:A:30:GLY:HA2	2.12	0.49
1:B:97:GLU:O	1:B:101:VAL:HG23	2.12	0.49
1:A:49:ASP:OD2	1:B:69:LYS:NZ	2.44	0.49
1:B:39:ASP:OD2	1:B:40:ALA:N	2.46	0.49
1:B:104:ASN:O	1:B:105:CYS:C	2.53	0.49
1:K:3:PHE:HE2	1:L:122:SER:HA	1.78	0.49
1:A:19:VAL:HG13	1:A:20:TRP:N	2.27	0.49
1:A:23:ILE:HG22	1:A:24:SER:N	2.27	0.49
1:B:4:GLN:O	1:B:7:LYS:N	2.46	0.49
1:L:68:PRO:O	1:L:69:LYS:C	2.53	0.49
1:A:2:LEU:O	1:A:3:PHE:C	2.56	0.48
1:L:40:ALA:O	1:L:43:ARG:HB2	2.13	0.48
1:L:113:TYR:CD1	1:L:113:TYR:C	2.91	0.48
1:A:2:LEU:O	1:A:6:ALA:N	2.30	0.48
1:A:4:GLN:NE2	1:A:73:TYR:O	2.47	0.48
1:L:73:TYR:HB3	1:L:92:LEU:HD22	1.95	0.48
1:L:102:ALA:O	1:L:105:CYS:HB3	2.13	0.48
1:A:43:ARG:O	1:A:44:CYS:C	2.55	0.48
1:K:37:PRO:C	1:K:39:ASP:N	2.72	0.48
1:A:121:SER:C	1:A:124:SER:H	2.21	0.48
1:B:69:LYS:HD2	1:B:69:LYS:O	2.14	0.48
1:K:2:LEU:HB2	1:L:31:TRP:CA	2.43	0.48
1:L:19:VAL:HG22	1:L:23:ILE:HD11	1.96	0.48
1:A:121:SER:C	1:A:124:SER:N	2.69	0.48
1:B:121:SER:O	1:B:125:ARG:N	2.41	0.48
1:K:24:SER:O	1:K:25:TYR:CB	2.62	0.48
1:K:42:ASP:O	1:K:45:CYS:HB2	2.14	0.48
1:L:4:GLN:HA	1:L:7:LYS:HG3	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:45:CYS:HA	2:L:134:ELD:H151	1.96	0.48
1:L:115:LYS:HA	1:L:118:LYS:HG3	1.96	0.48
1:K:1:ASN:C	1:K:3:PHE:H	2.21	0.48
1:A:3:PHE:O	1:A:6:ALA:HB3	2.14	0.48
1:B:36:THR:CG2	1:B:131:GLU:HB2	2.44	0.48
1:B:113:TYR:CD1	1:B:113:TYR:C	2.91	0.48
1:K:41:THR:O	1:K:44:CYS:HB2	2.13	0.48
1:A:110:LYS:O	1:A:113:TYR:HB2	2.14	0.47
1:B:9:ILE:HG12	1:B:106:PHE:HE2	1.79	0.47
1:B:101:VAL:O	1:B:104:ASN:HB2	2.12	0.47
1:B:121:SER:O	1:B:122:SER:C	2.56	0.47
1:L:3:PHE:O	1:L:7:LYS:CG	2.62	0.47
1:K:43:ARG:O	1:K:44:CYS:C	2.55	0.47
1:L:52:TYR:CD2	1:L:68:PRO:HB3	2.49	0.47
1:A:25:TYR:CD2	1:A:118:LYS:HG2	2.42	0.47
1:A:1:ASN:OD1	1:A:4:GLN:HG3	2.14	0.47
1:A:73:TYR:HB3	1:A:92:LEU:HB3	1.97	0.47
1:A:78:LYS:O	1:A:80:GLY:N	2.41	0.47
1:K:18:SER:O	1:K:19:VAL:C	2.58	0.47
1:A:91:CYS:O	1:A:92:LEU:C	2.57	0.47
1:A:121:SER:HB3	1:A:124:SER:HB3	1.96	0.47
1:B:4:GLN:C	1:B:6:ALA:N	2.70	0.47
1:B:29:CYS:O	1:B:30:GLY:C	2.58	0.47
1:K:70:LEU:HD11	1:L:49:ASP:CB	2.45	0.47
1:A:118:LYS:O	1:A:119:PHE:CB	2.62	0.47
1:A:49:ASP:CG	1:B:69:LYS:NZ	2.73	0.47
1:K:68:PRO:O	1:K:69:LYS:C	2.57	0.47
1:K:111:ASN:C	1:K:113:TYR:N	2.73	0.47
1:L:107:HIS:O	1:L:110:LYS:HG2	2.15	0.46
1:A:8:MET:HE3	1:A:99:ASP:CB	2.45	0.46
1:A:23:ILE:CG1	1:B:23:ILE:HD11	2.41	0.46
1:B:9:ILE:HG12	1:B:106:PHE:CE2	2.50	0.46
1:B:68:PRO:C	1:B:70:LEU:N	2.72	0.46
1:B:115:LYS:C	1:B:117:TYR:N	2.73	0.46
1:L:106:PHE:CD2	2:L:134:ELD:H183	2.51	0.46
1:L:106:PHE:CE2	2:L:134:ELD:H183	2.49	0.46
1:A:25:TYR:O	1:A:29:CYS:CB	2.44	0.46
1:K:8:MET:O	1:K:9:ILE:C	2.59	0.46
1:L:19:VAL:HA	1:L:22:TYR:HD2	1.80	0.46
1:B:91:CYS:HA	1:B:94:ASP:OD1	2.16	0.46
1:B:121:SER:O	1:B:125:ARG:CB	2.64	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:105:CYS:O	1:L:106:PHE:C	2.59	0.46
1:L:102:ALA:O	1:L:105:CYS:CB	2.64	0.46
1:L:118:LYS:HB3	1:L:119:PHE:CE2	2.49	0.46
1:A:78:LYS:C	1:A:80:GLY:H	2.24	0.46
1:B:1:ASN:OD1	1:B:2:LEU:N	2.48	0.46
1:A:1:ASN:N	1:A:4:GLN:OE1	2.46	0.46
1:A:31:TRP:CZ2	2:B:134:ELD:H121	2.51	0.46
1:K:5:PHE:O	1:K:8:MET:HB3	2.16	0.46
1:A:4:GLN:HE22	1:A:72:ILE:C	2.24	0.46
1:K:101:VAL:O	1:K:102:ALA:C	2.59	0.46
1:L:31:TRP:CE2	2:L:134:ELD:H62	2.51	0.46
1:B:26:GLY:HA3	1:B:42:ASP:OD1	2.16	0.46
1:B:104:ASN:C	1:B:106:PHE:N	2.69	0.46
1:L:46:PHE:HE1	1:L:50:CYS:SG	2.35	0.46
1:B:20:TRP:CE3	1:B:20:TRP:HA	2.52	0.46
1:B:97:GLU:OE1	1:B:97:GLU:HA	2.13	0.46
1:K:116:ASN:C	1:K:118:LYS:H	2.24	0.46
1:A:69:LYS:HG3	1:A:70:LEU:CD1	2.37	0.45
1:B:41:THR:CG2	1:B:109:ASN:HB2	2.45	0.45
1:K:16:ALA:C	1:K:18:SER:N	2.73	0.45
1:K:24:SER:HB3	1:K:119:PHE:CD1	2.51	0.45
1:L:26:GLY:HA3	1:L:42:ASP:OD1	2.16	0.45
1:A:1:ASN:OD1	1:A:4:GLN:N	2.42	0.45
1:A:5:PHE:O	1:A:6:ALA:C	2.58	0.45
1:A:23:ILE:O	1:A:30:GLY:CA	2.64	0.45
1:A:27:CYS:O	1:A:28:TYR:CD2	2.68	0.45
1:B:98:CYS:O	1:B:99:ASP:C	2.59	0.45
1:K:75:TYR:N	1:K:75:TYR:CD1	2.85	0.45
1:L:10:ASN:HD21	1:L:19:VAL:CG1	2.29	0.45
1:A:2:LEU:HB2	1:B:32:GLY:H	1.82	0.45
1:A:91:CYS:O	1:A:94:ASP:N	2.49	0.45
1:K:67:ASN:ND2	1:K:70:LEU:CD2	2.79	0.45
1:A:44:CYS:SG	1:A:105:CYS:C	3.00	0.45
1:B:25:TYR:HD1	1:B:42:ASP:HB2	1.80	0.45
1:K:94:ASP:HA	1:K:97:GLU:HG2	1.99	0.45
1:K:109:ASN:O	1:K:110:LYS:C	2.59	0.45
1:A:101:VAL:C	1:A:103:ALA:N	2.74	0.45
1:B:8:MET:SD	1:B:99:ASP:C	2.99	0.45
1:B:47:VAL:O	1:B:49:ASP:N	2.50	0.45
1:K:72:ILE:C	1:K:92:LEU:HD11	2.42	0.45
1:K:31:TRP:CH2	2:L:134:ELD:H121	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:32:GLY:H	1:L:2:LEU:HB2	1.81	0.45
1:K:44:CYS:SG	1:K:105:CYS:C	3.00	0.45
1:K:50:CYS:O	1:K:54:ARG:HG3	2.17	0.45
1:L:20:TRP:C	1:L:23:ILE:HG12	2.41	0.45
1:L:30:GLY:H	2:L:134:ELD:C12	2.30	0.45
1:K:5:PHE:O	1:K:5:PHE:CD1	2.70	0.45
1:A:10:ASN:O	1:A:14:GLY:C	2.60	0.45
1:B:95:ILE:C	1:B:97:GLU:N	2.72	0.45
1:A:46:PHE:O	1:A:49:ASP:N	2.51	0.45
1:K:2:LEU:HB2	1:L:31:TRP:O	2.17	0.45
1:A:6:ALA:O	1:A:7:LYS:C	2.60	0.44
1:A:28:TYR:OH	1:A:130:SER:HA	2.16	0.44
1:K:36:THR:HA	1:K:37:PRO:HD3	1.70	0.44
1:L:8:MET:SD	1:L:99:ASP:HB3	2.57	0.44
1:L:47:VAL:O	1:L:50:CYS:N	2.50	0.44
1:B:1:ASN:OD1	1:B:3:PHE:N	2.47	0.44
1:K:6:ALA:C	1:K:8:MET:N	2.73	0.44
1:K:37:PRO:C	1:K:39:ASP:H	2.25	0.44
1:K:104:ASN:O	1:K:105:CYS:C	2.60	0.44
1:A:27:CYS:O	1:A:28:TYR:HD2	2.00	0.44
1:B:4:GLN:O	1:B:6:ALA:N	2.51	0.44
1:K:20:TRP:C	1:K:22:TYR:H	2.24	0.44
1:K:25:TYR:HE2	1:K:114:ASN:O	2.01	0.44
1:L:19:VAL:CG1	1:L:20:TRP:H	2.26	0.44
1:A:94:ASP:O	1:A:96:CYS:N	2.51	0.44
1:K:1:ASN:ND2	1:K:1:ASN:O	2.48	0.44
1:L:10:ASN:O	1:L:11:GLY:C	2.60	0.44
1:A:52:TYR:CE2	1:A:68:PRO:HB2	2.52	0.44
1:B:1:ASN:CG	1:B:2:LEU:H	2.25	0.44
1:B:46:PHE:O	1:B:49:ASP:HB2	2.18	0.44
1:A:67:ASN:OD1	1:A:67:ASN:C	2.61	0.44
1:B:13:LEU:HB3	1:B:18:SER:CB	2.48	0.44
1:L:124:SER:O	1:L:126:CYS:N	2.51	0.44
1:B:43:ARG:O	1:B:44:CYS:C	2.60	0.44
1:B:102:ALA:O	1:B:104:ASN:N	2.51	0.44
1:K:2:LEU:N	1:K:2:LEU:HD23	2.32	0.44
1:B:23:ILE:CG2	1:B:24:SER:N	2.81	0.43
1:K:48:HIS:HE1	1:K:99:ASP:OD2	2.01	0.43
1:L:14:GLY:O	1:L:16:ALA:O	2.36	0.43
1:L:19:VAL:CG1	1:L:20:TRP:N	2.79	0.43
1:K:98:CYS:O	1:K:99:ASP:C	2.60	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:26:GLY:HA3	1:L:42:ASP:HB2	1.99	0.43
1:A:131:GLU:HG3	1:A:132:GLN:N	2.25	0.43
1:K:95:ILE:O	1:K:98:CYS:N	2.52	0.43
1:B:34:GLN:HG3	1:B:126:CYS:O	2.18	0.43
1:B:94:ASP:OD2	1:B:94:ASP:N	2.51	0.43
1:L:26:GLY:HA3	1:L:42:ASP:CB	2.48	0.43
1:L:45:CYS:HA	2:L:134:ELD:H172	2.00	0.43
1:B:5:PHE:CD1	1:B:99:ASP:HA	2.52	0.43
1:L:75:TYR:CD2	1:L:76:SER:N	2.87	0.43
1:B:77:PHE:HA	1:B:81:ASN:O	2.19	0.43
1:B:85:GLY:O	1:B:86:LYS:C	2.61	0.43
1:B:102:ALA:C	1:B:104:ASN:N	2.77	0.43
1:K:67:ASN:O	1:K:68:PRO:C	2.61	0.43
1:L:117:TYR:O	1:L:120:LEU:HG	2.19	0.43
1:A:28:TYR:CD1	1:A:46:PHE:HB2	2.48	0.43
1:B:5:PHE:CE1	1:B:102:ALA:HB3	2.54	0.43
1:K:93:ARG:O	1:K:96:CYS:N	2.52	0.43
1:L:61:CYS:SG	1:L:92:LEU:HD23	2.59	0.43
1:L:85:GLY:O	1:L:88:ASN:N	2.52	0.43
1:B:36:THR:HG23	1:B:131:GLU:HB2	2.01	0.43
1:L:45:CYS:O	1:L:46:PHE:C	2.60	0.43
1:K:34:GLN:O	1:K:126:CYS:HB3	2.19	0.43
1:B:50:CYS:O	1:B:53:GLY:N	2.52	0.42
1:B:91:CYS:O	1:B:94:ASP:HB2	2.19	0.42
1:B:13:LEU:HB3	1:B:18:SER:OG	2.19	0.42
1:B:5:PHE:HB2	1:B:99:ASP:CG	2.43	0.42
1:B:101:VAL:O	1:B:104:ASN:N	2.52	0.42
1:K:43:ARG:O	1:K:46:PHE:HB3	2.19	0.42
1:L:104:ASN:HD22	1:L:104:ASN:HA	1.62	0.42
1:L:126:CYS:HA	1:L:128:GLN:HE22	1.84	0.42
1:A:7:LYS:HA	1:A:10:ASN:HD22	1.83	0.42
1:B:34:GLN:CG	1:B:126:CYS:HB2	2.49	0.42
1:B:67:ASN:HA	1:B:68:PRO:HD3	1.81	0.42
1:L:31:TRP:O	1:L:32:GLY:O	2.36	0.42
1:L:35:GLY:CA	1:L:128:GLN:OE1	2.67	0.42
1:A:119:PHE:O	1:A:120:LEU:HD23	2.20	0.42
1:L:42:ASP:O	1:L:43:ARG:C	2.61	0.42
1:B:79:LYS:O	1:B:81:ASN:N	2.53	0.42
1:K:116:ASN:C	1:K:118:LYS:N	2.78	0.42
1:L:2:LEU:C	1:L:4:GLN:N	2.74	0.42
1:B:106:PHE:CE1	2:B:134:ELD:H171	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:40:ALA:O	1:A:43:ARG:N	2.53	0.42
1:A:44:CYS:SG	1:A:105:CYS:O	2.77	0.42
1:K:90:GLY:O	1:K:91:CYS:C	2.62	0.42
1:K:97:GLU:O	1:K:98:CYS:C	2.62	0.42
1:L:45:CYS:HA	2:L:134:ELD:C15	2.50	0.42
1:L:46:PHE:O	1:L:49:ASP:HB2	2.20	0.42
1:L:95:ILE:HG22	1:L:96:CYS:N	2.34	0.42
1:A:96:CYS:O	1:A:97:GLU:C	2.62	0.42
1:K:5:PHE:CD1	1:K:5:PHE:C	2.97	0.42
1:L:124:SER:C	1:L:126:CYS:H	2.28	0.42
1:A:48:HIS:O	1:A:49:ASP:C	2.63	0.41
1:A:80:GLY:O	1:A:81:ASN:ND2	2.53	0.41
1:A:88:ASN:HB3	1:A:93:ARG:HB2	2.02	0.41
1:K:70:LEU:HD13	1:K:70:LEU:N	2.35	0.41
1:L:101:VAL:HA	1:L:104:ASN:CB	2.38	0.41
1:A:7:LYS:O	1:A:8:MET:C	2.63	0.41
1:A:112:THR:O	1:A:112:THR:HG22	2.20	0.41
1:B:67:ASN:O	1:B:71:ALA:N	2.53	0.41
1:A:124:SER:HB2	1:A:127:ARG:CZ	2.50	0.41
1:B:61:CYS:O	1:B:68:PRO:HD3	2.20	0.41
1:B:125:ARG:HD2	1:B:125:ARG:HA	1.88	0.41
1:K:1:ASN:O	1:K:2:LEU:C	2.63	0.41
1:K:52:TYR:O	1:K:55:VAL:CG1	2.67	0.41
1:K:68:PRO:HG3	1:K:95:ILE:CD1	2.50	0.41
1:K:93:ARG:C	1:K:95:ILE:N	2.78	0.41
1:K:100:ARG:NH2	1:K:101:VAL:CG2	2.75	0.41
1:A:114:ASN:C	1:A:114:ASN:HD22	2.29	0.41
1:B:99:ASP:O	1:B:100:ARG:O	2.38	0.41
1:B:115:LYS:O	1:B:117:TYR:N	2.53	0.41
1:K:5:PHE:O	1:K:5:PHE:HD1	2.04	0.41
1:K:26:GLY:HA3	1:K:42:ASP:HB2	2.03	0.41
1:L:5:PHE:CE1	1:L:48:HIS:HD2	2.38	0.41
1:L:45:CYS:SG	2:L:134:ELD:H131	2.60	0.41
1:A:8:MET:HB3	1:A:82:ILE:HD11	2.01	0.41
1:A:52:TYR:CE1	1:A:68:PRO:HB3	2.55	0.41
1:L:14:GLY:O	1:L:16:ALA:C	2.62	0.41
1:K:45:CYS:O	1:K:46:PHE:C	2.62	0.41
1:L:98:CYS:O	1:L:102:ALA:N	2.48	0.41
1:A:12:LYS:HB3	1:A:13:LEU:HD12	2.03	0.41
1:K:10:ASN:HD21	1:K:19:VAL:HB	1.86	0.41
1:L:102:ALA:O	2:L:134:ELD:H182	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:46:PHE:CZ	1:B:131:GLU:O	2.74	0.41
1:B:50:CYS:O	1:B:51:CYS:C	2.64	0.41
1:B:92:LEU:HD12	1:B:92:LEU:N	2.36	0.41
1:L:8:MET:O	1:L:9:ILE:C	2.63	0.41
1:A:88:ASN:CB	1:A:93:ARG:HB2	2.51	0.41
1:B:105:CYS:O	1:B:109:ASN:OD1	2.37	0.41
1:K:2:LEU:HB2	1:L:31:TRP:HB3	2.03	0.41
1:K:13:LEU:HD12	1:K:13:LEU:N	2.36	0.41
1:K:19:VAL:HG13	1:K:20:TRP:N	2.35	0.41
1:K:118:LYS:O	1:K:119:PHE:C	2.63	0.41
1:L:95:ILE:O	1:L:96:CYS:C	2.63	0.41
1:L:106:PHE:CZ	2:L:134:ELD:H162	2.56	0.41
1:A:52:TYR:CZ	1:A:68:PRO:HB3	2.56	0.41
1:L:2:LEU:HA	1:L:5:PHE:HB3	2.03	0.41
1:L:10:ASN:O	1:L:14:GLY:N	2.53	0.41
1:L:39:ASP:OD2	1:L:41:THR:N	2.54	0.41
1:A:40:ALA:O	1:A:41:THR:C	2.63	0.40
1:K:4:GLN:NE2	1:K:73:TYR:CD1	2.89	0.40
1:K:67:ASN:HD22	1:L:53:GLY:HA3	1.86	0.40
1:B:114:ASN:O	1:B:115:LYS:C	2.64	0.40
1:K:6:ALA:O	1:K:7:LYS:C	2.63	0.40
1:L:5:PHE:CZ	1:L:48:HIS:CD2	3.08	0.40
1:L:72:ILE:HG23	1:L:72:ILE:O	2.22	0.40
1:B:104:ASN:HD22	1:B:104:ASN:HA	1.68	0.40
1:L:98:CYS:O	1:L:99:ASP:C	2.65	0.40
1:A:28:TYR:HH	1:A:130:SER:HA	1.87	0.40
1:A:31:TRP:CD1	1:A:31:TRP:N	2.90	0.40
1:A:98:CYS:O	1:A:99:ASP:C	2.64	0.40
1:L:44:CYS:HB3	2:L:134:ELD:C18	2.49	0.40
1:L:82:ILE:HG22	1:L:83:VAL:N	2.37	0.40
1:B:68:PRO:HA	1:B:71:ALA:HB3	2.03	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	120/122 (98%)	72 (60%)	33 (28%)	15 (12%)	0	1
1	B	120/122 (98%)	69 (58%)	33 (28%)	18 (15%)	0	1
1	K	120/122 (98%)	70 (58%)	34 (28%)	16 (13%)	0	1
1	L	120/122 (98%)	66 (55%)	35 (29%)	19 (16%)	0	1
All	All	480/488 (98%)	277 (58%)	135 (28%)	68 (14%)	0	1

All (68) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	46	PHE
1	A	55	VAL
1	A	127	ARG
1	B	16	ALA
1	B	17	PHE
1	B	26	GLY
1	B	47	VAL
1	K	25	TYR
1	K	92	LEU
1	K	115	LYS
1	L	16	ALA
1	L	19	VAL
1	L	76	SER
1	L	86	LYS
1	L	100	ARG
1	A	20	TRP
1	A	45	CYS
1	A	79	LYS
1	A	86	LYS
1	A	95	ILE
1	B	100	ARG
1	B	102	ALA
1	B	124	SER
1	K	38	LYS
1	K	119	PHE
1	K	122	SER
1	L	32	GLY
1	L	47	VAL
1	L	124	SER
1	L	125	ARG

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Mol	Chain	Res	Type
1	L	130	SER
1	A	11	GLY
1	A	25	TYR
1	A	131	GLU
1	B	86	LYS
1	K	11	GLY
1	K	75	TYR
1	K	79	LYS
1	K	98	CYS
1	K	112	THR
1	L	37	PRO
1	L	89	ASN
1	L	108	GLN
1	B	48	HIS
1	B	116	ASN
1	K	2	LEU
1	K	55	VAL
1	L	95	ILE
1	A	76	SER
1	B	37	PRO
1	B	80	GLY
1	B	85	GLY
1	B	101	VAL
1	K	132	GLN
1	L	120	LEU
1	A	100	ARG
1	B	127	ARG
1	L	14	GLY
1	A	80	GLY
1	B	55	VAL
1	L	26	GLY
1	A	72	ILE
1	K	33	GLY
1	L	9	ILE
1	L	68	PRO
1	K	80	GLY
1	B	14	GLY
1	B	68	PRO

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	104/104 (100%)	100 (96%)	4 (4%)	29	57
1	B	104/104 (100%)	91 (88%)	13 (12%)	4	19
1	K	104/104 (100%)	97 (93%)	7 (7%)	15	42
1	L	104/104 (100%)	95 (91%)	9 (9%)	9	32
All	All	416/416 (100%)	383 (92%)	33 (8%)	11	36

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	VAL
1	A	20	TRP
1	A	47	VAL
1	A	83	VAL
1	B	3	PHE
1	B	12	LYS
1	B	20	TRP
1	B	34	GLN
1	B	36	THR
1	B	56	ARG
1	B	77	PHE
1	B	83	VAL
1	B	94	ASP
1	B	97	GLU
1	B	124	SER
1	B	128	GLN
1	B	129	THR
1	K	1	ASN
1	K	2	LEU
1	K	41	THR
1	K	55	VAL
1	K	69	LYS
1	K	70	LEU
1	K	82	ILE

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Mol	Chain	Res	Type
1	L	3	PHE
1	L	7	LYS
1	L	39	ASP
1	L	41	THR
1	L	54	ARG
1	L	101	VAL
1	L	104	ASN
1	L	108	GLN
1	L	121	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	81	ASN
1	A	104	ASN
1	A	107	HIS
1	B	104	ASN
1	B	128	GLN
1	K	1	ASN
1	K	10	ASN
1	K	21	ASN
1	K	67	ASN
1	K	108	GLN
1	K	109	ASN
1	K	114	ASN
1	L	4	GLN
1	L	34	GLN
1	L	88	ASN
1	L	108	GLN
1	L	111	ASN
1	L	114	ASN
1	L	132	GLN

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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 ligands 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	ELD	L	134	-	19,19,19	0.43	0	19,19,19	1.65	3 (15%)
2	ELD	B	134	-	19,19,19	0.62	0	19,19,19	1.69	3 (15%)

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	ELD	L	134	-	-	11/17/17/17	-
2	ELD	B	134	-	-	11/17/17/17	-

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	134	ELD	O2-C1-C2	-5.42	104.68	121.04
2	L	134	ELD	O2-C1-C2	-5.25	105.20	121.04
2	B	134	ELD	C4-C3-C2	-3.59	99.94	113.13
2	L	134	ELD	C4-C3-C2	-3.30	101.01	113.13
2	L	134	ELD	C6-C5-C4	-2.92	99.63	114.37
2	B	134	ELD	C6-C5-C4	-2.47	101.88	114.37

There are no chirality outliers.

All (22) torsion outliers are listed below:

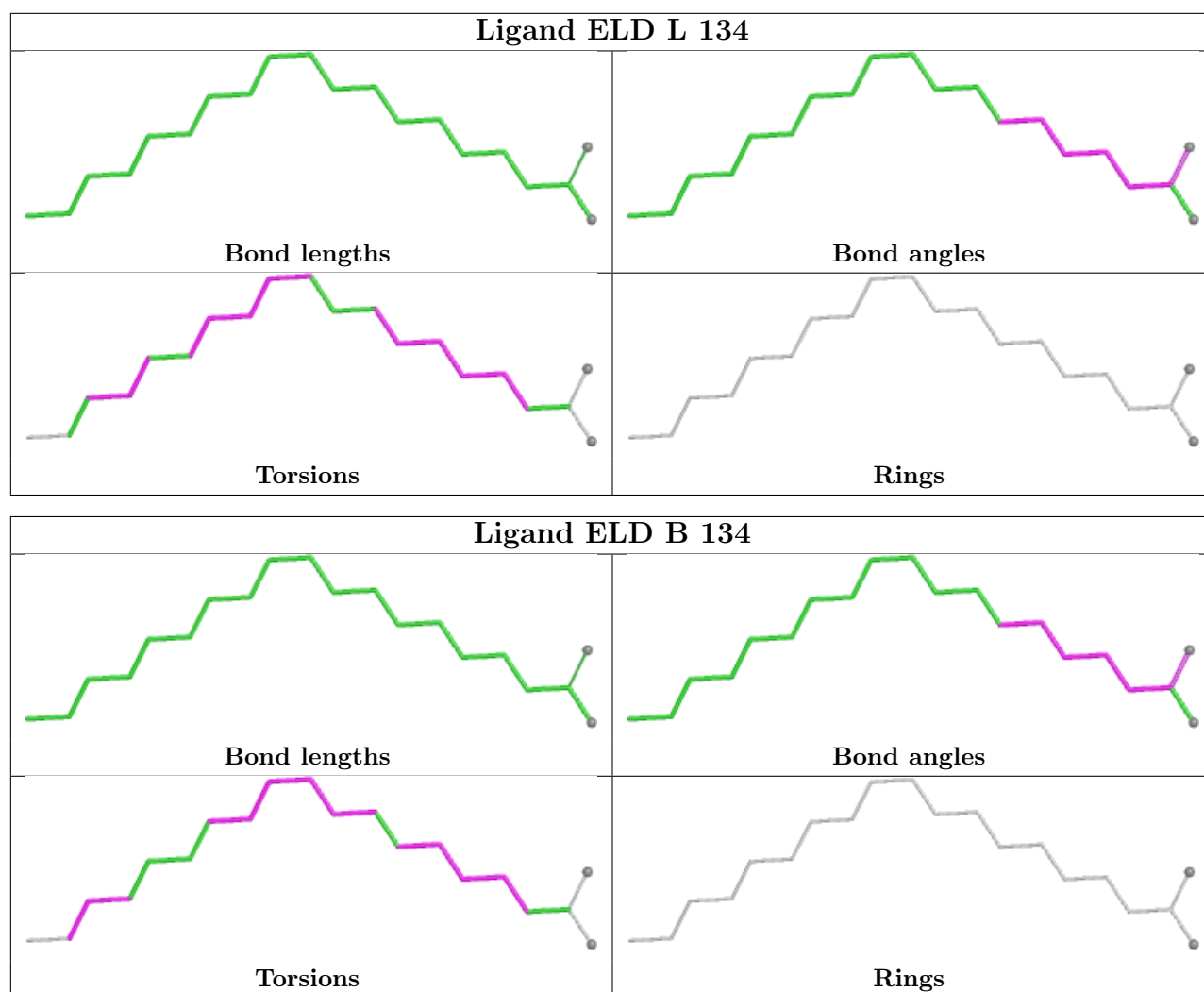
Mol	Chain	Res	Type	Atoms
2	L	134	ELD	C1-C2-C3-C4
2	B	134	ELD	C11-C10-C9-C8
2	L	134	ELD	C11-C10-C9-C8
2	L	134	ELD	C11-C12-C13-C14
2	B	134	ELD	C4-C5-C6-C7
2	L	134	ELD	C4-C5-C6-C7
2	B	134	ELD	C3-C4-C5-C6
2	B	134	ELD	C14-C15-C16-C17
2	L	134	ELD	C14-C15-C16-C17
2	L	134	ELD	C5-C6-C7-C8
2	L	134	ELD	C2-C3-C4-C5
2	B	134	ELD	C6-C7-C8-C9
2	B	134	ELD	C10-C11-C12-C13
2	L	134	ELD	C10-C11-C12-C13
2	B	134	ELD	C15-C16-C17-C18
2	L	134	ELD	C3-C4-C5-C6
2	B	134	ELD	C1-C2-C3-C4
2	B	134	ELD	C2-C3-C4-C5
2	L	134	ELD	C9-C10-C11-C12
2	L	134	ELD	C13-C14-C15-C16
2	B	134	ELD	C9-C10-C11-C12
2	B	134	ELD	C7-C8-C9-C10

There are no ring outliers.

2 monomers are involved in 28 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	L	134	ELD	16	0
2	B	134	ELD	12	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



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

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.