



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 12:40 PM UTC

PDB ID : 6NDB / pdb_00006ndb
Title : RHODOCETIN IN COMPLEX WITH THE INTEGRIN ALPHA2-A DOMAIN AND COBALT
Authors : Stetefeld, J.; McDougall, M.D.; Loewen, P.C.
Deposited on : 2018-12-13
Resolution : 3.20 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

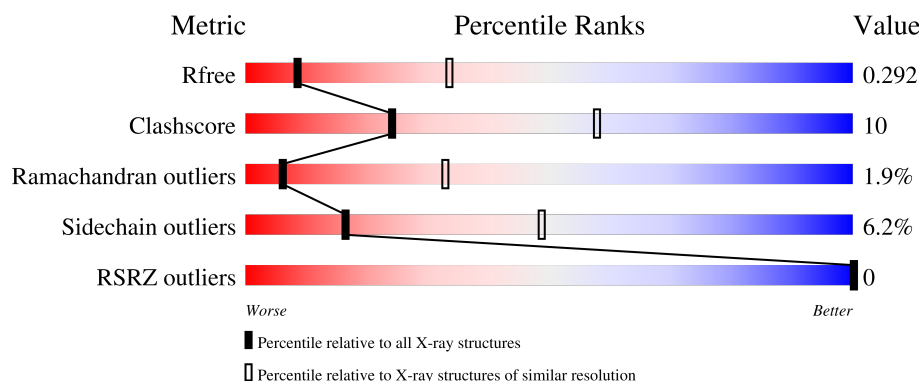
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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	1466 (3.20-3.20)
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (3.20-3.20)

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	135	70% 24% • •
1	D	135	72% 22% • • •
1	G	135	75% 19% • •
1	J	135	76% 19% • •
1	M	135	59% 35% • • •

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Mol	Chain	Length	Quality of chain
1	P	135	 65% 29% . . .
2	B	124	 69% 27% . . .
2	E	124	 67% 27% . .
2	H	124	 69% 27% . .
2	K	124	 69% 27% . . .
2	N	124	 61% 35% . .
2	Q	124	 67% 27% . . .
3	C	217	 72% 14% . 12%
3	F	217	 68% 18% . 12%
3	I	217	 67% 19% . 12%
3	L	217	 70% 17% 12%
3	O	217	 65% 21% . 12%
3	R	217	 69% 17% . 12%

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 21590 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 Snaclec rhodocetin subunit gamma.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	131	Total	C	N	O	S	0	0	0
			1070	672	189	200	9			
1	D	131	Total	C	N	O	S	0	0	0
			1070	672	189	200	9			
1	G	131	Total	C	N	O	S	0	0	0
			1070	672	189	200	9			
1	J	131	Total	C	N	O	S	0	0	0
			1070	672	189	200	9			
1	M	131	Total	C	N	O	S	0	0	0
			1070	672	189	200	9			
1	P	131	Total	C	N	O	S	0	0	0
			1070	672	189	200	9			

- Molecule 2 is a protein called Snaclec rhodocetin subunit delta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	122	Total	C	N	O	S	0	0	0
			1029	667	174	179	9			
2	E	122	Total	C	N	O	S	0	0	0
			1029	667	174	179	9			
2	H	122	Total	C	N	O	S	0	0	0
			1029	667	174	179	9			
2	K	122	Total	C	N	O	S	0	0	0
			1029	667	174	179	9			
2	N	122	Total	C	N	O	S	0	0	0
			1029	667	174	179	9			
2	Q	122	Total	C	N	O	S	0	0	0
			1029	667	174	179	9			

- Molecule 3 is a protein called Integrin alpha-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	191	Total	C	N	O	S	0	0	0
			1482	940	250	287	5			
3	F	191	Total	C	N	O	S	0	0	0
			1482	940	250	287	5			
3	I	191	Total	C	N	O	S	0	0	0
			1482	940	250	287	5			
3	L	191	Total	C	N	O	S	0	0	0
			1482	940	250	287	5			
3	O	191	Total	C	N	O	S	0	0	0
			1482	940	250	287	5			
3	R	191	Total	C	N	O	S	0	0	0
			1482	940	250	287	5			

There are 120 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	150	MET	-	expression tag	UNP P17301
C	151	GLY	-	expression tag	UNP P17301
C	152	SER	-	expression tag	UNP P17301
C	153	SER	-	expression tag	UNP P17301
C	154	HIS	-	expression tag	UNP P17301
C	155	HIS	-	expression tag	UNP P17301
C	156	HIS	-	expression tag	UNP P17301
C	157	HIS	-	expression tag	UNP P17301
C	158	HIS	-	expression tag	UNP P17301
C	159	HIS	-	expression tag	UNP P17301
C	160	SER	-	expression tag	UNP P17301
C	161	SER	-	expression tag	UNP P17301
C	162	GLY	-	expression tag	UNP P17301
C	163	LEU	-	expression tag	UNP P17301
C	164	VAL	-	expression tag	UNP P17301
C	165	PRO	-	expression tag	UNP P17301
C	166	ARG	-	expression tag	UNP P17301
C	167	GLY	-	expression tag	UNP P17301
C	168	GLY	-	expression tag	UNP P17301
C	169	SER	-	expression tag	UNP P17301
F	150	MET	-	expression tag	UNP P17301
F	151	GLY	-	expression tag	UNP P17301
F	152	SER	-	expression tag	UNP P17301
F	153	SER	-	expression tag	UNP P17301
F	154	HIS	-	expression tag	UNP P17301
F	155	HIS	-	expression tag	UNP P17301
F	156	HIS	-	expression tag	UNP P17301
F	157	HIS	-	expression tag	UNP P17301

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Chain	Residue	Modelled	Actual	Comment	Reference
F	158	HIS	-	expression tag	UNP P17301
F	159	HIS	-	expression tag	UNP P17301
F	160	SER	-	expression tag	UNP P17301
F	161	SER	-	expression tag	UNP P17301
F	162	GLY	-	expression tag	UNP P17301
F	163	LEU	-	expression tag	UNP P17301
F	164	VAL	-	expression tag	UNP P17301
F	165	PRO	-	expression tag	UNP P17301
F	166	ARG	-	expression tag	UNP P17301
F	167	GLY	-	expression tag	UNP P17301
F	168	GLY	-	expression tag	UNP P17301
F	169	SER	-	expression tag	UNP P17301
I	150	MET	-	expression tag	UNP P17301
I	151	GLY	-	expression tag	UNP P17301
I	152	SER	-	expression tag	UNP P17301
I	153	SER	-	expression tag	UNP P17301
I	154	HIS	-	expression tag	UNP P17301
I	155	HIS	-	expression tag	UNP P17301
I	156	HIS	-	expression tag	UNP P17301
I	157	HIS	-	expression tag	UNP P17301
I	158	HIS	-	expression tag	UNP P17301
I	159	HIS	-	expression tag	UNP P17301
I	160	SER	-	expression tag	UNP P17301
I	161	SER	-	expression tag	UNP P17301
I	162	GLY	-	expression tag	UNP P17301
I	163	LEU	-	expression tag	UNP P17301
I	164	VAL	-	expression tag	UNP P17301
I	165	PRO	-	expression tag	UNP P17301
I	166	ARG	-	expression tag	UNP P17301
I	167	GLY	-	expression tag	UNP P17301
I	168	GLY	-	expression tag	UNP P17301
I	169	SER	-	expression tag	UNP P17301
L	150	MET	-	expression tag	UNP P17301
L	151	GLY	-	expression tag	UNP P17301
L	152	SER	-	expression tag	UNP P17301
L	153	SER	-	expression tag	UNP P17301
L	154	HIS	-	expression tag	UNP P17301
L	155	HIS	-	expression tag	UNP P17301
L	156	HIS	-	expression tag	UNP P17301
L	157	HIS	-	expression tag	UNP P17301
L	158	HIS	-	expression tag	UNP P17301
L	159	HIS	-	expression tag	UNP P17301

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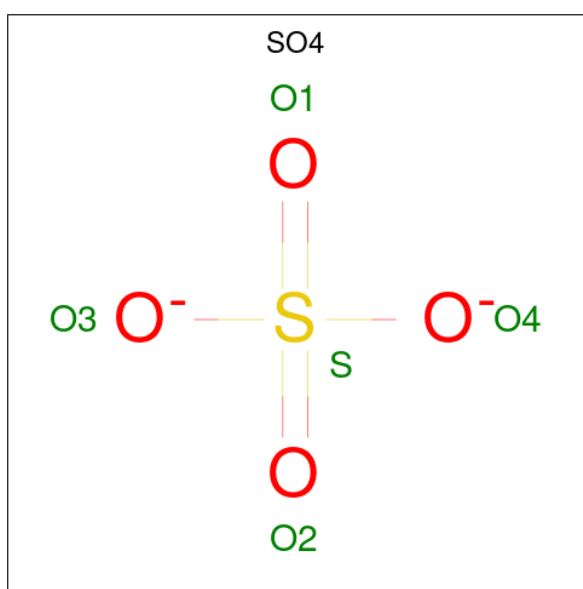
Chain	Residue	Modelled	Actual	Comment	Reference
L	160	SER	-	expression tag	UNP P17301
L	161	SER	-	expression tag	UNP P17301
L	162	GLY	-	expression tag	UNP P17301
L	163	LEU	-	expression tag	UNP P17301
L	164	VAL	-	expression tag	UNP P17301
L	165	PRO	-	expression tag	UNP P17301
L	166	ARG	-	expression tag	UNP P17301
L	167	GLY	-	expression tag	UNP P17301
L	168	GLY	-	expression tag	UNP P17301
L	169	SER	-	expression tag	UNP P17301
O	150	MET	-	expression tag	UNP P17301
O	151	GLY	-	expression tag	UNP P17301
O	152	SER	-	expression tag	UNP P17301
O	153	SER	-	expression tag	UNP P17301
O	154	HIS	-	expression tag	UNP P17301
O	155	HIS	-	expression tag	UNP P17301
O	156	HIS	-	expression tag	UNP P17301
O	157	HIS	-	expression tag	UNP P17301
O	158	HIS	-	expression tag	UNP P17301
O	159	HIS	-	expression tag	UNP P17301
O	160	SER	-	expression tag	UNP P17301
O	161	SER	-	expression tag	UNP P17301
O	162	GLY	-	expression tag	UNP P17301
O	163	LEU	-	expression tag	UNP P17301
O	164	VAL	-	expression tag	UNP P17301
O	165	PRO	-	expression tag	UNP P17301
O	166	ARG	-	expression tag	UNP P17301
O	167	GLY	-	expression tag	UNP P17301
O	168	GLY	-	expression tag	UNP P17301
O	169	SER	-	expression tag	UNP P17301
R	150	MET	-	expression tag	UNP P17301
R	151	GLY	-	expression tag	UNP P17301
R	152	SER	-	expression tag	UNP P17301
R	153	SER	-	expression tag	UNP P17301
R	154	HIS	-	expression tag	UNP P17301
R	155	HIS	-	expression tag	UNP P17301
R	156	HIS	-	expression tag	UNP P17301
R	157	HIS	-	expression tag	UNP P17301
R	158	HIS	-	expression tag	UNP P17301
R	159	HIS	-	expression tag	UNP P17301
R	160	SER	-	expression tag	UNP P17301
R	161	SER	-	expression tag	UNP P17301

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Chain	Residue	Modelled	Actual	Comment	Reference
R	162	GLY	-	expression tag	UNP P17301
R	163	LEU	-	expression tag	UNP P17301
R	164	VAL	-	expression tag	UNP P17301
R	165	PRO	-	expression tag	UNP P17301
R	166	ARG	-	expression tag	UNP P17301
R	167	GLY	-	expression tag	UNP P17301
R	168	GLY	-	expression tag	UNP P17301
R	169	SER	-	expression tag	UNP P17301

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total O S 5 4 1	0	0
4	A	1	Total O S 5 4 1	0	0
4	B	1	Total O S 5 4 1	0	0
4	D	1	Total O S 5 4 1	0	0
4	E	1	Total O S 5 4 1	0	0
4	E	1	Total O S 5 4 1	0	0
4	G	1	Total O S 5 4 1	0	0

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	H	1	Total O S 5 4 1	0	0
4	J	1	Total O S 5 4 1	0	0
4	K	1	Total O S 5 4 1	0	0
4	N	1	Total O S 5 4 1	0	0
4	O	1	Total O S 5 4 1	0	0
4	Q	1	Total O S 5 4 1	0	0
4	Q	1	Total O S 5 4 1	0	0
4	R	1	Total O S 5 4 1	0	0

- Molecule 5 is COBALT (II) ION (CCD ID: CO) (formula: Co).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	C	1	Total Co 1 1	0	0
5	F	1	Total Co 1 1	0	0
5	I	1	Total Co 1 1	0	0
5	L	1	Total Co 1 1	0	0
5	O	1	Total Co 1 1	0	0
5	R	1	Total Co 1 1	0	0

- Molecule 6 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	C	1	Total Na 1 1	0	0
6	F	1	Total Na 1 1	0	0
6	I	1	Total Na 1 1	0	0

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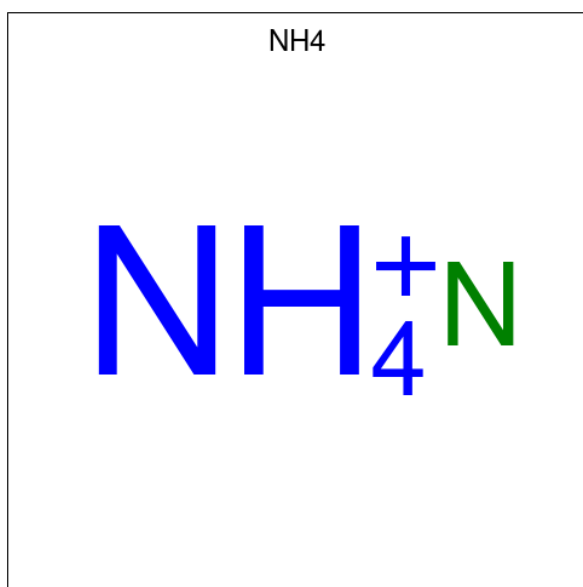
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	L	1	Total 1	Na 1	0	0
6	O	1	Total 1	Na 1	0	0
6	R	1	Total 1	Na 1	0	0

- Molecule 7 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	C	1	Total 1	Cl 1	0	0
7	E	1	Total 1	Cl 1	0	0
7	G	1	Total 1	Cl 1	0	0
7	H	1	Total 1	Cl 1	0	0
7	J	1	Total 1	Cl 1	0	0
7	K	1	Total 1	Cl 1	0	0
7	L	1	Total 1	Cl 1	0	0
7	M	1	Total 1	Cl 1	0	0
7	R	1	Total 1	Cl 1	0	0

- Molecule 8 is AMMONIUM ION (CCD ID: NH4) (formula: H₄N).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	D	1	Total N 1 1	0	0

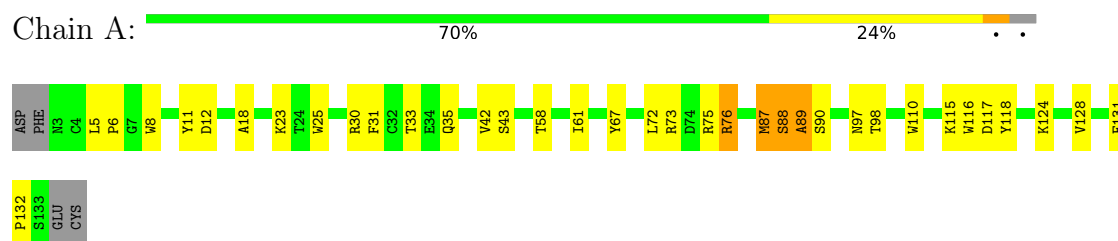
- Molecule 9 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	A	5	Total O 5 5	0	0
9	D	1	Total O 1 1	0	0
9	F	1	Total O 1 1	0	0

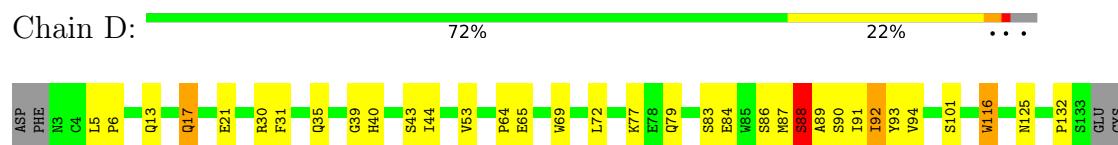
3 Residue-property plots

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 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.

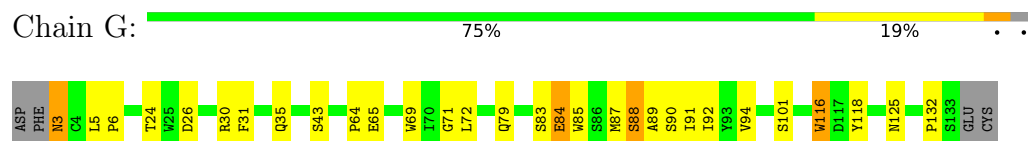
- Molecule 1: Snaclec rhodocetin subunit gamma



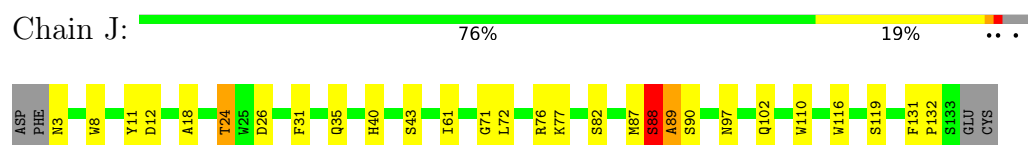
- Molecule 1: Snaclec rhodocetin subunit gamma



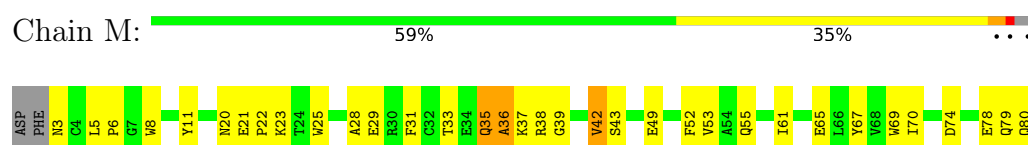
- Molecule 1: Snaclec rhodocetin subunit gamma



- Molecule 1: Snaclec rhodocetin subunit gamma



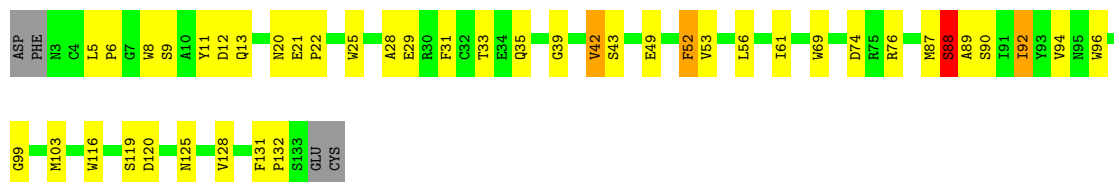
- Molecule 1: Snaclec rhodocetin subunit gamma





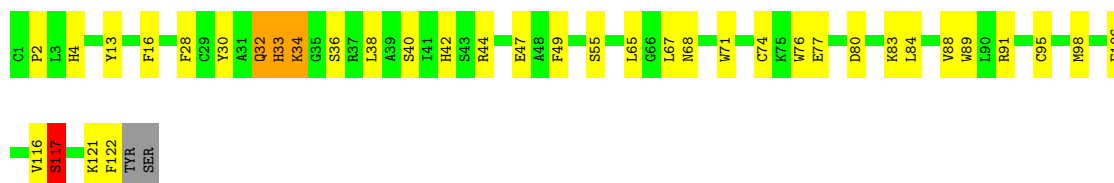
- Molecule 1: Snaclec rhodocetin subunit gamma

Chain P: 65% 29% ...



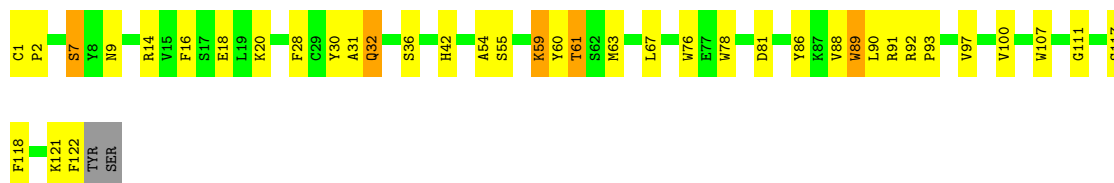
- Molecule 2: Snaclec rhodocetin subunit delta

Chain B: 69% 27% ...



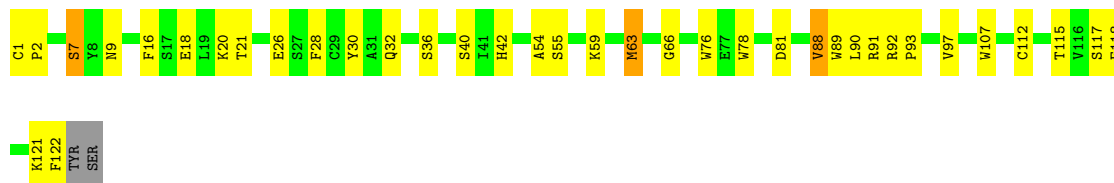
- Molecule 2: Snaclec rhodocetin subunit delta

Chain E: 67% 27% ..



- Molecule 2: Snaclec rhodocetin subunit delta

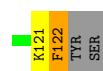
Chain H: 69% 27% ..



- Molecule 2: Snaclec rhodocetin subunit delta

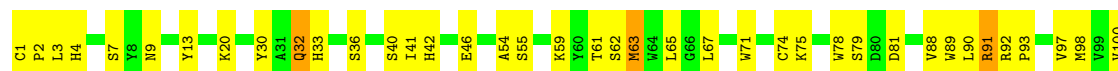
Chain K: 69% 27% ...





- Molecule 2: Snaclec rhodocetin subunit delta

Chain N: 61% 35% ..



- Molecule 2: Snaclec rhodocetin subunit delta

Chain Q: 67% 27% ...



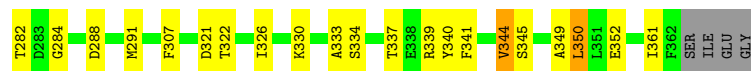
- Molecule 3: Integrin alpha-2

Chain C: 72% 14% 12%



- Molecule 3: Integrin alpha-2

Chain F: 68% 18% 12%



- Molecule 3: Integrin alpha-2

Chain I: 67% 19% 12%

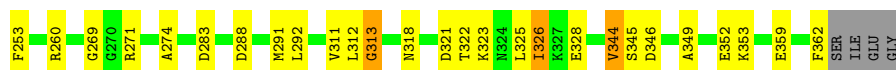
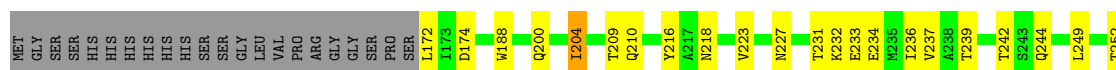




• Molecule 3: Integrin alpha-2



• Molecule 3: Integrin alpha-2



• Molecule 3: Integrin alpha-2



4 Data and refinement statistics

Property	Value	Source
Space group	P 41	Depositor
Cell constants a, b, c, α , β , γ	131.84Å 131.84Å 251.70Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.07 – 3.20 47.07 – 3.20	Depositor EDS
% Data completeness (in resolution range)	99.0 (47.07-3.20) 99.1 (47.07-3.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.14	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.98 (at 3.19Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.228 , 0.292 0.230 , 0.292	Depositor DCC
R_{free} test set	3554 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	62.1	Xtriage
Anisotropy	0.062	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 55.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.458 for h,-k,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	21590	wwPDB-VP
Average B, all atoms (Å ²)	67.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.31% 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: SO4, NA, NH4, CL, CO

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	1.02	0/1101	1.37	0/1491
1	D	1.05	0/1101	1.42	1/1491 (0.1%)
1	G	1.04	0/1101	1.41	0/1491
1	J	1.02	0/1101	1.38	0/1491
1	M	1.07	0/1101	1.44	0/1491
1	P	1.07	0/1101	1.44	1/1491 (0.1%)
2	B	1.02	0/1064	1.35	0/1439
2	E	0.99	0/1064	1.33	0/1439
2	H	0.97	0/1064	1.37	0/1439
2	K	1.02	0/1064	1.32	0/1439
2	N	1.04	0/1064	1.33	0/1439
2	Q	1.04	0/1064	1.34	2/1439 (0.1%)
3	C	1.05	0/1506	1.51	3/2040 (0.1%)
3	F	1.04	0/1506	1.48	0/2040
3	I	1.04	0/1506	1.49	1/2040 (0.0%)
3	L	1.04	0/1506	1.49	2/2040 (0.1%)
3	O	1.03	0/1506	1.47	2/2040 (0.1%)
3	R	1.02	0/1506	1.48	1/2040 (0.0%)
All	All	1.03	0/22026	1.42	13/29820 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	G	0	1
1	J	0	1
2	B	0	1
2	E	0	1

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Mol	Chain	#Chirality outliers	#Planarity outliers
All	All	0	5

There are no bond length outliers.

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	P	92	ILE	N-CA-C	-6.80	106.61	113.47
1	D	92	ILE	N-CA-C	-5.78	107.86	113.53
3	O	362	PHE	CA-CB-CG	5.75	119.56	113.80
3	L	362	PHE	CA-CB-CG	5.34	119.14	113.80
3	R	223	VAL	N-CA-C	-5.29	107.26	111.81
3	C	331	ALA	CA-C-N	5.25	127.26	120.70
3	C	331	ALA	C-N-CA	5.25	127.26	120.70
3	L	223	VAL	N-CA-C	-5.20	107.34	111.81
3	I	283	ASP	CA-CB-CG	5.14	117.74	112.60
3	O	223	VAL	N-CA-C	-5.06	107.46	111.81
3	C	223	VAL	N-CA-C	-5.05	107.81	111.90
2	Q	121	LYS	CA-C-N	5.00	130.70	121.70
2	Q	121	LYS	C-N-CA	5.00	130.70	121.70

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	132	PRO	Peptide
2	B	4	HIS	Peptide
2	E	89	TRP	Peptide
1	G	132	PRO	Peptide
1	J	88	SER	Peptide

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	1070	0	988	29	0
1	D	1070	0	988	17	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	G	1070	0	988	21	0
1	J	1070	0	988	26	0
1	M	1070	0	988	43	0
1	P	1070	0	988	37	0
2	B	1029	0	977	27	0
2	E	1029	0	977	33	0
2	H	1029	0	977	30	0
2	K	1029	0	977	24	0
2	N	1029	0	977	38	0
2	Q	1029	0	977	29	0
3	C	1482	0	1478	15	0
3	F	1482	0	1479	23	0
3	I	1482	0	1479	23	0
3	L	1482	0	1478	22	0
3	O	1482	0	1478	32	0
3	R	1482	0	1478	25	0
4	A	10	0	0	0	0
4	B	5	0	0	0	0
4	D	5	0	0	0	0
4	E	10	0	0	0	0
4	G	5	0	0	0	0
4	H	5	0	0	0	0
4	J	5	0	0	0	0
4	K	5	0	0	0	0
4	N	5	0	0	0	0
4	O	5	0	0	1	0
4	Q	10	0	0	0	0
4	R	5	0	0	0	0
5	C	1	0	0	0	0
5	F	1	0	0	0	0
5	I	1	0	0	0	0
5	L	1	0	0	0	0
5	O	1	0	0	0	0
5	R	1	0	0	0	0
6	C	1	0	0	0	0
6	F	1	0	0	0	0
6	I	1	0	0	0	0
6	L	1	0	0	0	0
6	O	1	0	0	0	0
6	R	1	0	0	0	0
7	C	1	0	0	1	0
7	E	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	G	1	0	0	0	0
7	H	1	0	0	0	0
7	J	1	0	0	0	0
7	K	1	0	0	0	0
7	L	1	0	0	0	0
7	M	1	0	0	0	0
7	R	1	0	0	0	0
8	D	1	0	0	0	0
9	A	5	0	0	0	0
9	D	1	0	0	0	0
9	F	1	0	0	0	0
All	All	21590	0	20660	422	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 10.

All (422) 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:M:87:MET:SD	2:N:40:SER:OG	2.35	0.85
1:M:89:ALA:HB1	2:N:42:HIS:NE2	1.93	0.84
1:M:29:GLU:CD	2:N:79:SER:HG	1.87	0.82
1:P:87:MET:SD	2:Q:40:SER:OG	2.37	0.82
2:Q:54:ALA:CB	2:Q:63:MET:HE1	2.10	0.81
2:Q:54:ALA:HB1	2:Q:63:MET:HE1	1.61	0.81
1:M:29:GLU:CD	2:N:79:SER:OG	2.23	0.81
2:E:90:LEU:HD12	2:E:90:LEU:C	2.06	0.81
2:N:32:GLN:HE21	2:N:32:GLN:CA	1.92	0.81
1:P:89:ALA:HB1	2:Q:42:HIS:CE1	2.15	0.81
2:N:54:ALA:CB	2:N:63:MET:HE1	2.13	0.78
2:H:20:LYS:HE2	2:H:28:PHE:CD2	2.19	0.77
1:M:87:MET:O	1:M:88:SER:HB2	1.84	0.76
1:P:87:MET:O	1:P:88:SER:HB2	1.84	0.75
2:H:90:LEU:HD12	2:H:90:LEU:C	2.11	0.74
3:R:204:ILE:HG12	3:R:235:MET:HE1	1.69	0.74
1:P:29:GLU:CD	2:Q:79:SER:OG	2.31	0.74
2:E:31:ALA:HB1	3:O:353:LYS:HE2	1.71	0.72
3:I:344:VAL:HG13	3:I:349:ALA:HB3	1.73	0.69
2:E:54:ALA:HB1	2:E:63:MET:HE1	1.74	0.69
3:R:204:ILE:HA	3:R:209:THR:O	1.94	0.68
1:M:29:GLU:OE2	2:N:79:SER:OG	2.12	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:29:GLU:OE1	2:N:79:SER:OG	2.11	0.67
2:E:20:LYS:HE2	2:E:28:PHE:CD2	2.29	0.67
1:G:116:TRP:O	2:H:89:TRP:HA	1.95	0.67
2:N:13:TYR:CE1	2:N:121:LYS:HG3	2.29	0.67
1:P:5:LEU:HB3	1:P:6:PRO:HD2	1.77	0.67
2:E:14:ARG:NH1	3:O:359:GLU:OE2	2.28	0.66
3:O:260:ARG:NH2	4:O:403:SO4:O4	2.26	0.66
1:D:79:GLN:O	1:D:101:SER:OG	2.10	0.66
1:D:31:PHE:O	1:D:35:GLN:HG2	1.96	0.66
2:E:61:THR:CG2	2:E:100:VAL:HB	2.26	0.65
3:O:231:THR:OG1	3:O:234:GLU:HB2	1.96	0.65
1:A:31:PHE:O	1:A:35:GLN:HG2	1.96	0.65
2:Q:90:LEU:HD12	2:Q:90:LEU:C	2.22	0.65
2:B:88:VAL:HG12	2:B:88:VAL:O	1.95	0.65
3:F:322:THR:O	3:F:326:ILE:HG12	1.96	0.65
3:F:344:VAL:HG13	3:F:349:ALA:HB3	1.78	0.65
1:A:87:MET:SD	2:B:40:SER:OG	2.48	0.65
1:A:115:LYS:HA	2:B:88:VAL:HG12	1.78	0.64
1:P:52:PHE:CD1	1:P:52:PHE:C	2.75	0.64
1:G:72:LEU:HD11	2:H:76:TRP:HB3	1.78	0.64
1:J:31:PHE:O	1:J:35:GLN:HG2	1.98	0.64
3:R:204:ILE:HG12	3:R:235:MET:CE	2.27	0.64
3:L:253:PHE:CZ	3:L:292:LEU:HG	2.33	0.64
1:M:20:ASN:HD21	1:M:61:ILE:HA	1.62	0.64
1:A:25:TRP:CZ2	1:A:73:ARG:HD3	2.34	0.62
1:P:29:GLU:CD	2:Q:79:SER:HG	2.06	0.62
2:H:88:VAL:HG12	2:H:88:VAL:O	1.98	0.62
2:N:90:LEU:C	2:N:90:LEU:HD12	2.24	0.62
2:N:32:GLN:HE21	2:N:32:GLN:HA	1.64	0.61
3:L:362:PHE:N	3:L:362:PHE:HD1	1.98	0.61
1:M:69:TRP:NE1	1:M:125:ASN:HB2	2.16	0.61
3:L:362:PHE:N	3:L:362:PHE:CD1	2.68	0.61
1:D:17:GLN:HE21	1:D:17:GLN:HA	1.63	0.61
3:I:194:PHE:CD1	3:I:350:LEU:HD22	2.35	0.61
1:G:87:MET:O	1:G:88:SER:HB3	2.00	0.60
2:Q:13:TYR:CE1	2:Q:121:LYS:HG3	2.36	0.60
1:A:30:ARG:O	1:A:33:THR:OG1	2.17	0.60
3:F:288:ASP:HB2	3:F:291:MET:HE3	1.82	0.60
3:C:253:PHE:CZ	3:C:292:LEU:HG	2.36	0.60
3:C:231:THR:OG1	3:C:234:GLU:HB2	2.01	0.60
3:C:288:ASP:HB2	3:C:291:MET:HE3	1.82	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:239:THR:O	3:C:242:THR:HG22	2.01	0.59
1:D:72:LEU:HD11	2:E:76:TRP:HB3	1.84	0.59
1:D:116:TRP:O	2:E:89:TRP:HA	2.01	0.59
3:I:349:ALA:O	3:I:352:GLU:HG2	2.01	0.59
1:J:116:TRP:O	2:K:89:TRP:HA	2.01	0.59
1:G:31:PHE:O	1:G:35:GLN:HG2	2.01	0.59
3:I:288:ASP:HB2	3:I:291:MET:HE3	1.83	0.59
3:F:194:PHE:CD1	3:F:350:LEU:HD22	2.38	0.58
1:J:87:MET:SD	2:K:40:SER:OG	2.59	0.58
1:A:43:SER:HB3	2:B:80:ASP:OD2	2.03	0.58
1:D:87:MET:O	1:D:88:SER:HB3	2.02	0.58
2:E:90:LEU:C	2:E:90:LEU:CD1	2.76	0.58
2:N:1:CYS:SG	2:N:7:SER:HB3	2.43	0.58
2:H:88:VAL:O	2:H:88:VAL:CG1	2.51	0.58
1:J:102:GLN:OE1	2:K:91:ARG:NH2	2.27	0.58
1:A:116:TRP:O	2:B:89:TRP:HA	2.04	0.58
2:H:16:PHE:HB3	2:H:18:GLU:OE1	2.04	0.57
1:M:70:ILE:HG22	2:N:78:TRP:HZ3	1.70	0.57
2:E:88:VAL:HG12	2:E:88:VAL:O	2.02	0.57
1:P:89:ALA:HB1	2:Q:42:HIS:NE2	2.19	0.57
1:A:75:ARG:HD2	2:B:77:GLU:OE2	2.04	0.57
1:J:43:SER:HB3	2:K:80:ASP:OD2	2.05	0.57
1:P:12:ASP:CG	1:P:13:GLN:H	2.12	0.57
3:C:184:SER:HB2	3:C:283:ASP:OD2	2.04	0.57
2:B:88:VAL:O	2:B:88:VAL:CG1	2.53	0.57
2:E:28:PHE:CE1	2:E:32:GLN:NE2	2.73	0.57
1:P:29:GLU:OE1	2:Q:79:SER:OG	2.22	0.57
1:A:87:MET:SD	2:B:40:SER:CB	2.93	0.56
3:I:172:LEU:HD12	3:I:207:THR:O	2.05	0.56
1:M:5:LEU:HB3	1:M:6:PRO:HD2	1.88	0.56
2:N:32:GLN:HE21	2:N:32:GLN:N	2.04	0.56
3:R:218:ASN:ND2	3:R:249:LEU:HD12	2.19	0.56
1:D:84:GLU:O	1:D:91:ILE:HD11	2.05	0.56
2:E:61:THR:HG21	2:E:100:VAL:HB	1.87	0.56
3:F:349:ALA:O	3:F:352:GLU:HG2	2.05	0.56
3:C:322:THR:O	3:C:326:ILE:HD13	2.05	0.56
3:C:329:ILE:N	3:C:329:ILE:HD13	2.21	0.56
3:L:288:ASP:HB2	3:L:291:MET:HE3	1.87	0.56
2:H:28:PHE:CD1	2:H:28:PHE:C	2.83	0.56
2:N:32:GLN:HA	2:N:32:GLN:NE2	2.20	0.55
1:A:115:LYS:HA	2:B:88:VAL:CG1	2.37	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:87:MET:HE3	2:K:38:LEU:O	2.05	0.55
1:P:31:PHE:O	1:P:35:GLN:HG2	2.07	0.55
3:L:239:THR:O	3:L:242:THR:HG22	2.07	0.55
1:J:72:LEU:HD11	2:K:76:TRP:HB3	1.89	0.55
2:E:16:PHE:HB3	2:E:18:GLU:OE1	2.06	0.55
1:J:87:MET:O	1:J:88:SER:HB3	2.07	0.55
2:H:36:SER:HA	2:H:121:LYS:O	2.07	0.54
2:E:90:LEU:HD12	2:E:91:ARG:N	2.21	0.54
3:F:330:LYS:O	3:F:333:ALA:HB3	2.08	0.54
1:M:89:ALA:HB1	2:N:42:HIS:CE1	2.43	0.54
1:A:110:TRP:HA	3:C:245:TYR:CE2	2.43	0.54
2:K:65:LEU:HD21	2:K:98:MET:HE3	1.89	0.54
2:Q:34:LYS:N	2:Q:34:LYS:HD2	2.23	0.54
1:A:25:TRP:CH2	1:A:73:ARG:HD3	2.43	0.54
3:F:204:ILE:HA	3:F:209:THR:O	2.08	0.54
2:E:92:ARG:HB2	2:E:93:PRO:CD	2.39	0.53
1:G:87:MET:SD	2:H:40:SER:OG	2.66	0.53
3:O:204:ILE:HA	3:O:209:THR:O	2.08	0.53
3:F:172:LEU:HD12	3:F:207:THR:O	2.09	0.53
1:D:69:TRP:NE1	1:D:125:ASN:HB2	2.23	0.53
2:N:40:SER:O	2:N:41:ILE:HD13	2.09	0.53
1:P:12:ASP:OD1	1:P:13:GLN:N	2.41	0.53
2:Q:61:THR:CG2	2:Q:100:VAL:HB	2.38	0.53
1:G:94:VAL:HA	2:H:107:TRP:CZ3	2.44	0.53
3:F:178:VAL:HA	3:F:214:ILE:O	2.08	0.53
1:P:20:ASN:ND2	1:P:61:ILE:HD12	2.24	0.53
1:P:28:ALA:O	1:P:31:PHE:HB3	2.08	0.53
3:L:231:THR:OG1	3:L:234:GLU:HB2	2.08	0.53
1:P:20:ASN:O	1:P:22:PRO:HD3	2.09	0.53
1:P:69:TRP:HA	1:P:69:TRP:CE3	2.44	0.53
1:M:87:MET:HG3	1:M:89:ALA:H	1.74	0.52
1:P:11:TYR:HB2	1:P:52:PHE:CE2	2.45	0.52
1:A:87:MET:SD	2:B:40:SER:HB3	2.49	0.52
1:M:35:GLN:O	1:M:36:ALA:HB3	2.09	0.52
2:N:54:ALA:HB1	2:N:63:MET:HE1	1.90	0.52
2:E:1:CYS:SG	2:E:7:SER:HB2	2.49	0.52
2:H:30:TYR:HA	2:H:36:SER:O	2.09	0.52
1:J:110:TRP:HA	3:L:245:TYR:CE2	2.45	0.52
3:L:322:THR:O	3:L:326:ILE:HD13	2.08	0.52
1:A:8:TRP:CZ3	1:A:131:PHE:CE2	2.97	0.52
1:G:69:TRP:NE1	1:G:125:ASN:HB2	2.25	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:O:239:THR:O	3:O:242:THR:HG22	2.09	0.52
2:B:44:ARG:O	2:B:47:GLU:HB3	2.09	0.52
1:A:87:MET:O	1:A:88:SER:CB	2.57	0.52
2:E:14:ARG:HG2	2:E:16:PHE:CE1	2.45	0.52
3:F:282:THR:HG22	3:F:284:GLY:H	1.75	0.52
1:M:33:THR:HA	1:M:39:GLY:O	2.10	0.52
1:P:20:ASN:HD21	1:P:61:ILE:HA	1.74	0.52
1:P:74:ASP:OD2	2:Q:75:LYS:N	2.43	0.52
3:R:283:ASP:CG	3:R:283:ASP:O	2.53	0.51
1:D:89:ALA:HB1	2:E:42:HIS:NE2	2.24	0.51
1:A:72:LEU:HD11	2:B:76:TRP:HB3	1.91	0.51
2:E:88:VAL:O	2:E:88:VAL:CG1	2.58	0.51
1:P:52:PHE:HE1	1:P:56:LEU:HD12	1.75	0.51
3:F:204:ILE:HB	3:F:232:LYS:CG	2.40	0.51
1:G:24:THR:HG22	1:G:26:ASP:H	1.75	0.51
3:R:288:ASP:HB2	3:R:291:MET:HE3	1.92	0.51
1:D:94:VAL:HA	2:E:107:TRP:CZ3	2.46	0.51
1:M:55:GLN:OE1	3:R:207:THR:HG22	2.11	0.51
3:O:227:ASN:HD22	3:O:269:GLY:CA	2.24	0.51
3:I:178:VAL:HA	3:I:214:ILE:O	2.10	0.51
3:I:231:THR:HB	3:I:234:GLU:HB2	1.93	0.51
3:L:362:PHE:HD1	3:L:362:PHE:H	1.57	0.51
3:I:330:LYS:O	3:I:333:ALA:HB3	2.11	0.50
3:O:344:VAL:CG1	3:O:346:ASP:O	2.60	0.50
2:Q:65:LEU:HD21	2:Q:98:MET:HB2	1.93	0.50
2:B:65:LEU:HD21	2:B:98:MET:HE3	1.93	0.50
2:Q:88:VAL:CG1	2:Q:88:VAL:O	2.59	0.50
2:B:32:GLN:O	2:B:33:HIS:HB2	2.12	0.50
2:N:20:LYS:O	2:N:115:THR:HA	2.12	0.50
2:N:61:THR:CG2	2:N:100:VAL:HB	2.42	0.50
3:O:188:TRP:CD2	3:O:244:GLN:HB2	2.47	0.50
3:R:239:THR:O	3:R:242:THR:HG22	2.12	0.50
3:F:218:ASN:HD21	3:F:248:ASP:H	1.58	0.50
3:O:233:GLU:O	3:O:237:VAL:HG23	2.12	0.50
2:N:88:VAL:CG1	2:N:88:VAL:O	2.59	0.50
1:P:29:GLU:OE2	2:Q:79:SER:OG	2.28	0.50
1:M:79:GLN:O	1:M:101:SER:OG	2.25	0.49
3:L:200:GLN:OE1	3:L:232:LYS:HE3	2.11	0.49
3:C:344:VAL:HG13	3:C:349:ALA:HB3	1.94	0.49
1:M:29:GLU:OE1	2:N:79:SER:CB	2.60	0.49
1:M:131:PHE:HD1	1:M:132:PRO:O	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:I:264:TYR:O	3:I:270:GLY:HA3	2.13	0.49
2:H:90:LEU:HD12	2:H:91:ARG:N	2.27	0.49
1:P:116:TRP:O	2:Q:89:TRP:HA	2.13	0.49
3:R:344:VAL:CG1	3:R:346:ASP:O	2.60	0.49
1:P:69:TRP:NE1	1:P:125:ASN:HB2	2.28	0.49
2:E:9:ASN:CG	2:E:9:ASN:O	2.55	0.49
1:G:64:PRO:HD2	1:G:65:GLU:OE1	2.12	0.49
2:K:68:ASN:ND2	2:K:95:CYS:SG	2.86	0.49
1:P:61:ILE:O	1:P:61:ILE:HG23	2.12	0.49
2:H:9:ASN:CG	2:H:9:ASN:O	2.55	0.48
1:M:80:GLN:O	1:M:80:GLN:HG2	2.12	0.48
2:E:28:PHE:CD1	2:E:28:PHE:C	2.90	0.48
1:M:20:ASN:O	1:M:22:PRO:HD3	2.13	0.48
2:E:97:VAL:O	2:E:107:TRP:HA	2.14	0.48
2:N:65:LEU:HD21	2:N:98:MET:HB2	1.96	0.48
3:R:344:VAL:HG13	3:R:349:ALA:HB3	1.94	0.48
1:G:79:GLN:O	1:G:101:SER:OG	2.24	0.48
3:I:239:THR:O	3:I:242:THR:HG22	2.13	0.48
1:M:69:TRP:C	1:M:70:ILE:HG13	2.37	0.48
2:H:90:LEU:C	2:H:90:LEU:CD1	2.81	0.48
1:J:87:MET:HE2	1:J:87:MET:HB3	1.79	0.48
2:K:33:HIS:HD2	2:K:122:PHE:O	1.96	0.48
1:P:94:VAL:HG13	1:P:96:TRP:CD1	2.49	0.48
2:H:63:MET:HB2	2:H:118:PHE:HA	1.95	0.48
2:K:68:ASN:HA	2:K:95:CYS:SG	2.54	0.48
2:K:116:VAL:C	2:K:117:SER:O	2.55	0.48
1:M:74:ASP:HB3	1:M:103:MET:HE3	1.96	0.48
3:O:200:GLN:NE2	3:O:236:ILE:CD1	2.77	0.48
3:R:322:THR:O	3:R:323:LYS:C	2.56	0.48
1:J:89:ALA:HB1	2:K:42:HIS:CE1	2.49	0.47
1:P:12:ASP:CG	1:P:13:GLN:N	2.70	0.47
1:D:43:SER:HA	2:E:78:TRP:CZ3	2.48	0.47
3:F:337:THR:O	3:F:341:PHE:HB3	2.14	0.47
2:K:44:ARG:O	2:K:47:GLU:HB3	2.15	0.47
3:O:288:ASP:HB2	3:O:291:MET:HE3	1.96	0.47
3:C:218:ASN:ND2	3:C:249:LEU:HD12	2.29	0.47
1:D:93:TYR:CE2	1:J:77:LYS:HB2	2.50	0.47
3:I:180:ASP:HB3	3:I:282:THR:HA	1.96	0.47
3:O:292:LEU:C	3:O:292:LEU:HD23	2.39	0.47
1:J:8:TRP:CZ3	1:J:131:PHE:CE2	3.02	0.47
3:O:322:THR:O	3:O:323:LYS:C	2.57	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:P:29:GLU:OE1	2:Q:79:SER:CB	2.62	0.47
1:J:18:ALA:HB1	1:J:61:ILE:HD11	1.97	0.47
1:P:49:GLU:O	1:P:53:VAL:HG23	2.14	0.47
3:R:254:GLY:O	3:R:257:GLN:HB3	2.15	0.47
1:D:5:LEU:HB3	1:D:6:PRO:HD2	1.97	0.47
2:H:20:LYS:O	2:H:115:THR:HA	2.15	0.47
1:J:87:MET:O	1:J:88:SER:CB	2.62	0.47
3:L:218:ASN:ND2	3:L:249:LEU:HD12	2.30	0.47
1:M:99:GLY:HA2	1:M:120:ASP:HB2	1.96	0.47
1:P:11:TYR:HB2	1:P:52:PHE:CD2	2.50	0.46
2:Q:30:TYR:HA	2:Q:36:SER:O	2.15	0.46
2:Q:32:GLN:CA	2:Q:32:GLN:HE21	2.27	0.46
1:M:78:GLU:OE1	1:M:78:GLU:N	2.44	0.46
3:O:227:ASN:ND2	3:O:269:GLY:CA	2.78	0.46
2:H:28:PHE:CE1	2:H:32:GLN:OE1	2.68	0.46
2:B:34:LYS:HE3	2:B:34:LYS:HA	1.98	0.46
1:G:3:ASN:N	1:G:3:ASN:ND2	2.64	0.46
3:O:200:GLN:NE2	3:O:236:ILE:HD13	2.31	0.46
2:B:68:ASN:HA	2:B:95:CYS:SG	2.55	0.46
3:F:307:PHE:HA	3:F:340:TYR:O	2.15	0.46
2:H:54:ALA:HB2	2:H:63:MET:HE1	1.97	0.46
1:A:18:ALA:HB1	1:A:61:ILE:HD11	1.98	0.46
3:F:204:ILE:HB	3:F:232:LYS:HG3	1.97	0.46
3:F:216:TYR:CD2	3:F:216:TYR:C	2.94	0.46
3:I:337:THR:O	3:I:341:PHE:HB3	2.16	0.46
1:M:11:TYR:HB2	1:M:52:PHE:CE2	2.50	0.46
3:O:325:LEU:O	3:O:328:GLU:N	2.49	0.46
3:O:344:VAL:HG13	3:O:349:ALA:HB3	1.98	0.46
3:R:216:TYR:CE2	3:R:252:THR:HA	2.50	0.46
2:E:63:MET:HB2	2:E:118:PHE:HA	1.98	0.46
3:L:184:SER:HB2	3:L:283:ASP:OD2	2.15	0.46
1:M:61:ILE:HG23	1:M:61:ILE:O	2.15	0.46
1:A:87:MET:HE3	2:B:38:LEU:O	2.16	0.46
1:A:97:ASN:HB2	2:B:106:PHE:HB2	1.96	0.46
1:M:8:TRP:CZ3	1:M:131:PHE:CE2	3.04	0.46
1:M:23:LYS:O	1:M:124:LYS:HA	2.16	0.46
2:B:84:LEU:HG	2:B:84:LEU:O	2.16	0.45
2:N:30:TYR:HA	2:N:36:SER:O	2.16	0.45
3:O:210:GLN:OE1	3:O:271:ARG:NE	2.49	0.45
3:O:227:ASN:HD22	3:O:269:GLY:HA2	1.81	0.45
1:J:40:HIS:CD2	2:K:80:ASP:HA	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:89:ALA:HB1	2:H:42:HIS:NE2	2.31	0.45
2:N:62:SER:HA	2:N:98:MET:O	2.17	0.45
1:G:87:MET:HG3	1:G:89:ALA:H	1.80	0.45
1:M:25:TRP:NE1	1:M:70:ILE:O	2.47	0.45
3:R:325:LEU:O	3:R:328:GLU:N	2.46	0.45
3:I:292:LEU:O	3:I:296:ILE:HG12	2.16	0.45
1:M:116:TRP:O	2:N:89:TRP:HA	2.16	0.45
1:P:99:GLY:HA2	1:P:120:ASP:HB2	1.98	0.45
1:A:67:TYR:HE2	1:A:117:ASP:HB3	1.81	0.45
1:J:11:TYR:O	1:J:12:ASP:C	2.59	0.45
1:M:49:GLU:O	1:M:53:VAL:HG23	2.17	0.45
1:M:69:TRP:HA	1:M:69:TRP:CE3	2.52	0.45
1:A:75:ARG:C	1:A:76:ARG:O	2.60	0.45
2:K:36:SER:HA	2:K:121:LYS:O	2.17	0.45
2:K:3:LEU:HD23	2:K:4:HIS:CE1	2.51	0.45
2:K:8:TYR:HB2	2:K:49:PHE:CE2	2.52	0.45
3:O:200:GLN:HE22	3:O:236:ILE:HD13	1.82	0.45
3:R:180:ASP:HB2	3:R:282:THR:HA	1.98	0.45
3:R:231:THR:HB	3:R:234:GLU:HB2	1.99	0.45
2:B:36:SER:HA	2:B:121:LYS:O	2.16	0.45
3:I:296:ILE:O	3:I:299:CYS:HB2	2.17	0.45
1:G:84:GLU:O	1:G:91:ILE:HD11	2.16	0.44
1:J:82:SER:HA	2:K:73:GLU:OE1	2.16	0.44
1:M:94:VAL:HG13	1:M:96:TRP:CD1	2.51	0.44
2:Q:108:PHE:CE1	3:R:318:ASN:HB3	2.52	0.44
2:B:71:TRP:HA	2:B:74:CYS:SG	2.57	0.44
2:H:97:VAL:O	2:H:107:TRP:HA	2.16	0.44
3:I:292:LEU:C	3:I:292:LEU:HD23	2.42	0.44
1:J:71:GLY:O	2:K:78:TRP:HA	2.17	0.44
2:Q:90:LEU:HD12	2:Q:91:ARG:N	2.31	0.44
2:B:30:TYR:HA	2:B:36:SER:O	2.17	0.44
3:O:325:LEU:O	3:O:326:ILE:C	2.59	0.44
2:E:30:TYR:HA	2:E:36:SER:O	2.17	0.44
2:E:86:TYR:CE2	2:E:88:VAL:HG21	2.52	0.44
1:A:110:TRP:HA	3:C:245:TYR:CD2	2.52	0.44
1:M:28:ALA:O	1:M:31:PHE:HB3	2.18	0.44
1:M:42:VAL:HB	1:M:128:VAL:HG12	1.99	0.44
2:Q:29:CYS:O	2:Q:32:GLN:HB2	2.18	0.44
2:E:36:SER:HA	2:E:121:LYS:O	2.16	0.44
1:P:42:VAL:HG12	1:P:128:VAL:O	2.18	0.44
3:R:283:ASP:OD1	3:R:283:ASP:C	2.61	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:42:VAL:HB	1:A:128:VAL:HG12	2.00	0.44
1:A:43:SER:CB	2:B:80:ASP:OD2	2.65	0.44
1:D:94:VAL:HA	2:E:107:TRP:CH2	2.53	0.44
1:D:39:GLY:O	1:D:40:HIS:CD2	2.71	0.43
2:E:59:LYS:HG2	2:E:60:TYR:CE2	2.53	0.43
2:K:21:THR:HA	2:K:114:LYS:O	2.18	0.43
2:E:54:ALA:CB	2:E:63:MET:HE1	2.46	0.43
2:N:92:ARG:HB2	2:N:93:PRO:CD	2.48	0.43
2:B:16:PHE:CE1	2:B:28:PHE:HE2	2.36	0.43
1:G:5:LEU:HB3	1:G:6:PRO:HD2	2.01	0.43
2:H:1:CYS:SG	2:H:7:SER:HB2	2.58	0.43
3:L:292:LEU:HD23	3:L:292:LEU:C	2.44	0.43
3:F:239:THR:O	3:F:242:THR:HG22	2.19	0.43
1:M:65:GLU:OE2	1:M:126:PRO:HD3	2.18	0.43
3:O:218:ASN:ND2	3:O:249:LEU:HD12	2.33	0.43
1:P:25:TRP:O	1:P:25:TRP:CE3	2.72	0.43
1:P:74:ASP:HB3	1:P:103:MET:HE3	1.99	0.43
2:Q:40:SER:O	2:Q:41:ILE:HD13	2.19	0.43
1:A:87:MET:O	1:A:88:SER:HB2	2.18	0.43
1:D:64:PRO:HD2	1:D:65:GLU:OE1	2.18	0.43
1:J:43:SER:CB	2:K:80:ASP:OD2	2.66	0.43
3:I:339:ARG:O	3:I:360:GLN:NE2	2.52	0.43
2:N:108:PHE:CE1	3:O:318:ASN:HB3	2.54	0.43
3:O:283:ASP:HA	3:O:311:VAL:O	2.17	0.43
1:P:8:TRP:CZ3	1:P:131:PHE:CE2	3.06	0.43
3:R:325:LEU:O	3:R:326:ILE:C	2.61	0.43
2:H:92:ARG:HB2	2:H:93:PRO:CD	2.49	0.43
3:I:217:ALA:C	3:I:249:LEU:O	2.62	0.43
2:K:84:LEU:O	2:K:84:LEU:HG	2.19	0.43
3:O:216:TYR:CE2	3:O:252:THR:HA	2.54	0.43
3:R:253:PHE:CE1	3:R:292:LEU:HG	2.53	0.43
3:L:192:LYS:O	3:L:196:GLU:HG3	2.18	0.43
3:O:312:LEU:HG	3:O:344:VAL:O	2.19	0.43
3:R:188:TRP:CD2	3:R:244:GLN:HB2	2.54	0.43
1:J:110:TRP:HA	3:L:245:TYR:CD2	2.54	0.42
1:M:67:TYR:HA	1:M:107:LEU:O	2.19	0.42
2:N:67:LEU:HD12	2:N:67:LEU:HA	1.89	0.42
3:O:312:LEU:O	3:O:313:GLY:C	2.61	0.42
3:O:344:VAL:HG12	3:O:346:ASP:O	2.19	0.42
3:L:173:ILE:O	3:L:271:ARG:NH2	2.52	0.42
1:A:23:LYS:O	1:A:124:LYS:HA	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:107:TRP:N	2:E:107:TRP:CD1	2.86	0.42
3:F:204:ILE:HB	3:F:232:LYS:HG2	2.01	0.42
2:N:63:MET:HB2	2:N:118:PHE:HA	2.00	0.42
2:N:71:TRP:HA	2:N:74:CYS:SG	2.60	0.42
2:Q:90:LEU:C	2:Q:90:LEU:CD1	2.92	0.42
3:R:305:LEU:HD12	3:R:305:LEU:HA	1.78	0.42
3:C:277:VAL:HG22	3:C:305:LEU:HB3	2.01	0.42
2:H:28:PHE:CD1	2:H:28:PHE:O	2.72	0.42
3:L:175:VAL:O	3:L:211:VAL:HA	2.19	0.42
3:C:339:ARG:NH2	7:C:403:CL:CL	2.89	0.42
2:N:13:TYR:OH	2:N:46:GLU:OE2	2.33	0.42
1:P:52:PHE:C	1:P:52:PHE:HD1	2.26	0.42
2:Q:1:CYS:SG	2:Q:7:SER:HB3	2.59	0.42
3:R:233:GLU:O	3:R:237:VAL:HG23	2.20	0.42
1:J:131:PHE:HD1	1:J:132:PRO:O	2.02	0.42
2:N:54:ALA:HB2	2:N:63:MET:HE1	1.98	0.42
1:G:43:SER:HA	2:H:78:TRP:CE3	2.55	0.42
3:I:307:PHE:HA	3:I:340:TYR:O	2.20	0.42
1:M:35:GLN:H	1:M:35:GLN:HG2	1.60	0.42
1:M:99:GLY:HA2	1:M:120:ASP:CG	2.45	0.42
2:Q:70:PRO:HG2	2:Q:109:ASN:ND2	2.35	0.42
3:F:179:CYS:O	3:F:215:GLN:HA	2.20	0.42
3:F:206:PRO:O	3:F:271:ARG:NH2	2.53	0.42
1:G:94:VAL:HA	2:H:107:TRP:CH2	2.54	0.42
1:P:131:PHE:HD1	1:P:132:PRO:O	2.03	0.42
1:G:71:GLY:O	2:H:78:TRP:HA	2.20	0.41
1:M:42:VAL:HG13	1:M:70:ILE:HG23	2.02	0.41
1:A:11:TYR:O	1:A:12:ASP:C	2.63	0.41
1:P:33:THR:HA	1:P:39:GLY:O	2.20	0.41
2:Q:107:TRP:CD1	2:Q:107:TRP:N	2.87	0.41
3:C:322:THR:O	3:C:326:ILE:CD1	2.68	0.41
2:H:26:GLU:HA	2:H:26:GLU:OE1	2.21	0.41
1:J:24:THR:HG22	1:J:26:ASP:H	1.86	0.41
3:L:277:VAL:HG22	3:L:305:LEU:HB3	2.01	0.41
3:O:172:LEU:O	3:O:274:ALA:HA	2.20	0.41
3:R:344:VAL:HG12	3:R:346:ASP:O	2.20	0.41
1:A:89:ALA:HB1	2:B:42:HIS:CE1	2.55	0.41
3:F:264:TYR:O	3:F:270:GLY:HA3	2.21	0.41
1:G:87:MET:O	1:G:88:SER:CB	2.68	0.41
3:L:188:TRP:CD2	3:L:244:GLN:HB2	2.56	0.41
3:L:289:GLY:O	3:L:292:LEU:HB2	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:44:ILE:HD11	1:D:53:VAL:HG11	2.02	0.41
1:J:18:ALA:HB1	1:J:61:ILE:CD1	2.51	0.41
1:J:97:ASN:HB2	2:K:106:PHE:HB2	2.02	0.41
3:L:289:GLY:C	3:L:291:MET:N	2.79	0.41
2:N:90:LEU:C	2:N:90:LEU:CD1	2.93	0.41
2:N:97:VAL:O	2:N:107:TRP:HA	2.20	0.41
3:C:357:LEU:O	3:C:361:ILE:HG12	2.20	0.41
3:F:218:ASN:H	3:F:218:ASN:ND2	2.18	0.41
3:I:172:LEU:CD1	3:I:207:THR:O	2.69	0.41
3:I:314:TYR:O	3:I:318:ASN:ND2	2.49	0.41
1:M:74:ASP:OD2	2:N:75:LYS:N	2.54	0.41
1:A:5:LEU:HB3	1:A:6:PRO:HD2	2.03	0.41
1:G:69:TRP:CD1	1:G:125:ASN:HB2	2.56	0.41
2:H:21:THR:HG23	2:H:112:CYS:O	2.19	0.41
3:R:181:GLU:HG3	3:R:181:GLU:O	2.20	0.41
3:F:173:ILE:HD12	3:F:361:ILE:CG2	2.50	0.41
3:I:356:THR:HG22	3:I:360:GLN:OE1	2.21	0.41
3:L:283:ASP:OD1	3:L:283:ASP:C	2.64	0.41
3:O:174:ASP:OD2	3:O:274:ALA:HB3	2.21	0.41
3:O:253:PHE:CE1	3:O:292:LEU:HG	2.55	0.41
1:M:37:LYS:O	1:M:38:ARG:HB2	2.21	0.41
2:B:116:VAL:C	2:B:117:SER:O	2.63	0.40
1:G:85:TRP:HE3	2:H:66:GLY:O	2.04	0.40
3:I:349:ALA:O	3:I:350:LEU:C	2.64	0.40
1:J:89:ALA:HB1	2:K:42:HIS:NE2	2.35	0.40
2:B:13:TYR:CD2	2:B:49:PHE:HE2	2.39	0.40
2:E:86:TYR:CZ	2:E:88:VAL:HG21	2.56	0.40
3:I:296:ILE:O	3:I:299:CYS:N	2.54	0.40
2:Q:59:LYS:H	2:Q:59:LYS:HE3	1.85	0.40
2:N:3:LEU:O	2:N:4:HIS:HB2	2.21	0.40
2:N:90:LEU:HD12	2:N:91:ARG:N	2.37	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	129/135 (96%)	112 (87%)	14 (11%)	3 (2%)	5	29
1	D	129/135 (96%)	119 (92%)	8 (6%)	2 (2%)	7	36
1	G	129/135 (96%)	115 (89%)	13 (10%)	1 (1%)	16	50
1	J	129/135 (96%)	111 (86%)	14 (11%)	4 (3%)	3	22
1	M	129/135 (96%)	111 (86%)	16 (12%)	2 (2%)	7	36
1	P	129/135 (96%)	114 (88%)	13 (10%)	2 (2%)	7	36
2	B	120/124 (97%)	104 (87%)	12 (10%)	4 (3%)	3	21
2	E	120/124 (97%)	102 (85%)	16 (13%)	2 (2%)	7	35
2	H	120/124 (97%)	106 (88%)	11 (9%)	3 (2%)	4	27
2	K	120/124 (97%)	101 (84%)	16 (13%)	3 (2%)	4	27
2	N	120/124 (97%)	101 (84%)	15 (12%)	4 (3%)	3	21
2	Q	120/124 (97%)	107 (89%)	7 (6%)	6 (5%)	1	13
3	C	189/217 (87%)	176 (93%)	12 (6%)	1 (0%)	24	59
3	F	189/217 (87%)	176 (93%)	12 (6%)	1 (0%)	24	59
3	I	189/217 (87%)	173 (92%)	13 (7%)	3 (2%)	7	36
3	L	189/217 (87%)	175 (93%)	11 (6%)	3 (2%)	7	36
3	O	189/217 (87%)	171 (90%)	16 (8%)	2 (1%)	11	43
3	R	189/217 (87%)	172 (91%)	13 (7%)	4 (2%)	5	31
All	All	2628/2856 (92%)	2346 (89%)	232 (9%)	50 (2%)	6	33

All (50) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	E	2	PRO
1	G	88	SER
2	H	2	PRO
1	M	88	SER
2	N	117	SER
1	P	88	SER
2	Q	3	LEU
2	Q	117	SER
1	A	76	ARG
1	A	88	SER
2	B	33	HIS

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Mol	Chain	Res	Type
3	C	313	GLY
1	D	88	SER
3	F	334	SER
3	I	286	SER
3	I	334	SER
1	J	88	SER
1	J	89	ALA
2	K	117	SER
3	L	313	GLY
2	N	2	PRO
3	O	313	GLY
2	Q	9	ASN
1	A	89	ALA
2	B	117	SER
3	L	217	ALA
1	M	36	ALA
1	P	76	ARG
2	B	2	PRO
1	D	132	PRO
2	H	117	SER
1	J	24	THR
1	J	76	ARG
3	L	290	SER
2	N	9	ASN
2	N	33	HIS
2	Q	33	HIS
2	B	83	LYS
3	I	217	ALA
2	K	2	PRO
2	K	83	LYS
3	R	217	ALA
3	R	323	LYS
2	E	111	GLY
3	R	326	ILE
2	H	88	VAL
3	O	326	ILE
2	Q	2	PRO
2	Q	88	VAL
3	R	313	GLY

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	114/118 (97%)	109 (96%)	5 (4%)	25	59
1	D	114/118 (97%)	103 (90%)	11 (10%)	8	32
1	G	114/118 (97%)	106 (93%)	8 (7%)	14	45
1	J	114/118 (97%)	111 (97%)	3 (3%)	40	69
1	M	114/118 (97%)	108 (95%)	6 (5%)	20	53
1	P	114/118 (97%)	105 (92%)	9 (8%)	11	40
2	B	110/112 (98%)	103 (94%)	7 (6%)	16	48
2	E	110/112 (98%)	101 (92%)	9 (8%)	10	39
2	H	110/112 (98%)	104 (94%)	6 (6%)	19	52
2	K	110/112 (98%)	100 (91%)	10 (9%)	9	34
2	N	110/112 (98%)	103 (94%)	7 (6%)	16	48
2	Q	110/112 (98%)	100 (91%)	10 (9%)	9	34
3	C	160/181 (88%)	147 (92%)	13 (8%)	11	39
3	F	160/181 (88%)	150 (94%)	10 (6%)	16	48
3	I	160/181 (88%)	150 (94%)	10 (6%)	16	48
3	L	160/181 (88%)	152 (95%)	8 (5%)	22	55
3	O	160/181 (88%)	154 (96%)	6 (4%)	29	62
3	R	160/181 (88%)	155 (97%)	5 (3%)	35	66
All	All	2304/2466 (93%)	2161 (94%)	143 (6%)	16	49

All (143) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	58	THR
1	A	87	MET
1	A	90	SER
1	A	98	THR
1	A	118	TYR
2	B	32	GLN

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Mol	Chain	Res	Type
2	B	34	LYS
2	B	55	SER
2	B	67	LEU
2	B	91	ARG
2	B	117	SER
2	B	122	PHE
3	C	172	LEU
3	C	184	SER
3	C	208	LYS
3	C	273	SER
3	C	303	ASN
3	C	312	LEU
3	C	329	ILE
3	C	334	SER
3	C	337	THR
3	C	339	ARG
3	C	343	ASN
3	C	344	VAL
3	C	362	PHE
1	D	13	GLN
1	D	17	GLN
1	D	21	GLU
1	D	30	ARG
1	D	77	LYS
1	D	83	SER
1	D	86	SER
1	D	88	SER
1	D	90	SER
1	D	92	ILE
1	D	116	TRP
2	E	7	SER
2	E	32	GLN
2	E	55	SER
2	E	59	LYS
2	E	61	THR
2	E	67	LEU
2	E	81	ASP
2	E	117	SER
2	E	122	PHE
3	F	184	SER
3	F	204	ILE
3	F	207	THR

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Mol	Chain	Res	Type
3	F	231	THR
3	F	265	SER
3	F	321	ASP
3	F	339	ARG
3	F	344	VAL
3	F	345	SER
3	F	350	LEU
1	G	3	ASN
1	G	30	ARG
1	G	83	SER
1	G	84	GLU
1	G	90	SER
1	G	92	ILE
1	G	116	TRP
1	G	118	TYR
2	H	7	SER
2	H	55	SER
2	H	59	LYS
2	H	63	MET
2	H	81	ASP
2	H	122	PHE
3	I	184	SER
3	I	192	LYS
3	I	202	LEU
3	I	232	LYS
3	I	283	ASP
3	I	321	ASP
3	I	343	ASN
3	I	344	VAL
3	I	345	SER
3	I	350	LEU
1	J	3	ASN
1	J	90	SER
1	J	119	SER
2	K	3	LEU
2	K	17	SER
2	K	32	GLN
2	K	37	ARG
2	K	52	LYS
2	K	79	SER
2	K	81	ASP
2	K	91	ARG

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Mol	Chain	Res	Type
2	K	117	SER
2	K	122	PHE
3	L	184	SER
3	L	273	SER
3	L	303	ASN
3	L	321	ASP
3	L	334	SER
3	L	343	ASN
3	L	345	SER
3	L	362	PHE
1	M	3	ASN
1	M	21	GLU
1	M	35	GLN
1	M	42	VAL
1	M	43	SER
1	M	88	SER
2	N	32	GLN
2	N	55	SER
2	N	59	LYS
2	N	63	MET
2	N	81	ASP
2	N	91	ARG
2	N	122	PHE
3	O	204	ILE
3	O	232	LYS
3	O	321	ASP
3	O	344	VAL
3	O	345	SER
3	O	352	GLU
1	P	9	SER
1	P	21	GLU
1	P	42	VAL
1	P	43	SER
1	P	52	PHE
1	P	88	SER
1	P	90	SER
1	P	92	ILE
1	P	119	SER
2	Q	17	SER
2	Q	32	GLN
2	Q	34	LYS
2	Q	59	LYS

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Mol	Chain	Res	Type
2	Q	67	LEU
2	Q	88	VAL
2	Q	91	ARG
2	Q	119	VAL
2	Q	120	CYS
2	Q	122	PHE
3	R	231	THR
3	R	305	LEU
3	R	321	ASP
3	R	344	VAL
3	R	352	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (36) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	60	ASN
2	B	69	ASN
3	C	298	GLN
3	C	318	ASN
1	D	13	GLN
1	D	17	GLN
1	D	59	ASN
1	D	60	ASN
2	E	32	GLN
2	E	109	ASN
3	F	241	GLN
3	F	303	ASN
3	F	343	ASN
1	G	3	ASN
1	G	60	ASN
1	G	79	GLN
3	I	210	GLN
3	I	241	GLN
3	I	303	ASN
3	I	343	ASN
1	J	79	GLN
2	K	9	ASN
2	K	33	HIS
2	K	68	ASN
2	K	109	ASN
3	L	225	ASN
3	L	251	ASN

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Mol	Chain	Res	Type
1	M	55	GLN
1	M	60	ASN
2	N	32	GLN
3	O	200	GLN
3	O	303	ASN
3	O	318	ASN
2	Q	32	GLN
3	R	218	ASN
3	R	318	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 37 ligands modelled in this entry, 21 are monoatomic and 1 is modelled with single atom - leaving 15 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	SO4	Q	201	-	4,4,4	0.30	0	6,6,6	0.08	0
4	SO4	E	203	5,6	4,4,4	0.30	0	6,6,6	0.30	0
4	SO4	G	202	-	4,4,4	0.27	0	6,6,6	0.17	0
4	SO4	J	201	-	4,4,4	0.26	0	6,6,6	0.12	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	SO4	N	201	5,6	4,4,4	0.33	0	6,6,6	0.09	0
4	SO4	R	402	-	4,4,4	0.34	0	6,6,6	0.09	0
4	SO4	H	202	5,6	4,4,4	0.34	0	6,6,6	0.23	0
4	SO4	D	201	-	4,4,4	0.33	0	6,6,6	0.10	0
4	SO4	K	202	6	4,4,4	0.35	0	6,6,6	0.09	0
4	SO4	Q	202	6	4,4,4	0.32	0	6,6,6	0.11	0
4	SO4	A	201	-	4,4,4	0.28	0	6,6,6	0.13	0
4	SO4	O	403	-	4,4,4	0.31	0	6,6,6	0.05	0
4	SO4	E	201	-	4,4,4	0.31	0	6,6,6	0.10	0
4	SO4	B	201	6	4,4,4	0.37	0	6,6,6	0.15	0
4	SO4	A	202	-	4,4,4	0.32	0	6,6,6	0.09	0

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	O	403	SO4	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	A	131/135 (97%)	-1.39	0 100 100	43, 61, 85, 91	0
1	D	131/135 (97%)	-1.41	0 100 100	36, 57, 87, 103	0
1	G	131/135 (97%)	-1.38	0 100 100	35, 59, 92, 102	0
1	J	131/135 (97%)	-1.43	0 100 100	43, 59, 83, 97	0
1	M	131/135 (97%)	-1.27	0 100 100	39, 74, 123, 140	0
1	P	131/135 (97%)	-1.28	0 100 100	42, 71, 125, 134	0
2	B	122/124 (98%)	-1.26	0 100 100	47, 75, 107, 124	0
2	E	122/124 (98%)	-1.35	0 100 100	40, 65, 97, 119	0
2	H	122/124 (98%)	-1.33	0 100 100	40, 65, 95, 117	0
2	K	122/124 (98%)	-1.29	0 100 100	46, 73, 107, 127	0
2	N	122/124 (98%)	-1.31	0 100 100	39, 71, 103, 128	0
2	Q	122/124 (98%)	-1.31	0 100 100	44, 72, 103, 122	0
3	C	191/217 (88%)	-1.40	0 100 100	42, 65, 91, 145	0
3	F	191/217 (88%)	-1.36	0 100 100	33, 62, 95, 114	0
3	I	191/217 (88%)	-1.41	0 100 100	32, 62, 95, 122	0
3	L	191/217 (88%)	-1.37	0 100 100	42, 63, 92, 143	0
3	O	191/217 (88%)	-1.43	0 100 100	39, 60, 91, 115	0
3	R	191/217 (88%)	-1.42	0 100 100	38, 59, 93, 117	0
All	All	2664/2856 (93%)	-1.36	0 100 100	32, 64, 99, 145	0

There are no RSRZ outliers to report.

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 ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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
4	SO4	A	202	5/5	0.95	0.08	153,157,186,208	0
7	CL	G	201	1/1	0.96	0.08	74,74,74,74	0
4	SO4	N	201	5/5	0.98	0.07	99,99,113,126	0
4	SO4	O	403	5/5	0.98	0.04	93,96,117,126	0
4	SO4	Q	202	5/5	0.98	0.07	96,107,124,130	0
4	SO4	R	402	5/5	0.98	0.04	81,88,101,101	5
4	SO4	A	201	5/5	0.98	0.09	79,86,90,96	0
7	CL	K	201	1/1	0.98	0.04	70,70,70,70	0
7	CL	L	403	1/1	0.98	0.09	70,70,70,70	0
7	CL	M	201	1/1	0.98	0.03	59,59,59,59	0
7	CL	R	403	1/1	0.98	0.03	69,69,69,69	0
4	SO4	D	201	5/5	0.99	0.04	50,68,74,89	0
4	SO4	Q	201	5/5	0.99	0.04	77,77,96,97	0
4	SO4	E	201	5/5	0.99	0.04	80,85,95,104	0
4	SO4	E	203	5/5	0.99	0.04	68,72,85,98	0
6	NA	C	402	1/1	0.99	0.03	35,35,35,35	0
6	NA	I	402	1/1	0.99	0.02	24,24,24,24	0
6	NA	L	402	1/1	0.99	0.02	28,28,28,28	0
6	NA	O	402	1/1	0.99	0.03	49,49,49,49	0
6	NA	R	404	1/1	0.99	0.03	50,50,50,50	0
7	CL	C	403	1/1	0.99	0.02	66,66,66,66	0
7	CL	E	202	1/1	0.99	0.07	62,62,62,62	0
4	SO4	G	202	5/5	0.99	0.03	55,64,75,84	0
7	CL	H	201	1/1	0.99	0.03	69,69,69,69	0
7	CL	J	202	1/1	0.99	0.12	80,80,80,80	0
4	SO4	H	202	5/5	0.99	0.04	65,77,100,104	0
4	SO4	J	201	5/5	0.99	0.04	62,69,79,87	0
4	SO4	K	202	5/5	0.99	0.05	69,75,97,99	0
4	SO4	B	201	5/5	0.99	0.03	65,70,103,119	0
8	NH4	D	202	1/1	0.99	0.06	31,31,31,31	0
5	CO	F	401	1/1	1.00	0.02	63,63,63,63	0
5	CO	I	401	1/1	1.00	0.02	60,60,60,60	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	CO	L	401	1/1	1.00	0.02	62,62,62,62	0
5	CO	O	401	1/1	1.00	0.03	89,89,89,89	0
5	CO	R	401	1/1	1.00	0.02	88,88,88,88	0
5	CO	C	401	1/1	1.00	0.01	63,63,63,63	0
6	NA	F	402	1/1	1.00	0.02	21,21,21,21	0

6.5 Other polymers [i](#)

There are no such residues in this entry.