



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 5, 2026 – 03:21 AM UTC

PDB ID : 6NDA / pdb_00006nda
Title : RHODOCETIN IN COMPLEX WITH THE INTEGRIN ALPHA2-A DOMAIN AND CADMIUM
Authors : Stetefeld, J.; McDougall, M.D.; Loewen, P.C.
Deposited on : 2018-12-13
Resolution : 3.15 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

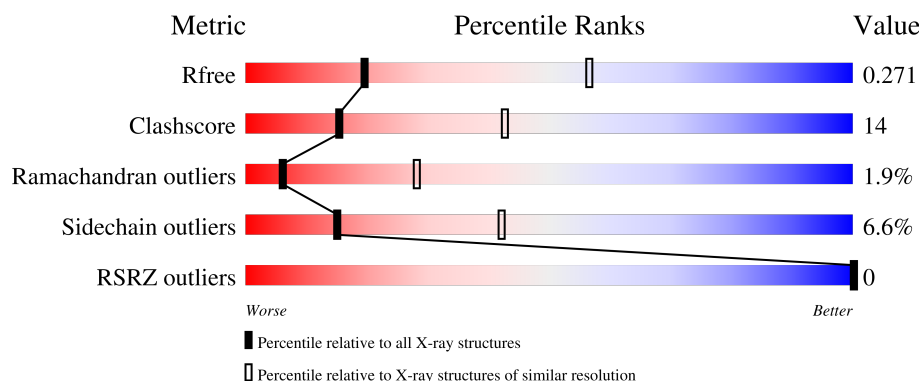
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

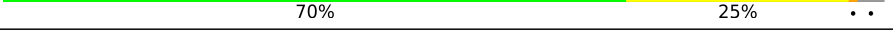
The reported resolution of this entry is 3.15 Å.

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.



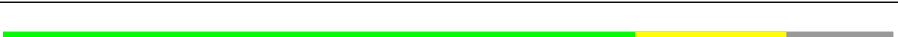
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	2361 (3.20-3.12)
Clashscore	190562	2486 (3.20-3.12)
Ramachandran outliers	187476	2405 (3.20-3.12)
Sidechain outliers	187428	2404 (3.20-3.12)
RSRZ outliers	180081	2361 (3.20-3.12)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	135	 65% 31% . .
1	D	135	 64% 30% . .
1	G	135	 70% 20% 7% .
1	J	135	 70% 25% . .
1	M	135	 59% 33% . .

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Mol	Chain	Length	Quality of chain
1	P	135	 56%36% . . .
2	B	124	 60%35% . .
2	E	124	 64%32% . .
2	H	124	 70%26% . .
2	K	124	 60%35% . .
2	N	124	 52%41% . . .
2	Q	124	 60%32%5% . .
3	C	217	 61%25% . 12%
3	F	217	 63%20%5% 12%
3	I	217	 57%27% . 12%
3	L	217	 71%17% 12%
3	O	217	 62%25% . 12%
3	R	217	 62%24% . 12%

2 Entry composition [i](#)

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 1482	C 940	N 250	O 287	S 5	0	0	0
3	F	191	Total 1482	C 940	N 250	O 287	S 5	0	0	0
3	I	191	Total 1482	C 940	N 250	O 287	S 5	0	0	0
3	L	191	Total 1482	C 940	N 250	O 287	S 5	0	0	0
3	O	191	Total 1482	C 940	N 250	O 287	S 5	0	0	0
3	R	191	Total 1482	C 940	N 250	O 287	S 5	0	0	0

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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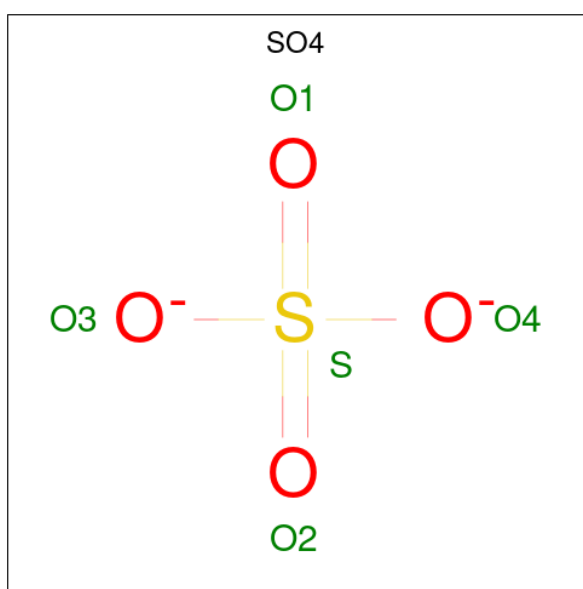
Chain	Residue	Modelled	Actual	Comment	Reference
L	160	SER	-	expression tag	UNP P17301
L	161	SER	-	expression tag	UNP P17301
L	162	GLY	-	expression tag	UNP P17301
L	163	LEU	-	expression tag	UNP P17301
L	164	VAL	-	expression tag	UNP P17301
L	165	PRO	-	expression tag	UNP P17301
L	166	ARG	-	expression tag	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	C	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	F	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 CADMIUM ION (CCD ID: CD) (formula: Cd).

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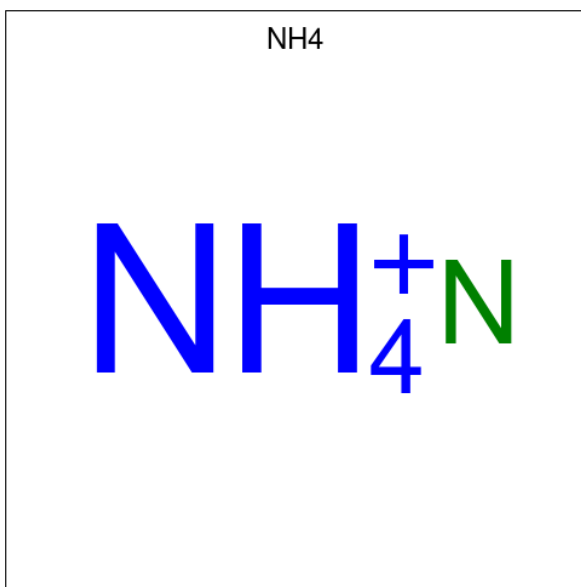
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	L	1	Total 1	Na 1	0	0
6	O	1	Total 1	Na 1	0	0
6	R	1	Total 1	Na 1	0	0

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	C	2	Total 2	Cl 2	0	0
7	E	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

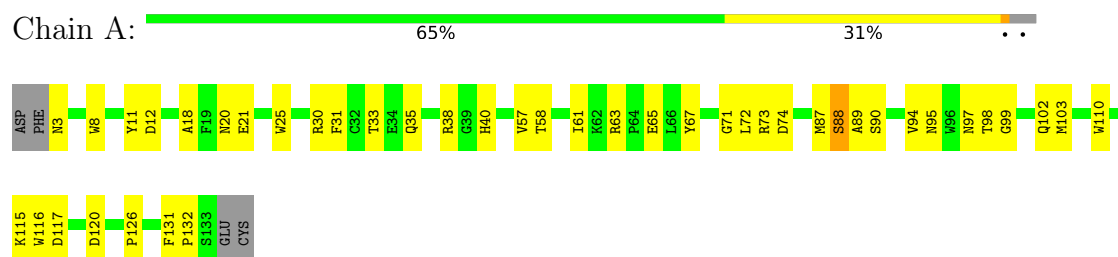
- Molecule 9 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	A	4	Total O 4 4	0	0
9	D	1	Total O 1 1	0	0
9	F	1	Total O 1 1	0	0
9	G	1	Total O 1 1	0	0

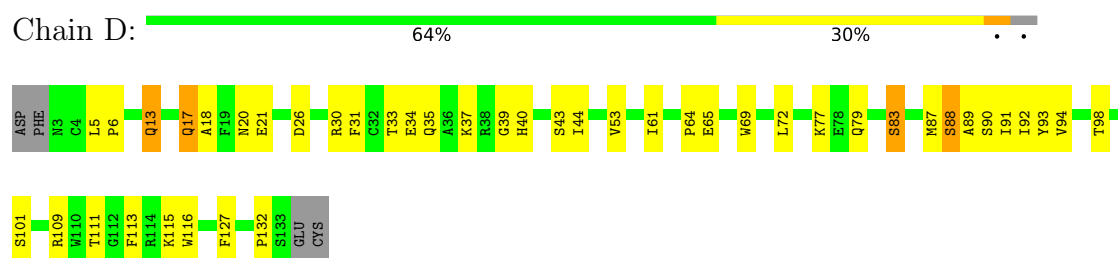
3 Residue-property plots

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.

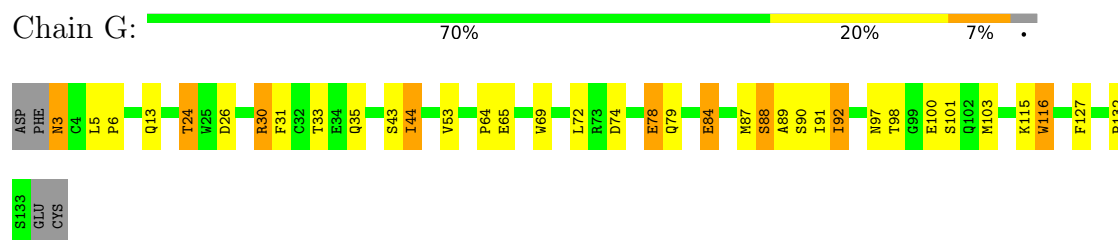
• Molecule 1: Snaclec rhodocetin subunit gamma



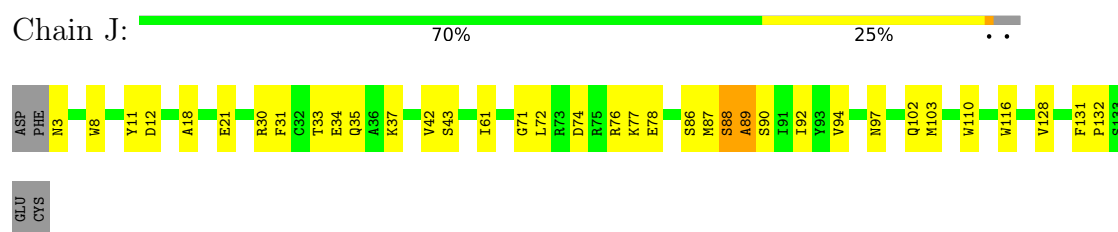
• Molecule 1: Snaclec rhodocetin subunit gamma



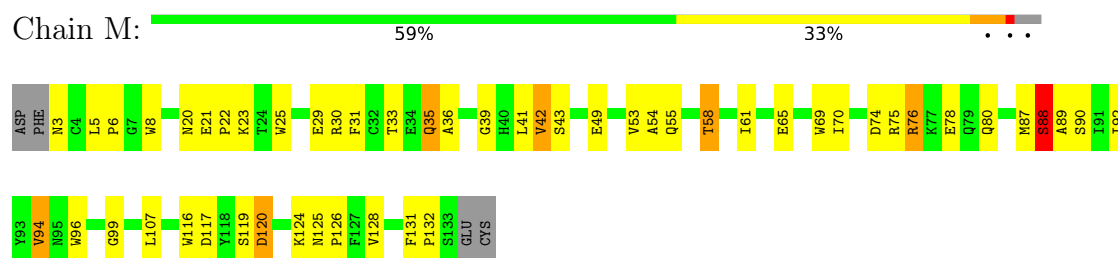
• Molecule 1: Snaclec rhodocetin subunit gamma



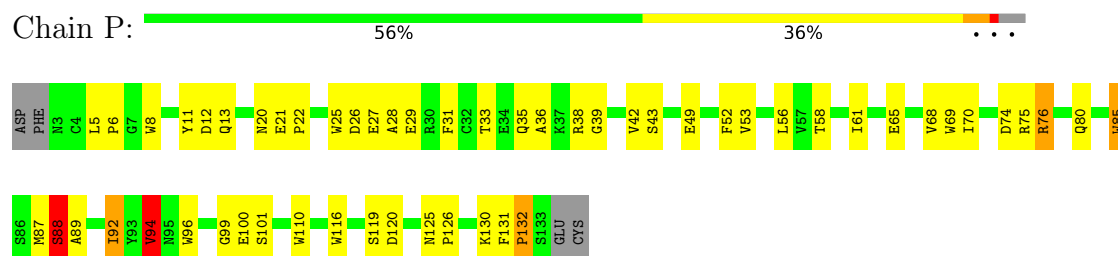
• Molecule 1: Snaclec rhodocetin subunit gamma



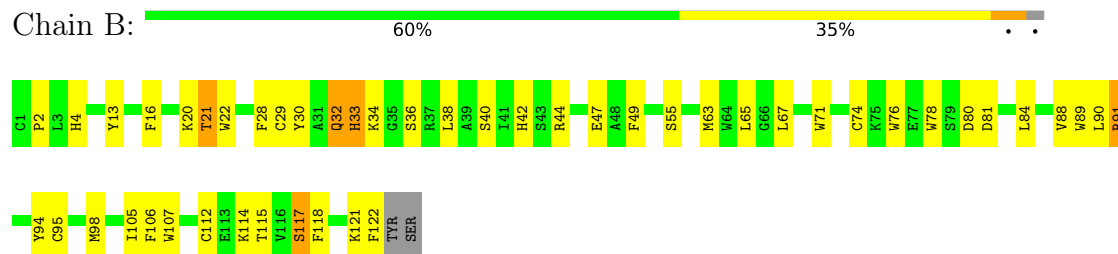
- Molecule 1: Snaclec rhodocetin subunit gamma



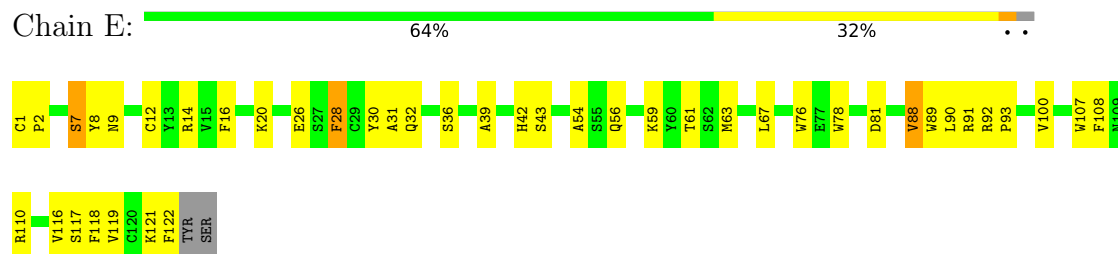
- Molecule 1: Snaclec rhodocetin subunit gamma



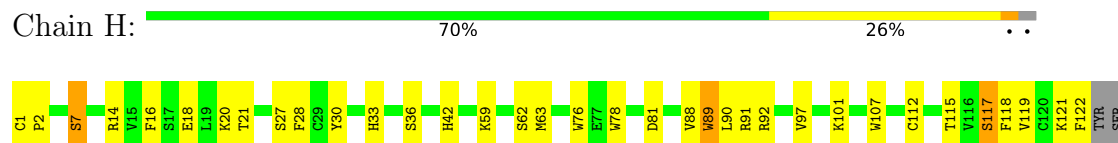
- Molecule 2: Snaclec rhodocetin subunit delta



- Molecule 2: Snaclec rhodocetin subunit delta



- Molecule 2: Snaclec rhodocetin subunit delta



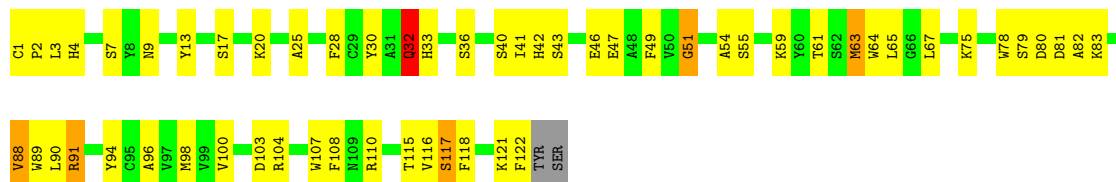
- Molecule 2: Snaclec rhodocetin subunit delta





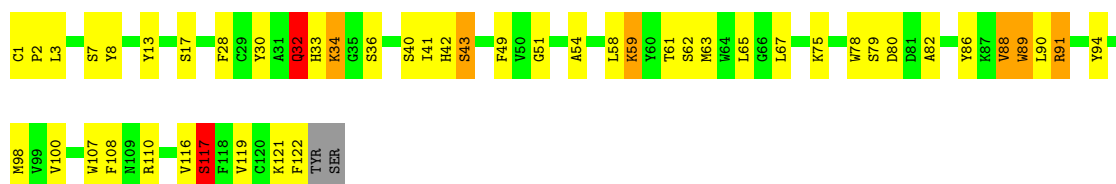
- Molecule 2: Snaclec rhodocetin subunit delta

Chain N: 52% 41% . . .



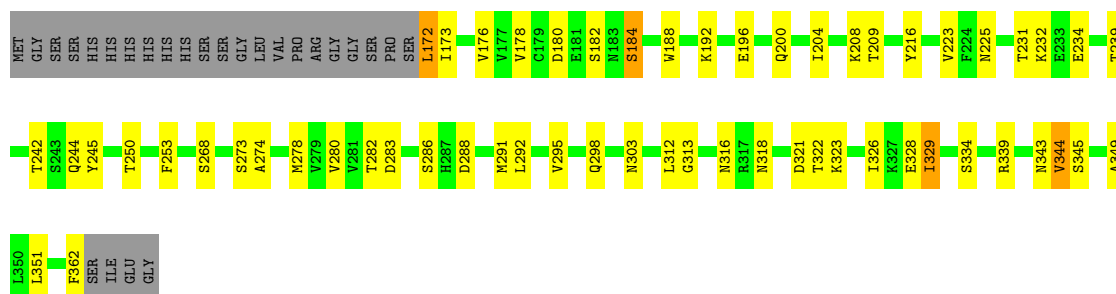
- Molecule 2: Snaclec rhodocetin subunit delta

Chain Q: 60% 32% 5% . .



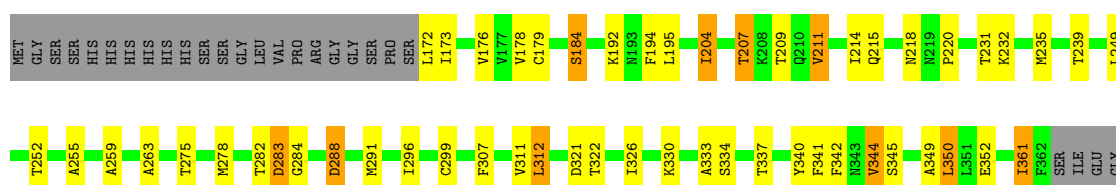
- Molecule 3: Integrin alpha-2

Chain C: 61% 25% . 12%

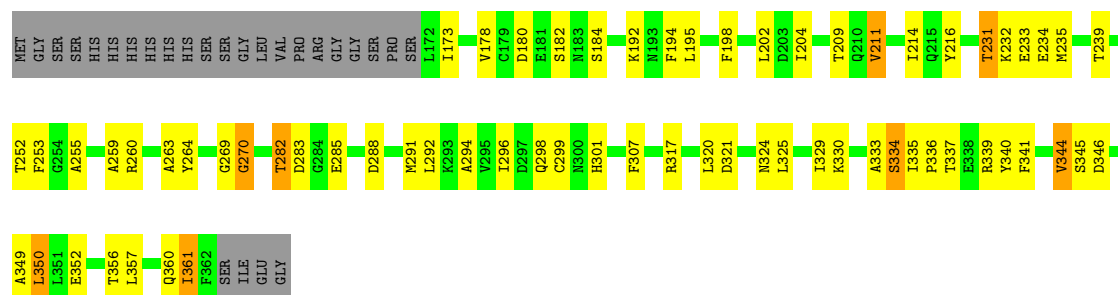


- Molecule 3: Integrin alpha-2

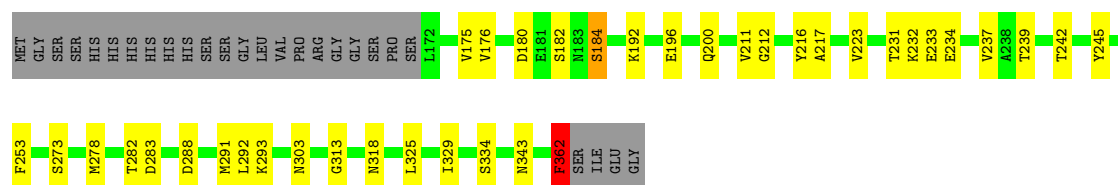
Chain F: 63% 20% 5% 12%



Chain I: 57% 27% 2% 12%



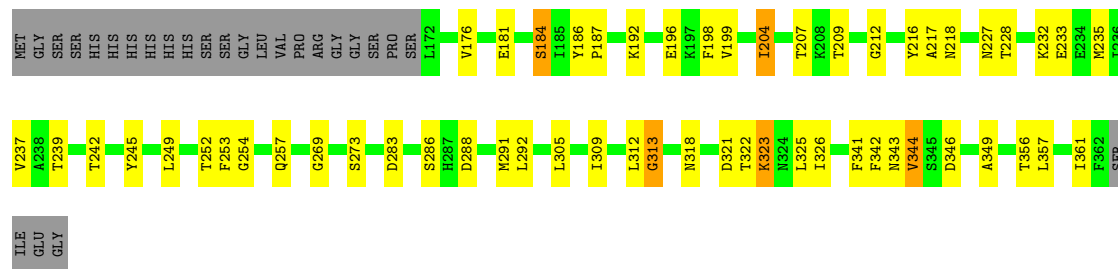
Chain L: 71% 17% 12%



Chain O: 62% 25% 12%



Chain R: 62% 24% • 12%



4 Data and refinement statistics

Property	Value	Source
Space group	P 41	Depositor
Cell constants a, b, c, α , β , γ	130.53Å 130.53Å 250.92Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	92.47 – 3.15 92.47 – 3.15	Depositor EDS
% Data completeness (in resolution range)	93.7 (92.47-3.15) 93.7 (92.47-3.15)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.91 (at 3.13Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.212 , 0.271 0.215 , 0.271	Depositor DCC
R_{free} test set	3406 reflections (4.71%)	wwPDB-VP
Wilson B-factor (Å ²)	50.3	Xtriage
Anisotropy	0.290	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 51.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.458 for h,-k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	21590	wwPDB-VP
Average B, all atoms (Å ²)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.11% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, CD, CL, NH4, NA

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	1.03	0/1101	1.39	0/1491
1	D	1.08	0/1101	1.40	1/1491 (0.1%)
1	G	1.06	1/1101 (0.1%)	1.40	0/1491
1	J	1.05	0/1101	1.41	1/1491 (0.1%)
1	M	1.10	0/1101	1.47	2/1491 (0.1%)
1	P	1.08	0/1101	1.46	3/1491 (0.2%)
2	B	1.02	0/1064	1.39	0/1439
2	E	1.00	0/1064	1.36	0/1439
2	H	0.99	0/1064	1.37	0/1439
2	K	1.03	0/1064	1.35	0/1439
2	N	1.05	0/1064	1.33	0/1439
2	Q	1.07	0/1064	1.34	0/1439
3	C	1.05	0/1506	1.56	1/2040 (0.0%)
3	F	1.07	0/1506	1.56	5/2040 (0.2%)
3	I	1.08	0/1506	1.56	5/2040 (0.2%)
3	L	1.04	0/1506	1.54	2/2040 (0.1%)
3	O	1.04	0/1506	1.51	2/2040 (0.1%)
3	R	1.04	0/1506	1.47	3/2040 (0.1%)
All	All	1.05	1/22026 (0.0%)	1.45	25/29820 (0.1%)

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	P	0	1
2	B	0	1
2	E	0	1

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Mol	Chain	#Chirality outliers	#Planarity outliers
2	H	0	1
2	N	0	1
2	Q	0	2
All	All	0	9

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	G	44	ILE	C-O	5.08	1.29	1.24

The worst 5 of 25 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	P	92	ILE	N-CA-C	-6.77	106.64	113.47
3	F	211	VAL	CA-C-N	6.74	126.23	121.65
3	F	211	VAL	C-N-CA	6.74	126.23	121.65
3	F	220	PRO	CA-C-O	-6.00	114.75	122.13
3	I	211	VAL	CA-C-N	5.88	125.65	121.65

There are no chirality outliers.

5 of 9 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	132	PRO	Peptide
2	B	4	HIS	Peptide
2	E	39	ALA	Peptide
1	G	132	PRO	Peptide
2	H	89	TRP	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	38	0
1	D	1070	0	988	31	0
1	G	1070	0	988	26	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	J	1070	0	988	33	0
1	M	1070	0	988	40	0
1	P	1070	0	988	55	0
2	B	1029	0	977	37	0
2	E	1029	0	977	42	0
2	H	1029	0	977	34	0
2	K	1029	0	977	31	0
2	N	1029	0	977	49	0
2	Q	1029	0	977	46	0
3	C	1482	0	1477	32	0
3	F	1482	0	1477	34	0
3	I	1482	0	1477	38	0
3	L	1482	0	1477	23	0
3	O	1482	0	1477	34	0
3	R	1482	0	1477	37	0
4	A	10	0	0	0	0
4	C	5	0	0	1	0
4	D	5	0	0	0	0
4	E	5	0	0	0	0
4	F	5	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	1	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	1	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	1	0
7	C	2	0	0	0	0
7	E	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
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	4	0	0	0	0
9	D	1	0	0	0	0
9	F	1	0	0	0	0
9	G	1	0	0	0	0
All	All	21590	0	20652	572	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 14.

The worst 5 of 572 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)
2:H:20:LYS:HE2	2:H:28:PHE:CD2	2.01	0.95
2:Q:54:ALA:CB	2:Q:63:MET:HE1	2.02	0.90
2:Q:54:ALA:HB1	2:Q:63:MET:HE1	1.56	0.86
2:E:20:LYS:HE2	2:E:28:PHE:CD2	2.11	0.86
1:M:89:ALA:HB1	2:N:42:HIS:NE2	1.92	0.84

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%)	16 (12%)	1 (1%)	16 46

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	129/135 (96%)	112 (87%)	15 (12%)	2 (2%)	7	32
1	G	129/135 (96%)	115 (89%)	12 (9%)	2 (2%)	7	32
1	J	129/135 (96%)	115 (89%)	12 (9%)	2 (2%)	7	32
1	M	129/135 (96%)	109 (84%)	15 (12%)	5 (4%)	2	14
1	P	129/135 (96%)	111 (86%)	16 (12%)	2 (2%)	7	32
2	B	120/124 (97%)	101 (84%)	15 (12%)	4 (3%)	3	17
2	E	120/124 (97%)	102 (85%)	16 (13%)	2 (2%)	7	30
2	H	120/124 (97%)	108 (90%)	11 (9%)	1 (1%)	16	46
2	K	120/124 (97%)	102 (85%)	13 (11%)	5 (4%)	2	13
2	N	120/124 (97%)	100 (83%)	13 (11%)	7 (6%)	1	8
2	Q	120/124 (97%)	99 (82%)	15 (12%)	6 (5%)	1	10
3	C	189/217 (87%)	173 (92%)	15 (8%)	1 (0%)	24	55
3	F	189/217 (87%)	171 (90%)	16 (8%)	2 (1%)	11	39
3	I	189/217 (87%)	175 (93%)	12 (6%)	2 (1%)	11	39
3	L	189/217 (87%)	173 (92%)	14 (7%)	2 (1%)	11	39
3	O	189/217 (87%)	172 (91%)	16 (8%)	1 (0%)	24	55
3	R	189/217 (87%)	176 (93%)	10 (5%)	3 (2%)	7	32
All	All	2628/2856 (92%)	2326 (88%)	252 (10%)	50 (2%)	6	28

5 of 50 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	33	HIS
2	E	2	PRO
1	G	88	SER
3	I	334	SER
1	M	76	ARG

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	54
1	D	114/118 (97%)	105 (92%)	9 (8%)	11	36
1	G	114/118 (97%)	105 (92%)	9 (8%)	11	36
1	J	114/118 (97%)	110 (96%)	4 (4%)	32	59
1	M	114/118 (97%)	104 (91%)	10 (9%)	9	31
1	P	114/118 (97%)	105 (92%)	9 (8%)	11	36
2	B	110/112 (98%)	102 (93%)	8 (7%)	13	38
2	E	110/112 (98%)	104 (94%)	6 (6%)	19	47
2	H	110/112 (98%)	104 (94%)	6 (6%)	19	47
2	K	110/112 (98%)	100 (91%)	10 (9%)	9	30
2	N	110/112 (98%)	101 (92%)	9 (8%)	10	34
2	Q	110/112 (98%)	100 (91%)	10 (9%)	9	30
3	C	160/181 (88%)	145 (91%)	15 (9%)	8	29
3	F	160/181 (88%)	149 (93%)	11 (7%)	14	40
3	I	160/181 (88%)	147 (92%)	13 (8%)	11	35
3	L	160/181 (88%)	153 (96%)	7 (4%)	25	54
3	O	160/181 (88%)	157 (98%)	3 (2%)	50	69
3	R	160/181 (88%)	152 (95%)	8 (5%)	22	50
All	All	2304/2466 (93%)	2152 (93%)	152 (7%)	15	42

5 of 152 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	N	55	SER
2	Q	122	PHE
2	N	88	VAL
1	P	88	SER
3	R	344	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 34 such sidechains are listed below:

Mol	Chain	Res	Type
2	N	32	GLN
3	O	200	GLN
1	P	14	HIS
2	E	109	ASN

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Mol	Chain	Res	Type
2	E	32	GLN

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	K	202	-	4,4,4	0.33	0	6,6,6	0.21	0
4	SO4	O	403	-	4,4,4	0.31	0	6,6,6	0.13	0
4	SO4	D	201	-	4,4,4	0.33	0	6,6,6	0.11	0
4	SO4	Q	201	-	4,4,4	0.26	0	6,6,6	0.10	0
4	SO4	H	202	5,6	4,4,4	0.40	0	6,6,6	0.31	0
4	SO4	E	201	-	4,4,4	0.27	0	6,6,6	0.11	0
4	SO4	A	202	-	4,4,4	0.32	0	6,6,6	0.11	0
4	SO4	J	201	-	4,4,4	0.28	0	6,6,6	0.10	0
4	SO4	A	201	-	4,4,4	0.28	0	6,6,6	0.10	0
4	SO4	G	201	-	4,4,4	0.33	0	6,6,6	0.12	0
4	SO4	N	201	5,6	4,4,4	0.22	0	6,6,6	0.21	0
4	SO4	F	402	5,6	4,4,4	0.35	0	6,6,6	0.39	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	SO4	R	402	-	4,4,4	0.30	0	6,6,6	0.15	0
4	SO4	Q	202	-	4,4,4	0.35	0	6,6,6	0.10	0
4	SO4	C	402	-	4,4,4	0.34	0	6,6,6	0.27	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.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	O	403	SO4	1	0
4	Q	202	SO4	1	0
4	C	402	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 ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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%)	-1.52	0 100 100	29, 49, 72, 79	0
1	D	131/135 (97%)	-1.57	0 100 100	26, 44, 74, 86	0
1	G	131/135 (97%)	-1.53	0 100 100	26, 45, 79, 88	0
1	J	131/135 (97%)	-1.51	0 100 100	30, 49, 71, 80	0
1	M	131/135 (97%)	-1.33	0 100 100	31, 63, 108, 139	0
1	P	131/135 (97%)	-1.29	0 100 100	33, 63, 111, 131	0
2	B	122/124 (98%)	-1.33	0 100 100	34, 62, 89, 109	0
2	E	122/124 (98%)	-1.46	0 100 100	28, 56, 84, 105	0
2	H	122/124 (98%)	-1.45	0 100 100	30, 57, 85, 104	0
2	K	122/124 (98%)	-1.42	0 100 100	32, 59, 92, 107	0
2	N	122/124 (98%)	-1.38	0 100 100	30, 57, 90, 107	0
2	Q	122/124 (98%)	-1.42	0 100 100	31, 58, 90, 112	0
3	C	191/217 (88%)	-1.46	0 100 100	29, 51, 78, 123	0
3	F	191/217 (88%)	-1.48	0 100 100	20, 48, 82, 99	0
3	I	191/217 (88%)	-1.48	0 100 100	24, 48, 80, 114	0
3	L	191/217 (88%)	-1.49	0 100 100	27, 52, 77, 116	0
3	O	191/217 (88%)	-1.50	0 100 100	28, 49, 81, 111	0
3	R	191/217 (88%)	-1.52	0 100 100	29, 48, 79, 113	0
All	All	2664/2856 (93%)	-1.46	0 100 100	20, 52, 87, 139	0

There are no RSRZ outliers to report.

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	SO4	A	202	5/5	0.96	0.07	119,134,145,177	0
7	CL	J	202	1/1	0.97	0.12	77,77,77,77	0
4	SO4	R	402	5/5	0.98	0.05	68,81,90,91	5
7	CL	L	403	1/1	0.98	0.05	61,61,61,61	0
4	SO4	E	201	5/5	0.99	0.06	56,65,74,74	0
4	SO4	F	402	5/5	0.99	0.04	48,52,83,84	0
4	SO4	G	201	5/5	0.99	0.04	50,58,64,70	0
4	SO4	H	202	5/5	0.99	0.03	50,57,82,85	0
4	SO4	J	201	5/5	0.99	0.04	63,66,84,84	0
4	SO4	K	202	5/5	0.99	0.03	60,64,90,91	0
4	SO4	N	201	5/5	0.99	0.04	63,76,86,99	0
4	SO4	O	403	5/5	0.99	0.03	81,82,99,106	0
4	SO4	Q	201	5/5	0.99	0.04	64,65,82,83	0
4	SO4	Q	202	5/5	0.99	0.05	69,71,91,98	0
4	SO4	A	201	5/5	0.99	0.04	65,72,80,84	0
7	CL	C	404	1/1	0.99	0.05	53,53,53,53	0
7	CL	C	405	1/1	0.99	0.02	60,60,60,60	0
7	CL	E	202	1/1	0.99	0.09	56,56,56,56	0
4	SO4	C	402	5/5	0.99	0.03	47,57,83,96	0
7	CL	K	201	1/1	0.99	0.04	49,49,49,49	0
4	SO4	D	201	5/5	0.99	0.04	43,61,68,79	0
7	CL	M	201	1/1	0.99	0.03	51,51,51,51	0
7	CL	R	403	1/1	0.99	0.02	55,55,55,55	0
8	NH4	D	202	1/1	0.99	0.06	13,13,13,13	0
6	NA	L	402	1/1	1.00	0.02	14,14,14,14	0
6	NA	O	402	1/1	1.00	0.02	19,19,19,19	0
6	NA	R	404	1/1	1.00	0.03	11,11,11,11	0
5	CD	C	401	1/1	1.00	0.00	40,40,40,40	0
5	CD	F	401	1/1	1.00	0.01	35,35,35,35	0
5	CD	I	401	1/1	1.00	0.00	33,33,33,33	0
7	CL	H	201	1/1	1.00	0.03	55,55,55,55	0
5	CD	L	401	1/1	1.00	0.01	42,42,42,42	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	CD	O	401	1/1	1.00	0.00	45,45,45,45	0
5	CD	R	401	1/1	1.00	0.01	45,45,45,45	0
6	NA	C	403	1/1	1.00	0.01	8,8,8,8	0
6	NA	F	403	1/1	1.00	0.03	19,19,19,19	0
6	NA	I	402	1/1	1.00	0.03	13,13,13,13	0

6.5 Other polymers [i](#)

There are no such residues in this entry.