



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 07:58 PM UTC

PDB ID : 6NDF / pdb_00006ndf
Title : RHODOCETIN IN COMPLEX WITH THE INTEGRIN ALPHA2-A DOMAIN WITH STRONTIUM
Authors : Stetefeld, J.; McDougall, M.D.; Loewen, P.C.
Deposited on : 2018-12-13
Resolution : 3.05 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

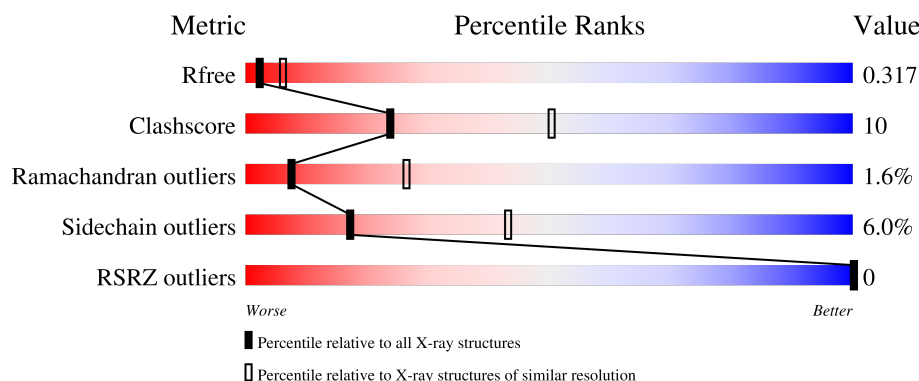
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

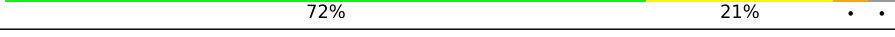
The reported resolution of this entry is 3.05 Å.

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.



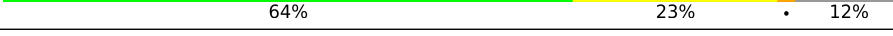
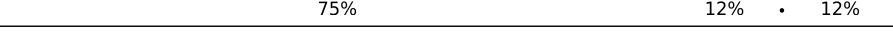
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	2469 (3.10-3.02)
Clashscore	190562	2569 (3.10-3.02)
Ramachandran outliers	187476	2424 (3.10-3.02)
Sidechain outliers	187428	2423 (3.10-3.02)
RSRZ outliers	180081	2469 (3.10-3.02)

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	 73% 21% 5% 1% 1%
1	D	135	 72% 24% 2% 1% 1%
1	G	135	 69% 23% 5% 3% 1%
1	J	135	 72% 21% 5% 1% 1%
1	M	135	 65% 28% 5% 2% 1%

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Mol	Chain	Length	Quality of chain
1	P	135	
2	B	124	
2	E	124	
2	H	124	
2	K	124	
2	N	124	
2	Q	124	
3	C	217	
3	F	217	
3	I	217	
3	L	217	
3	O	217	
3	R	217	

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
4	SO4	O	403	-	-	X	-

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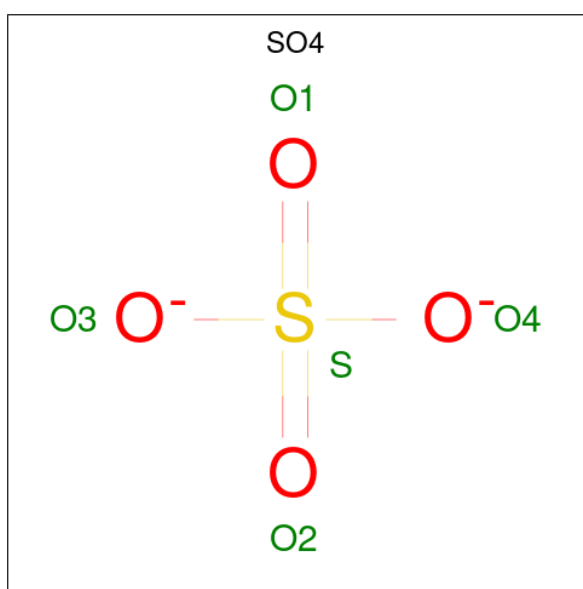
Chain	Residue	Modelled	Actual	Comment	Reference
L	160	SER	-	expression tag	UNP P17301
L	161	SER	-	expression tag	UNP P17301
L	162	GLY	-	expression tag	UNP P17301
L	163	LEU	-	expression tag	UNP P17301
L	164	VAL	-	expression tag	UNP P17301
L	165	PRO	-	expression tag	UNP P17301
L	166	ARG	-	expression tag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

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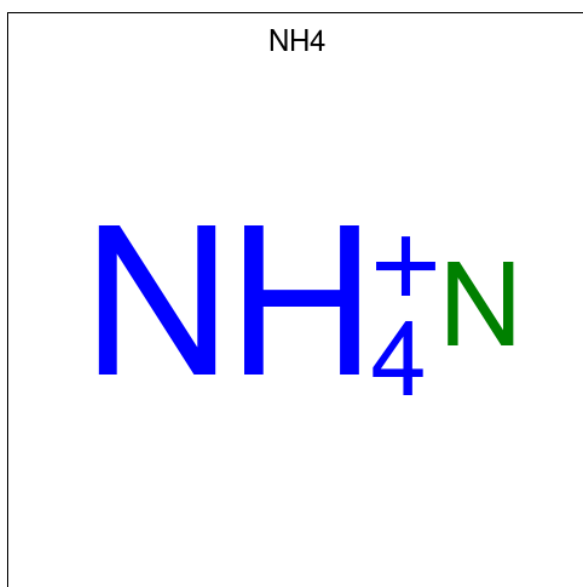
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	L	1	Total 1	Na 1	0	0
6	O	1	Total 1	Na 1	0	0
6	R	1	Total 1	Na 1	0	0

- 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

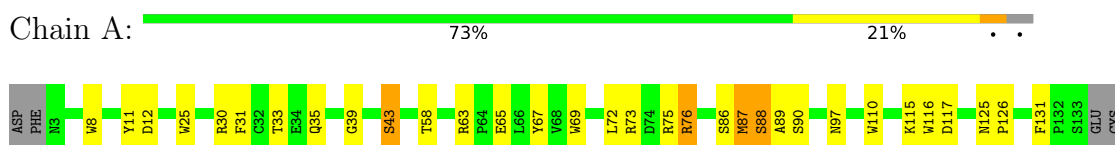
- Molecule 9 is water.

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

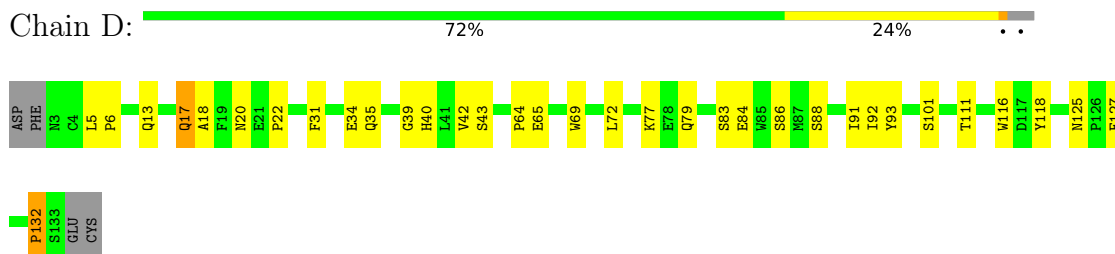
3 Residue-property plots

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.

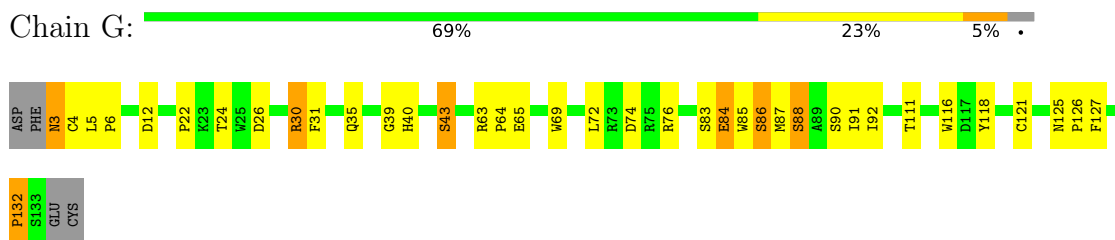
- Molecule 1: Snaclec rhodocetin subunit gamma



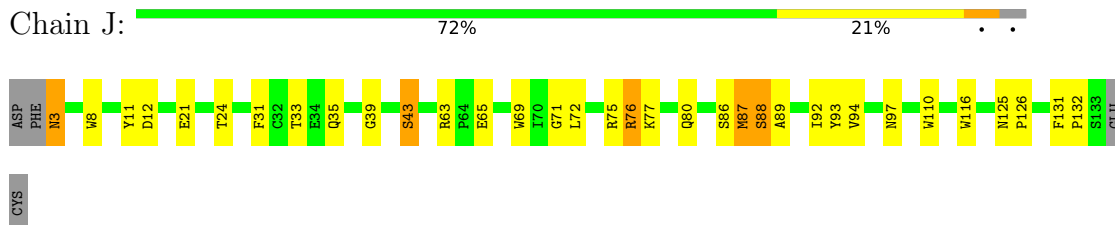
- Molecule 1: Snaclec rhodocetin subunit gamma



- Molecule 1: Snaclec rhodocetin subunit gamma

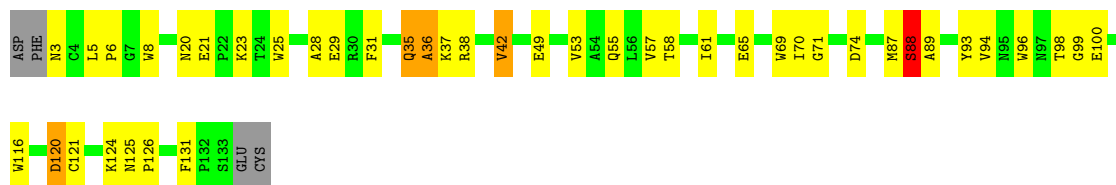


- Molecule 1: Snaclec rhodocetin subunit gamma



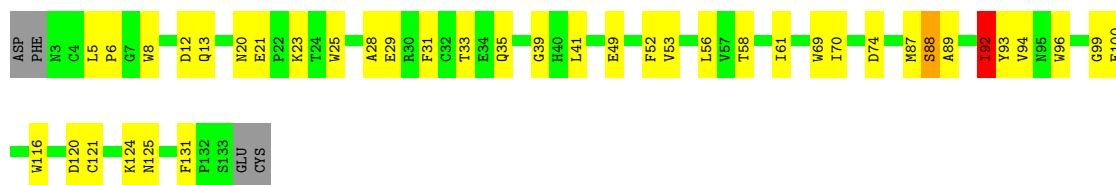
- Molecule 1: Snaclec rhodocetin subunit gamma

Chain M:  65% 28% ...



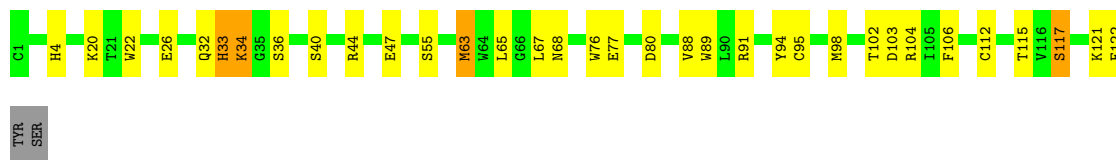
- Molecule 1: Snaclec rhodocetin subunit gamma

Chain P:  67% 28% ..



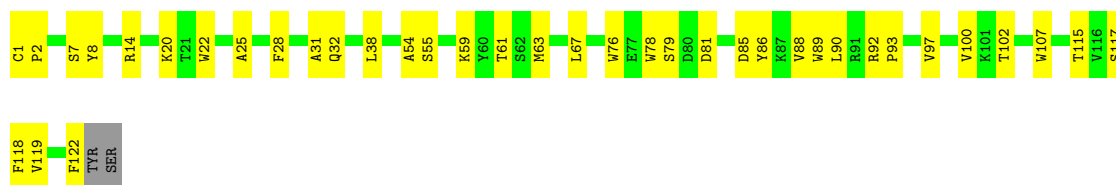
- Molecule 2: Snaclec rhodocetin subunit delta

Chain B:  71% 24% ..



- Molecule 2: Snaclec rhodocetin subunit delta

Chain E:  68% 31% .



- Molecule 2: Snaclec rhodocetin subunit delta

Chain H:  77% 19% ...



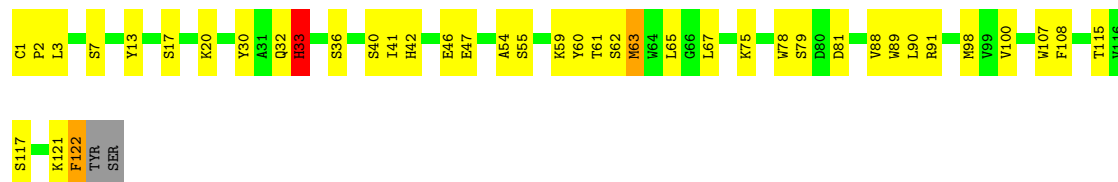
- Molecule 2: Snaclec rhodocetin subunit delta

Chain K:  69% 29% ..



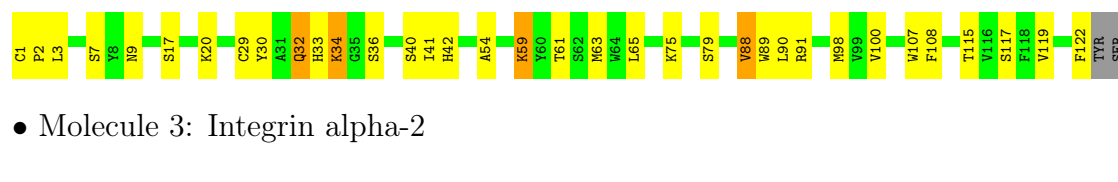
- Molecule 2: Snaclec rhodocetin subunit delta

Chain N:



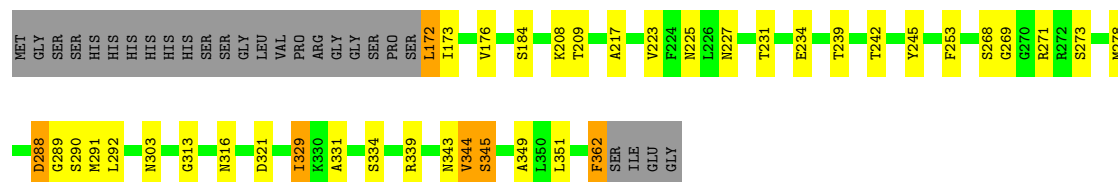
- Molecule 2: Snaclec rhodocetin subunit delta

Chain Q:



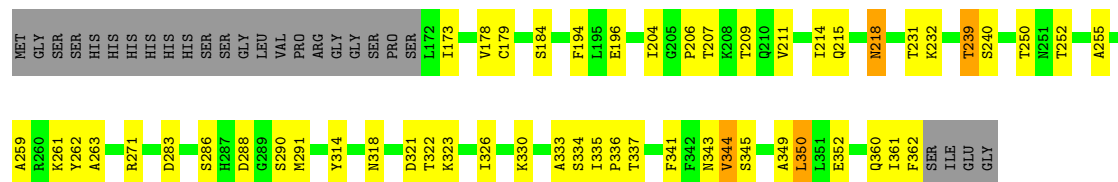
- Molecule 3: Integrin alpha-2

Chain C:



- Molecule 3: Integrin alpha-2

Chain F:



- Molecule 3: Integrin alpha-2

Chain I:





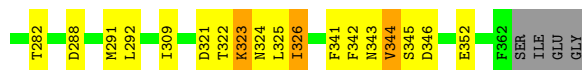
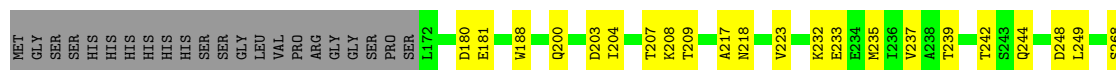
- Molecule 3: Integrin alpha-2



- 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	132.40Å 132.40Å 252.99Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.50 – 3.05 48.50 – 3.05	Depositor EDS
% Data completeness (in resolution range)	98.4 (48.50-3.05) 98.4 (48.50-3.05)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.45 (at 3.07Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.257 , 0.317 0.259 , 0.317	Depositor DCC
R_{free} test set	3998 reflections (4.84%)	wwPDB-VP
Wilson B-factor (Å ²)	67.1	Xtriage
Anisotropy	0.047	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 76.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.36$	Xtriage
Estimated twinning fraction	0.469 for h,-k,-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	21590	wwPDB-VP
Average B, all atoms (Å ²)	78.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.22% 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: SR, NH4, CL, SO4, 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	1.04	0/1101	1.37	0/1491
1	D	1.10	0/1101	1.46	1/1491 (0.1%)
1	G	1.10	0/1101	1.46	0/1491
1	J	1.05	0/1101	1.41	1/1491 (0.1%)
1	M	1.11	0/1101	1.43	0/1491
1	P	1.11	0/1101	1.44	1/1491 (0.1%)
2	B	1.04	0/1064	1.39	0/1439
2	E	1.01	0/1064	1.37	1/1439 (0.1%)
2	H	1.00	0/1064	1.39	1/1439 (0.1%)
2	K	1.05	0/1064	1.37	0/1439
2	N	1.06	0/1064	1.39	0/1439
2	Q	1.06	0/1064	1.37	0/1439
3	C	1.04	0/1506	1.53	5/2040 (0.2%)
3	F	1.06	0/1506	1.47	3/2040 (0.1%)
3	I	1.06	0/1506	1.48	3/2040 (0.1%)
3	L	1.04	0/1506	1.51	2/2040 (0.1%)
3	O	1.04	0/1506	1.52	2/2040 (0.1%)
3	R	1.03	0/1506	1.51	2/2040 (0.1%)
All	All	1.05	0/22026	1.45	22/29820 (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	D	0	1
1	G	0	1
2	B	0	1
2	H	0	1
All	All	0	4

There are no bond length outliers.

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	O	223	VAL	N-CA-C	-7.06	105.49	111.56
3	L	362	PHE	CA-CB-CG	6.32	120.12	113.80
3	R	223	VAL	N-CA-C	-6.31	105.67	111.67
3	I	283	ASP	CA-CB-CG	6.18	118.78	112.60
3	F	283	ASP	CA-CB-CG	5.93	118.53	112.60
1	P	92	ILE	N-CA-C	-5.67	107.74	113.47
2	E	85	ASP	N-CA-C	-5.63	105.14	111.28
1	J	92	ILE	N-CA-C	-5.53	107.42	111.90
3	C	223	VAL	N-CA-C	-5.47	107.11	111.81
3	C	362	PHE	CA-CB-CG	5.43	119.23	113.80
3	C	288	ASP	CA-CB-CG	5.37	117.97	112.60
2	H	85	ASP	N-CA-C	-5.33	105.47	111.28
3	R	248	ASP	CA-CB-CG	5.28	117.88	112.60
3	I	211	VAL	CA-C-N	5.27	125.82	122.18
3	I	211	VAL	C-N-CA	5.27	125.82	122.18
1	D	92	ILE	N-CA-C	-5.22	108.42	113.53
3	L	223	VAL	N-CA-C	-5.14	107.39	111.81
3	O	248	ASP	CA-CB-CG	5.10	117.70	112.60
3	F	211	VAL	CA-C-N	5.08	125.69	122.18
3	F	211	VAL	C-N-CA	5.08	125.69	122.18
3	C	331	ALA	CA-C-N	5.05	127.03	120.56
3	C	331	ALA	C-N-CA	5.05	127.03	120.56

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	4	HIS	Peptide
1	D	132	PRO	Peptide
1	G	132	PRO	Peptide
2	H	85	ASP	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	23	0
1	G	1070	0	988	27	0
1	J	1070	0	988	33	0
1	M	1070	0	988	41	0
1	P	1070	0	988	40	0
2	B	1029	0	977	26	0
2	E	1029	0	977	28	0
2	H	1029	0	977	24	0
2	K	1029	0	977	23	0
2	N	1029	0	977	33	0
2	Q	1029	0	977	29	0
3	C	1482	0	1478	19	0
3	F	1482	0	1477	26	0
3	I	1482	0	1477	22	0
3	L	1482	0	1478	21	0
3	O	1482	0	1477	31	0
3	R	1482	0	1477	22	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	2	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	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	E	1	0	0	0	0
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	F	1	0	0	0	0
9	L	1	0	0	0	0
All	All	21590	0	20654	426	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 (426) 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:29:GLU:OE2	2:N:79:SER:OG	1.81	0.98
2:H:20:LYS:HE2	2:H:28:PHE:CD2	2.06	0.91
1:M:29:GLU:CD	2:N:79:SER:HG	1.82	0.87
1:A:87:MET:SD	2:B:40:SER:HB3	2.15	0.86
1:P:29:GLU:CD	2:Q:79:SER:OG	2.19	0.86
1:G:84:GLU:O	1:G:91:ILE:HD11	1.76	0.84
1:M:29:GLU:CD	2:N:79:SER:OG	2.21	0.84
2:Q:54:ALA:HB1	2:Q:63:MET:HE1	1.59	0.84
2:E:90:LEU:HD12	2:E:90:LEU:C	2.04	0.83
2:Q:54:ALA:CB	2:Q:63:MET:HE1	2.10	0.82
2:H:90:LEU:HD12	2:H:90:LEU:C	2.05	0.81
3:F:361:ILE:O	3:F:362:PHE:O	2.00	0.79
2:E:20:LYS:HE2	2:E:28:PHE:CD2	2.17	0.79
2:Q:90:LEU:HD12	2:Q:90:LEU:C	2.08	0.78
1:P:29:GLU:OE2	2:Q:79:SER:OG	1.99	0.78
2:B:88:VAL:HG12	2:B:88:VAL:O	1.84	0.76
1:D:84:GLU:O	1:D:91:ILE:HD11	1.86	0.76
3:I:344:VAL:HG13	3:I:349:ALA:HB3	1.68	0.75
2:N:1:CYS:SG	2:N:7:SER:HB3	2.27	0.75
2:Q:1:CYS:SG	2:Q:7:SER:HB3	2.28	0.74
1:P:87:MET:O	1:P:88:SER:HB2	1.88	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:116:TRP:O	2:K:89:TRP:HA	1.86	0.74
1:A:87:MET:SD	2:B:40:SER:CB	2.76	0.73
1:J:87:MET:SD	2:K:40:SER:HB3	2.28	0.73
3:O:260:ARG:NH2	4:O:403:SO4:O4	2.21	0.72
2:N:90:LEU:C	2:N:90:LEU:HD12	2.14	0.72
2:N:54:ALA:CB	2:N:63:MET:HE1	2.19	0.72
1:D:31:PHE:O	1:D:35:GLN:HG2	1.89	0.71
1:M:87:MET:O	1:M:88:SER:HB2	1.90	0.71
1:J:87:MET:SD	2:K:40:SER:OG	2.46	0.71
1:J:87:MET:SD	2:K:40:SER:CB	2.79	0.71
3:F:349:ALA:O	3:F:352:GLU:HG2	1.91	0.71
3:I:361:ILE:O	3:I:362:PHE:O	2.09	0.71
1:J:86:SER:OG	2:K:26:GLU:OE2	2.09	0.70
3:C:344:VAL:HG13	3:C:349:ALA:HB3	1.72	0.69
1:G:31:PHE:O	1:G:35:GLN:HG2	1.93	0.69
1:J:87:MET:O	1:J:88:SER:HB3	1.92	0.69
1:D:39:GLY:O	1:D:40:HIS:CD2	2.46	0.68
3:I:349:ALA:O	3:I:352:GLU:HG2	1.92	0.68
1:M:55:GLN:OE1	3:R:207:THR:HG22	1.92	0.68
2:E:54:ALA:HB1	2:E:63:MET:HE1	1.74	0.68
1:P:29:GLU:CD	2:Q:79:SER:HG	1.99	0.68
2:E:31:ALA:HB1	3:O:353:LYS:HE2	1.76	0.67
3:F:322:THR:O	3:F:326:ILE:HG12	1.95	0.67
3:L:351:LEU:HD23	3:L:351:LEU:O	1.96	0.66
3:L:231:THR:OG1	3:L:234:GLU:HB2	1.96	0.66
3:C:173:ILE:HB	3:C:209:THR:HG22	1.78	0.66
1:M:20:ASN:HD21	1:M:61:ILE:HA	1.61	0.65
1:P:5:LEU:HB3	1:P:6:PRO:HD2	1.78	0.65
1:A:115:LYS:HA	2:B:88:VAL:HG12	1.79	0.65
3:L:173:ILE:HB	3:L:209:THR:HG22	1.78	0.64
1:A:116:TRP:O	2:B:89:TRP:HA	1.97	0.64
3:O:239:THR:O	3:O:242:THR:HG22	1.98	0.64
2:B:88:VAL:O	2:B:88:VAL:CG1	2.45	0.64
1:P:89:ALA:HB1	2:Q:42:HIS:CE1	2.32	0.64
3:R:322:THR:O	3:R:323:LYS:C	2.41	0.64
3:C:231:THR:OG1	3:C:234:GLU:HB2	1.98	0.63
1:A:86:SER:OG	2:B:26:GLU:OE2	2.16	0.63
1:G:39:GLY:O	1:G:40:HIS:CD2	2.52	0.62
3:L:253:PHE:CZ	3:L:292:LEU:HG	2.34	0.62
3:C:351:LEU:O	3:C:351:LEU:HD23	1.99	0.62
1:M:89:ALA:HB1	2:N:42:HIS:NE2	2.15	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:72:LEU:HD11	2:H:76:TRP:HB3	1.80	0.62
3:R:204:ILE:HA	3:R:209:THR:O	2.01	0.61
2:E:90:LEU:C	2:E:90:LEU:CD1	2.73	0.61
1:J:8:TRP:CZ3	1:J:131:PHE:CE2	2.88	0.61
1:A:43:SER:HB3	2:B:80:ASP:OD2	2.00	0.61
1:D:79:GLN:O	1:D:101:SER:OG	2.15	0.60
3:R:181:GLU:O	3:R:181:GLU:HG3	2.00	0.60
1:P:87:MET:SD	2:Q:40:SER:OG	2.47	0.60
2:B:44:ARG:O	2:B:47:GLU:HB3	2.01	0.60
1:A:8:TRP:CZ3	1:A:131:PHE:CE2	2.88	0.60
2:E:88:VAL:HG12	2:E:88:VAL:O	2.01	0.60
1:A:87:MET:SD	2:B:40:SER:OG	2.60	0.60
3:F:344:VAL:HG13	3:F:349:ALA:HB3	1.83	0.60
1:J:31:PHE:O	1:J:35:GLN:HG2	2.01	0.60
3:C:253:PHE:CZ	3:C:292:LEU:HG	2.37	0.60
1:A:31:PHE:O	1:A:35:GLN:HG2	2.01	0.59
1:P:49:GLU:O	1:P:53:VAL:HG23	2.03	0.59
3:C:173:ILE:HD12	3:C:209:THR:HG22	1.84	0.59
1:G:87:MET:O	1:G:88:SER:HB3	2.02	0.59
1:D:72:LEU:HD11	2:E:76:TRP:HB3	1.85	0.59
2:N:40:SER:O	2:N:41:ILE:HD13	2.03	0.58
2:Q:40:SER:O	2:Q:41:ILE:HD13	2.02	0.58
1:M:87:MET:SD	2:N:40:SER:OG	2.50	0.58
1:M:49:GLU:O	1:M:53:VAL:HG23	2.03	0.58
2:H:88:VAL:HG12	2:H:88:VAL:O	2.02	0.58
2:H:90:LEU:C	2:H:90:LEU:CD1	2.74	0.58
1:P:52:PHE:CD1	1:P:52:PHE:C	2.82	0.57
2:E:14:ARG:NH1	3:O:359:GLU:OE2	2.38	0.57
1:M:116:TRP:O	2:N:89:TRP:HA	2.04	0.57
2:Q:88:VAL:CG1	2:Q:88:VAL:O	2.52	0.57
1:A:115:LYS:HA	2:B:88:VAL:CG1	2.35	0.57
3:O:322:THR:O	3:O:323:LYS:C	2.48	0.57
1:J:43:SER:HB3	2:K:80:ASP:OD2	2.04	0.57
1:P:12:ASP:OD1	1:P:13:GLN:N	2.38	0.57
3:F:288:ASP:HB2	3:F:291:MET:HE3	1.85	0.56
3:O:322:THR:O	3:O:324:ASN:N	2.38	0.56
1:A:30:ARG:O	1:A:33:THR:OG1	2.18	0.56
2:N:13:TYR:OH	2:N:46:GLU:OE2	2.17	0.56
2:K:20:LYS:O	2:K:115:THR:HA	2.05	0.56
3:R:233:GLU:O	3:R:237:VAL:HG23	2.05	0.56
2:N:88:VAL:CG1	2:N:88:VAL:O	2.53	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:Q:90:LEU:C	2:Q:90:LEU:CD1	2.79	0.56
1:G:24:THR:HG22	1:G:26:ASP:H	1.71	0.56
2:K:44:ARG:O	2:K:47:GLU:HB3	2.06	0.56
3:O:231:THR:OG1	3:O:234:GLU:HB2	2.05	0.56
2:B:103:ASP:O	2:B:104:ARG:HB3	2.06	0.56
3:L:288:ASP:HB2	3:L:291:MET:HE3	1.88	0.56
1:P:31:PHE:O	1:P:35:GLN:HG2	2.06	0.56
1:M:5:LEU:HB3	1:M:6:PRO:HD2	1.89	0.55
1:M:99:GLY:HA2	1:M:120:ASP:HB2	1.89	0.55
3:C:288:ASP:HB2	3:C:291:MET:HE3	1.89	0.55
2:K:8:TYR:HB2	2:K:49:PHE:CE2	2.41	0.55
3:L:362:PHE:CD1	3:L:362:PHE:N	2.75	0.55
2:E:61:THR:CG2	2:E:100:VAL:HB	2.36	0.55
3:O:188:TRP:CD2	3:O:244:GLN:HB2	2.42	0.55
3:O:200:GLN:HE22	3:O:236:ILE:HD13	1.72	0.55
3:F:330:LYS:O	3:F:333:ALA:HB3	2.07	0.55
1:G:64:PRO:HD2	1:G:65:GLU:OE1	2.07	0.55
3:C:225:ASN:OD1	3:C:268:SER:HB2	2.08	0.54
3:F:361:ILE:C	3:F:362:PHE:O	2.50	0.54
3:C:173:ILE:HD12	3:C:209:THR:CG2	2.38	0.54
1:D:93:TYR:CE2	1:J:77:LYS:HB2	2.43	0.54
2:H:28:PHE:CD1	2:H:28:PHE:C	2.85	0.54
3:O:233:GLU:O	3:O:237:VAL:HG23	2.08	0.54
3:O:242:THR:O	3:O:242:THR:HG23	2.07	0.54
3:R:341:PHE:CZ	3:R:343:ASN:HB3	2.42	0.54
3:I:239:THR:O	3:I:242:THR:HG22	2.09	0.54
1:D:43:SER:HA	2:E:78:TRP:CZ3	2.43	0.53
1:A:25:TRP:CZ2	1:A:73:ARG:HD3	2.43	0.53
3:I:288:ASP:HB2	3:I:291:MET:HE3	1.90	0.53
2:Q:34:LYS:N	2:Q:34:LYS:HD2	2.24	0.53
3:R:239:THR:O	3:R:242:THR:HG22	2.09	0.53
2:K:65:LEU:HD21	2:K:98:MET:HE3	1.91	0.53
2:H:20:LYS:HE2	2:H:28:PHE:HD2	1.65	0.53
2:H:54:ALA:HB2	2:H:63:MET:HE1	1.91	0.53
1:J:131:PHE:HD1	1:J:132:PRO:O	1.91	0.53
1:A:97:ASN:HB2	2:B:106:PHE:HB2	1.91	0.53
1:G:3:ASN:N	1:G:3:ASN:ND2	2.55	0.53
1:M:28:ALA:O	1:M:31:PHE:N	2.42	0.53
1:P:87:MET:HG3	1:P:89:ALA:H	1.73	0.53
1:G:5:LEU:HB3	1:G:6:PRO:HD2	1.91	0.52
2:H:97:VAL:O	2:H:107:TRP:HA	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:5:LEU:HB3	1:D:6:PRO:HD2	1.91	0.52
1:D:69:TRP:NE1	1:D:125:ASN:HB2	2.24	0.52
1:M:70:ILE:HG22	1:M:71:GLY:H	1.74	0.52
3:R:204:ILE:HG12	3:R:235:MET:HE1	1.92	0.52
3:R:218:ASN:ND2	3:R:249:LEU:HD12	2.24	0.52
1:G:116:TRP:O	2:H:89:TRP:HA	2.09	0.52
3:O:218:ASN:ND2	3:O:249:LEU:HD12	2.25	0.52
1:M:93:TYR:OH	2:N:47:GLU:OE2	2.28	0.52
2:B:34:LYS:HE3	2:B:34:LYS:HA	1.92	0.52
3:L:362:PHE:N	3:L:362:PHE:HD1	2.06	0.52
1:G:69:TRP:NE1	1:G:125:ASN:HB2	2.24	0.52
1:M:70:ILE:HG22	2:N:78:TRP:HZ3	1.75	0.51
2:E:28:PHE:CD1	2:E:28:PHE:C	2.87	0.51
1:M:25:TRP:NE1	1:M:70:ILE:O	2.42	0.51
1:M:31:PHE:O	1:M:31:PHE:CD1	2.64	0.51
2:N:88:VAL:O	2:N:88:VAL:HG12	2.11	0.51
3:O:261:LYS:NZ	4:O:403:SO4:O3	2.32	0.51
1:A:75:ARG:C	1:A:76:ARG:O	2.53	0.51
3:L:285:GLU:OE1	3:L:324:ASN:HB2	2.11	0.51
3:L:362:PHE:HD1	3:L:362:PHE:H	1.59	0.51
3:R:204:ILE:HG12	3:R:235:MET:CE	2.41	0.51
3:F:194:PHE:CD1	3:F:350:LEU:HD22	2.46	0.51
1:J:72:LEU:HD11	2:K:76:TRP:HB3	1.93	0.51
1:P:28:ALA:O	1:P:31:PHE:HB3	2.11	0.51
1:P:74:ASP:OD2	2:Q:75:LYS:O	2.29	0.51
3:R:344:VAL:CG1	3:R:346:ASP:O	2.59	0.51
3:F:204:ILE:HB	3:F:232:LYS:CG	2.41	0.50
3:F:218:ASN:HD22	3:F:218:ASN:H	1.59	0.50
2:N:90:LEU:C	2:N:90:LEU:CD1	2.85	0.50
2:K:22:TRP:CZ2	2:K:68:ASN:HB2	2.45	0.50
1:M:89:ALA:HB1	2:N:42:HIS:CE1	2.45	0.50
2:H:107:TRP:N	2:H:107:TRP:CD1	2.76	0.50
1:G:22:PRO:HA	1:G:125:ASN:O	2.12	0.50
3:C:172:LEU:HB3	3:C:173:ILE:HG13	1.93	0.50
3:C:329:ILE:N	3:C:329:ILE:HD13	2.26	0.50
3:L:316:ASN:ND2	3:L:345:SER:O	2.44	0.50
1:D:22:PRO:HA	1:D:125:ASN:O	2.12	0.50
2:N:33:HIS:HD2	2:N:122:PHE:O	1.95	0.50
3:R:344:VAL:HG12	3:R:346:ASP:O	2.11	0.50
1:G:69:TRP:CD1	1:G:125:ASN:HB2	2.47	0.49
1:J:97:ASN:HB2	2:K:106:PHE:HB2	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:Q:90:LEU:HD12	2:Q:90:LEU:O	2.11	0.49
1:D:34:GLU:HG3	1:D:34:GLU:O	2.12	0.49
2:E:86:TYR:CE2	2:E:88:VAL:HG21	2.46	0.49
1:P:25:TRP:HB3	1:P:121:CYS:SG	2.52	0.49
2:E:107:TRP:N	2:E:107:TRP:CD1	2.80	0.49
2:N:65:LEU:HD21	2:N:98:MET:HB2	1.94	0.49
3:L:289:GLY:O	3:L:292:LEU:N	2.37	0.49
3:O:204:ILE:HA	3:O:209:THR:O	2.12	0.49
1:P:74:ASP:OD2	2:Q:75:LYS:N	2.45	0.49
2:Q:59:LYS:HE3	2:Q:59:LYS:H	1.77	0.49
2:H:21:THR:HG23	2:H:112:CYS:O	2.13	0.49
3:I:330:LYS:O	3:I:333:ALA:HB3	2.11	0.49
1:J:80:GLN:CD	1:J:80:GLN:H	2.21	0.49
1:A:11:TYR:HA	3:I:291:MET:HE2	1.95	0.49
3:F:314:TYR:O	3:F:318:ASN:ND2	2.45	0.49
1:P:116:TRP:O	2:Q:89:TRP:HA	2.12	0.49
2:H:88:VAL:O	2:H:88:VAL:CG1	2.61	0.49
3:O:309:ILE:HA	3:O:342:PHE:O	2.12	0.49
3:F:252:THR:O	3:F:255:ALA:HB3	2.12	0.49
1:D:42:VAL:HG22	1:D:43:SER:O	2.13	0.48
3:L:289:GLY:C	3:L:291:MET:H	2.21	0.48
2:E:54:ALA:CB	2:E:63:MET:HE1	2.41	0.48
3:I:179:CYS:O	3:I:215:GLN:HA	2.13	0.48
2:Q:61:THR:CG2	2:Q:100:VAL:HB	2.43	0.48
2:B:63:MET:HA	2:B:117:SER:O	2.13	0.48
1:D:111:THR:HG21	2:E:90:LEU:CD2	2.44	0.48
1:M:87:MET:HG3	1:M:89:ALA:H	1.78	0.48
1:G:63:ARG:HG3	1:G:126:PRO:HG3	1.95	0.48
2:H:36:SER:HA	2:H:121:LYS:O	2.12	0.48
1:D:39:GLY:O	1:D:40:HIS:CG	2.66	0.48
3:R:309:ILE:HA	3:R:342:PHE:O	2.12	0.48
1:J:89:ALA:HB1	2:K:42:HIS:CE1	2.48	0.48
3:L:318:ASN:O	3:L:320:LEU:HG	2.14	0.48
1:M:87:MET:HE2	1:M:89:ALA:HB2	1.95	0.48
3:R:180:ASP:HB2	3:R:282:THR:HA	1.94	0.48
3:R:200:GLN:NE2	3:R:232:LYS:HE3	2.29	0.48
1:J:3:ASN:N	1:J:3:ASN:ND2	2.62	0.48
1:A:11:TYR:CD1	1:A:12:ASP:N	2.82	0.48
2:N:61:THR:CG2	2:N:100:VAL:HB	2.44	0.48
3:I:339:ARG:HG2	3:I:360:GLN:HE21	1.79	0.47
1:M:35:GLN:O	1:M:36:ALA:HB3	2.13	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:Q:65:LEU:HD21	2:Q:98:MET:HB2	1.97	0.47
2:E:22:TRP:O	2:E:25:ALA:HB3	2.14	0.47
1:J:71:GLY:O	2:K:78:TRP:HA	2.15	0.47
1:M:55:GLN:OE1	3:R:207:THR:CG2	2.60	0.47
1:P:69:TRP:HA	1:P:69:TRP:CE3	2.49	0.47
3:R:288:ASP:HB2	3:R:291:MET:HE3	1.96	0.47
1:A:87:MET:O	1:A:88:SER:CB	2.62	0.47
2:B:94:TYR:C	2:B:112:CYS:SG	2.98	0.47
2:E:88:VAL:O	2:E:88:VAL:CG1	2.63	0.47
3:F:204:ILE:HA	3:F:209:THR:O	2.14	0.47
1:M:25:TRP:HB3	1:M:121:CYS:SG	2.54	0.47
1:A:25:TRP:CH2	1:A:73:ARG:HD3	2.50	0.47
2:E:38:LEU:HD23	2:E:119:VAL:O	2.14	0.47
3:F:206:PRO:O	3:F:271:ARG:NH2	2.48	0.47
1:G:85:TRP:HE3	2:H:66:GLY:O	1.98	0.47
1:M:74:ASP:OD2	2:N:75:LYS:N	2.47	0.47
1:P:20:ASN:HD21	1:P:61:ILE:HA	1.79	0.47
1:P:69:TRP:NE1	1:P:125:ASN:HB2	2.30	0.47
2:K:68:ASN:HA	2:K:95:CYS:SG	2.55	0.47
3:O:200:GLN:NE2	3:O:236:ILE:HD13	2.30	0.47
3:R:203:ASP:OD2	3:R:208:LYS:HD2	2.14	0.47
1:G:86:SER:HB3	2:H:66:GLY:CA	2.45	0.47
1:J:33:THR:HA	1:J:39:GLY:O	2.15	0.47
2:B:36:SER:HA	2:B:121:LYS:O	2.14	0.46
1:J:11:TYR:CD1	1:J:12:ASP:N	2.84	0.46
1:P:8:TRP:CZ3	1:P:131:PHE:CE2	3.03	0.46
2:B:20:LYS:O	2:B:115:THR:HA	2.15	0.46
1:D:20:ASN:O	1:D:22:PRO:HD3	2.16	0.46
1:P:52:PHE:HE1	1:P:56:LEU:HD12	1.81	0.46
2:Q:20:LYS:O	2:Q:115:THR:HA	2.16	0.46
1:A:69:TRP:NE1	1:A:125:ASN:HB2	2.31	0.46
3:I:349:ALA:O	3:I:350:LEU:C	2.58	0.46
3:F:204:ILE:HB	3:F:232:LYS:HG2	1.98	0.46
3:R:188:TRP:CD2	3:R:244:GLN:HB2	2.51	0.46
3:I:314:TYR:O	3:I:318:ASN:ND2	2.47	0.46
1:M:28:ALA:O	1:M:31:PHE:HB3	2.16	0.46
2:Q:107:TRP:CD1	2:Q:107:TRP:N	2.83	0.46
2:K:103:ASP:O	2:K:104:ARG:HB3	2.15	0.46
1:M:37:LYS:O	1:M:38:ARG:HB2	2.15	0.46
3:O:173:ILE:O	3:O:209:THR:HA	2.15	0.46
1:D:116:TRP:O	2:E:89:TRP:HA	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:69:TRP:NE1	1:J:125:ASN:HB2	2.31	0.46
1:J:75:ARG:C	1:J:76:ARG:O	2.59	0.46
1:J:80:GLN:CD	1:J:80:GLN:N	2.74	0.46
1:M:69:TRP:NE1	1:M:125:ASN:HB2	2.31	0.46
1:A:11:TYR:HA	3:I:291:MET:CE	2.46	0.45
2:B:68:ASN:HA	2:B:95:CYS:SG	2.57	0.45
1:M:42:VAL:HG13	1:M:70:ILE:HG23	1.98	0.45
1:A:33:THR:HA	1:A:39:GLY:O	2.16	0.45
2:K:63:MET:HA	2:K:117:SER:O	2.16	0.45
1:M:23:LYS:O	1:M:124:LYS:HA	2.17	0.45
1:A:63:ARG:HB3	1:A:65:GLU:OE1	2.17	0.45
1:A:67:TYR:HE2	1:A:117:ASP:HB3	1.81	0.45
3:O:325:LEU:O	3:O:326:ILE:C	2.58	0.45
1:D:83:SER:O	1:D:91:ILE:HG13	2.17	0.45
3:F:337:THR:O	3:F:341:PHE:HB3	2.17	0.45
1:P:29:GLU:OE1	2:Q:79:SER:OG	2.35	0.45
3:I:292:LEU:O	3:I:296:ILE:HG12	2.17	0.45
3:O:344:VAL:CG1	3:O:346:ASP:O	2.65	0.45
1:G:43:SER:HB3	2:H:80:ASP:OD2	2.16	0.45
3:I:337:THR:O	3:I:341:PHE:HB3	2.16	0.45
3:O:200:GLN:NE2	3:O:236:ILE:CD1	2.80	0.45
1:P:28:ALA:O	1:P:31:PHE:N	2.50	0.45
1:A:75:ARG:HD2	2:B:77:GLU:OE2	2.17	0.45
2:H:63:MET:HB2	2:H:118:PHE:HA	1.98	0.45
3:L:343:ASN:O	3:L:343:ASN:ND2	2.49	0.45
1:P:20:ASN:ND2	1:P:61:ILE:HD12	2.31	0.45
2:E:97:VAL:O	2:E:107:TRP:HA	2.17	0.45
2:B:32:GLN:O	2:B:33:HIS:HB2	2.16	0.44
2:Q:88:VAL:O	2:Q:88:VAL:HG12	2.17	0.44
1:J:87:MET:O	1:J:88:SER:CB	2.62	0.44
1:M:8:TRP:CZ3	1:M:131:PHE:CE2	3.05	0.44
1:D:64:PRO:HD2	1:D:65:GLU:OE1	2.16	0.44
3:F:178:VAL:HA	3:F:214:ILE:O	2.16	0.44
3:F:196:GLU:OE1	3:F:240:SER:HB3	2.17	0.44
3:O:326:ILE:O	3:O:327:LYS:C	2.60	0.44
1:P:12:ASP:CG	1:P:13:GLN:H	2.24	0.44
2:E:20:LYS:O	2:E:115:THR:HA	2.16	0.44
3:I:178:VAL:HA	3:I:214:ILE:O	2.17	0.44
2:N:13:TYR:CE1	2:N:121:LYS:HG3	2.53	0.44
3:O:200:GLN:HE22	3:O:236:ILE:CD1	2.30	0.44
3:O:242:THR:O	3:O:242:THR:CG2	2.66	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:P:99:GLY:HA2	1:P:120:ASP:HB2	2.00	0.44
1:A:75:ARG:O	1:A:76:ARG:O	2.35	0.44
1:M:69:TRP:HA	1:M:69:TRP:CE3	2.52	0.44
3:O:210:GLN:OE1	3:O:271:ARG:NE	2.50	0.44
2:H:22:TRP:O	2:H:25:ALA:HB3	2.17	0.44
2:Q:30:TYR:HA	2:Q:36:SER:O	2.18	0.44
3:R:325:LEU:O	3:R:326:ILE:C	2.61	0.44
2:N:20:LYS:O	2:N:115:THR:HA	2.18	0.43
2:B:103:ASP:O	2:B:103:ASP:OD1	2.36	0.43
2:E:1:CYS:SG	2:E:7:SER:HB3	2.59	0.43
1:J:93:TYR:O	1:J:94:VAL:HG13	2.18	0.43
2:N:54:ALA:HB1	2:N:63:MET:HE1	1.98	0.43
1:D:69:TRP:CD1	1:D:125:ASN:HB2	2.53	0.43
3:I:180:ASP:CB	3:I:282:THR:HA	2.48	0.43
3:L:289:GLY:C	3:L:291:MET:N	2.76	0.43
3:O:203:ASP:OD2	3:O:208:LYS:HD2	2.18	0.43
1:P:89:ALA:HB1	2:Q:42:HIS:NE2	2.33	0.43
1:G:83:SER:O	1:G:91:ILE:HG13	2.18	0.43
1:M:74:ASP:OD2	2:N:75:LYS:O	2.36	0.43
3:C:239:THR:O	3:C:242:THR:HG22	2.18	0.43
1:D:43:SER:HA	2:E:78:TRP:CE3	2.54	0.43
1:G:74:ASP:OD1	1:G:76:ARG:HG2	2.19	0.43
3:O:344:VAL:HG12	3:O:346:ASP:O	2.19	0.43
2:B:22:TRP:CH2	2:B:68:ASN:HB2	2.54	0.43
3:I:312:LEU:HB2	3:I:316:ASN:ND2	2.33	0.43
1:J:8:TRP:CZ3	1:J:131:PHE:CD2	3.05	0.43
1:M:94:VAL:HG13	1:M:96:TRP:CD1	2.53	0.43
1:A:110:TRP:HA	3:C:245:TYR:CE2	2.54	0.43
3:L:253:PHE:CE1	3:L:292:LEU:HG	2.54	0.43
3:F:179:CYS:O	3:F:215:GLN:HA	2.19	0.43
1:G:39:GLY:O	1:G:40:HIS:CG	2.72	0.43
1:J:87:MET:HE3	2:K:38:LEU:O	2.18	0.43
1:A:72:LEU:HD11	2:B:76:TRP:HB3	2.01	0.43
3:F:261:LYS:HD3	3:F:262:TYR:CZ	2.54	0.43
1:G:87:MET:SD	2:H:40:SER:OG	2.70	0.43
3:I:361:ILE:C	3:I:362:PHE:O	2.62	0.43
1:M:31:PHE:CD1	1:M:31:PHE:C	2.97	0.43
1:G:111:THR:HG21	2:H:90:LEU:CD2	2.49	0.42
2:K:22:TRP:CH2	2:K:68:ASN:HB2	2.54	0.42
3:L:292:LEU:HD23	3:L:292:LEU:C	2.44	0.42
2:B:65:LEU:HD21	2:B:98:MET:HE3	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:53:VAL:O	1:M:57:VAL:HG23	2.20	0.42
1:G:86:SER:HB3	2:H:66:GLY:HA3	2.00	0.42
1:J:87:MET:HE2	1:J:87:MET:HB3	1.83	0.42
3:L:249:LEU:HD22	3:L:287:HIS:ND1	2.33	0.42
2:E:90:LEU:HD12	2:E:90:LEU:O	2.19	0.42
3:I:196:GLU:OE1	3:I:240:SER:HB3	2.20	0.42
2:N:107:TRP:CD1	2:N:107:TRP:N	2.86	0.42
3:C:316:ASN:ND2	3:C:345:SER:O	2.52	0.42
1:D:17:GLN:HE21	1:D:18:ALA:H	1.68	0.42
1:G:24:THR:HG23	1:G:121:CYS:O	2.19	0.42
2:H:85:ASP:OD1	2:H:85:ASP:N	2.52	0.42
3:I:325:LEU:O	3:I:329:ILE:HG12	2.19	0.42
3:I:292:LEU:C	3:I:292:LEU:HD23	2.44	0.42
3:O:180:ASP:HB2	3:O:282:THR:HA	2.00	0.42
3:R:322:THR:O	3:R:324:ASN:N	2.53	0.42
3:C:173:ILE:O	3:C:271:ARG:NH2	2.53	0.42
2:E:61:THR:HG21	2:E:100:VAL:HB	2.01	0.42
1:J:110:TRP:HA	3:L:245:TYR:CE2	2.55	0.42
2:Q:29:CYS:O	2:Q:32:GLN:HB2	2.20	0.42
3:L:289:GLY:O	3:L:291:MET:N	2.53	0.41
1:M:69:TRP:C	1:M:70:ILE:HG13	2.44	0.41
1:D:17:GLN:HE21	1:D:17:GLN:HA	1.86	0.41
2:H:86:TYR:CE2	2:H:88:VAL:HG21	2.54	0.41
3:O:288:ASP:HB2	3:O:291:MET:HE3	2.02	0.41
2:E:92:ARG:HB2	2:E:93:PRO:CD	2.50	0.41
1:G:30:ARG:HG2	1:G:30:ARG:HH21	1.84	0.41
1:P:12:ASP:CG	1:P:13:GLN:N	2.79	0.41
1:P:25:TRP:NE1	1:P:70:ILE:O	2.44	0.41
2:B:22:TRP:CZ2	2:B:68:ASN:HB2	2.55	0.41
1:J:87:MET:HG3	1:J:89:ALA:CB	2.50	0.41
2:N:3:LEU:O	2:N:3:LEU:HG	2.19	0.41
1:P:92:ILE:O	1:P:92:ILE:CG2	2.68	0.41
2:E:61:THR:HG22	2:E:100:VAL:HB	2.02	0.41
2:E:63:MET:HB2	2:E:118:PHE:HA	2.03	0.41
3:I:357:LEU:O	3:I:361:ILE:HG12	2.20	0.41
1:J:76:ARG:NH1	2:K:75:LYS:HD2	2.35	0.41
1:M:65:GLU:OE2	1:M:126:PRO:HD3	2.20	0.41
1:P:23:LYS:O	1:P:124:LYS:HA	2.21	0.41
1:P:33:THR:HA	1:P:39:GLY:O	2.20	0.41
1:M:31:PHE:O	1:M:35:GLN:HG2	2.21	0.41
2:N:30:TYR:HA	2:N:36:SER:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:P:12:ASP:OD1	1:P:12:ASP:C	2.64	0.41
3:F:259:ALA:O	3:F:263:ALA:HB3	2.20	0.41
1:J:11:TYR:O	1:J:12:ASP:C	2.63	0.41
2:N:67:LEU:HD12	2:N:67:LEU:HA	1.87	0.41
3:C:227:ASN:ND2	3:C:269:GLY:CA	2.83	0.41
1:J:63:ARG:HB3	1:J:65:GLU:OE2	2.20	0.41
2:K:21:THR:HA	2:K:114:LYS:O	2.20	0.41
1:P:31:PHE:CD1	1:P:31:PHE:C	2.99	0.41
1:A:8:TRP:CZ3	1:A:131:PHE:CD2	3.08	0.41
1:A:87:MET:HG3	1:A:89:ALA:HB2	2.02	0.41
3:C:176:VAL:O	3:C:278:MET:HA	2.21	0.41
3:C:289:GLY:C	3:C:291:MET:H	2.28	0.41
3:F:204:ILE:HB	3:F:232:LYS:HG3	2.02	0.41
3:F:218:ASN:H	3:F:218:ASN:ND2	2.17	0.41
1:G:30:ARG:HG2	1:G:30:ARG:NH2	2.36	0.41
2:K:94:TYR:C	2:K:112:CYS:SG	3.04	0.41
1:M:29:GLU:OE1	2:N:79:SER:OG	2.38	0.41
2:N:60:TYR:HD2	3:O:321:ASP:OD1	2.04	0.41
1:P:29:GLU:OE1	2:Q:79:SER:CB	2.69	0.41
1:P:41:LEU:HD23	1:P:41:LEU:HA	1.96	0.41
1:P:87:MET:HE2	1:P:89:ALA:HB2	2.02	0.41
1:D:69:TRP:HB2	1:D:127:PHE:HB3	2.03	0.41
3:F:173:ILE:HD12	3:F:361:ILE:CG2	2.50	0.41
3:F:335:ILE:HG13	3:F:336:PRO:HA	2.02	0.41
3:O:292:LEU:C	3:O:292:LEU:HD23	2.46	0.41
3:R:292:LEU:C	3:R:292:LEU:HD23	2.46	0.41
3:O:181:GLU:O	3:O:181:GLU:HG3	2.22	0.40
1:P:100:GLU:CD	2:Q:108:PHE:HB2	2.46	0.40
3:F:196:GLU:HG3	3:F:239:THR:HG22	2.04	0.40
1:G:69:TRP:HB2	1:G:127:PHE:HB3	2.03	0.40
1:P:94:VAL:HG13	1:P:96:TRP:CD1	2.56	0.40
3:C:289:GLY:C	3:C:291:MET:N	2.79	0.40
1:J:3:ASN:N	1:J:3:ASN:HD22	2.20	0.40
3:L:291:MET:O	3:L:292:LEU:C	2.64	0.40
2:N:62:SER:HA	2:N:98:MET:O	2.21	0.40
1:M:100:GLU:CD	2:N:108:PHE:HB2	2.46	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%)	109 (84%)	17 (13%)	3 (2%)	5	19
1	D	129/135 (96%)	113 (88%)	15 (12%)	1 (1%)	16	42
1	G	129/135 (96%)	112 (87%)	15 (12%)	2 (2%)	7	26
1	J	129/135 (96%)	108 (84%)	17 (13%)	4 (3%)	3	14
1	M	129/135 (96%)	114 (88%)	12 (9%)	3 (2%)	5	19
1	P	129/135 (96%)	111 (86%)	17 (13%)	1 (1%)	16	42
2	B	120/124 (97%)	105 (88%)	14 (12%)	1 (1%)	16	42
2	E	120/124 (97%)	105 (88%)	13 (11%)	2 (2%)	7	25
2	H	120/124 (97%)	105 (88%)	13 (11%)	2 (2%)	7	25
2	K	120/124 (97%)	100 (83%)	19 (16%)	1 (1%)	16	42
2	N	120/124 (97%)	103 (86%)	14 (12%)	3 (2%)	4	17
2	Q	120/124 (97%)	104 (87%)	11 (9%)	5 (4%)	2	10
3	C	189/217 (87%)	175 (93%)	11 (6%)	3 (2%)	7	26
3	F	189/217 (87%)	178 (94%)	8 (4%)	3 (2%)	7	26
3	I	189/217 (87%)	175 (93%)	13 (7%)	1 (0%)	24	52
3	L	189/217 (87%)	169 (89%)	18 (10%)	2 (1%)	11	34
3	O	189/217 (87%)	174 (92%)	13 (7%)	2 (1%)	11	34
3	R	189/217 (87%)	175 (93%)	11 (6%)	3 (2%)	7	26
All	All	2628/2856 (92%)	2335 (89%)	251 (10%)	42 (2%)	7	26

All (42) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	76	ARG
1	G	88	SER
2	H	2	PRO
2	N	2	PRO

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Mol	Chain	Res	Type
3	O	323	LYS
2	Q	3	LEU
2	Q	117	SER
3	R	323	LYS
1	A	88	SER
2	B	33	HIS
2	E	2	PRO
1	J	76	ARG
1	J	88	SER
2	K	3	LEU
3	L	290	SER
1	M	88	SER
2	N	33	HIS
2	N	117	SER
2	Q	2	PRO
2	Q	33	HIS
3	C	217	ALA
1	D	132	PRO
3	F	286	SER
3	F	334	SER
2	H	8	TYR
3	I	334	SER
3	L	217	ALA
1	P	88	SER
3	C	290	SER
2	E	8	TYR
1	M	36	ALA
1	M	120	ASP
3	R	326	ILE
1	J	24	THR
1	J	126	PRO
3	F	323	LYS
3	O	326	ILE
2	Q	9	ASN
3	R	217	ALA
3	C	313	GLY
1	A	126	PRO
1	G	132	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	114/118 (97%)	110 (96%)	4 (4%)	32	59
1	D	114/118 (97%)	108 (95%)	6 (5%)	20	47
1	G	114/118 (97%)	104 (91%)	10 (9%)	9	30
1	J	114/118 (97%)	110 (96%)	4 (4%)	32	59
1	M	114/118 (97%)	107 (94%)	7 (6%)	17	42
1	P	114/118 (97%)	110 (96%)	4 (4%)	32	59
2	B	110/112 (98%)	102 (93%)	8 (7%)	13	36
2	E	110/112 (98%)	101 (92%)	9 (8%)	10	32
2	H	110/112 (98%)	105 (96%)	5 (4%)	24	52
2	K	110/112 (98%)	102 (93%)	8 (7%)	13	36
2	N	110/112 (98%)	101 (92%)	9 (8%)	10	32
2	Q	110/112 (98%)	102 (93%)	8 (7%)	13	36
3	C	160/181 (88%)	147 (92%)	13 (8%)	11	33
3	F	160/181 (88%)	147 (92%)	13 (8%)	11	33
3	I	160/181 (88%)	148 (92%)	12 (8%)	12	36
3	L	160/181 (88%)	153 (96%)	7 (4%)	25	53
3	O	160/181 (88%)	154 (96%)	6 (4%)	29	57
3	R	160/181 (88%)	155 (97%)	5 (3%)	35	61
All	All	2304/2466 (93%)	2166 (94%)	138 (6%)	17	43

All (138) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	43	SER
1	A	58	THR
1	A	87	MET
1	A	90	SER
2	B	34	LYS
2	B	55	SER

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Mol	Chain	Res	Type
2	B	63	MET
2	B	67	LEU
2	B	91	ARG
2	B	102	THR
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	321	ASP
3	C	329	ILE
3	C	334	SER
3	C	339	ARG
3	C	343	ASN
3	C	344	VAL
3	C	345	SER
3	C	362	PHE
1	D	13	GLN
1	D	17	GLN
1	D	77	LYS
1	D	86	SER
1	D	88	SER
1	D	118	TYR
2	E	32	GLN
2	E	55	SER
2	E	59	LYS
2	E	67	LEU
2	E	79	SER
2	E	81	ASP
2	E	102	THR
2	E	117	SER
2	E	122	PHE
3	F	184	SER
3	F	207	THR
3	F	218	ASN
3	F	231	THR
3	F	239	THR
3	F	250	THR
3	F	290	SER
3	F	321	ASP

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Mol	Chain	Res	Type
3	F	343	ASN
3	F	344	VAL
3	F	345	SER
3	F	350	LEU
3	F	360	GLN
1	G	3	ASN
1	G	4	CYS
1	G	12	ASP
1	G	30	ARG
1	G	43	SER
1	G	84	GLU
1	G	86	SER
1	G	90	SER
1	G	92	ILE
1	G	118	TYR
2	H	21	THR
2	H	55	SER
2	H	59	LYS
2	H	63	MET
2	H	79	SER
3	I	184	SER
3	I	192	LYS
3	I	231	THR
3	I	232	LYS
3	I	283	ASP
3	I	290	SER
3	I	321	ASP
3	I	343	ASN
3	I	344	VAL
3	I	345	SER
3	I	350	LEU
3	I	360	GLN
1	J	3	ASN
1	J	21	GLU
1	J	43	SER
1	J	87	MET
2	K	17	SER
2	K	37	ARG
2	K	67	LEU
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	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	58	THR
1	M	88	SER
1	M	98	THR
2	N	17	SER
2	N	32	GLN
2	N	33	HIS
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	207	THR
3	O	321	ASP
3	O	334	SER
3	O	344	VAL
3	O	345	SER
3	O	352	GLU
1	P	21	GLU
1	P	58	THR
1	P	92	ILE
1	P	93	TYR
2	Q	17	SER
2	Q	32	GLN
2	Q	34	LYS
2	Q	59	LYS
2	Q	88	VAL
2	Q	91	ARG
2	Q	119	VAL

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Mol	Chain	Res	Type
2	Q	122	PHE
3	R	268	SER
3	R	321	ASP
3	R	344	VAL
3	R	345	SER
3	R	352	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (48) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	60	ASN
2	B	56	GLN
3	C	200	GLN
3	C	210	GLN
3	C	298	GLN
3	C	343	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	200	GLN
3	F	210	GLN
3	F	218	ASN
3	F	225	ASN
3	F	241	GLN
3	F	257	GLN
3	F	303	ASN
3	F	316	ASN
3	F	343	ASN
1	G	40	HIS
1	G	60	ASN
1	G	79	GLN
2	H	109	ASN
3	I	200	GLN
3	I	210	GLN
3	I	219	ASN
3	I	303	ASN
3	I	316	ASN
3	I	360	GLN
1	J	60	ASN

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Mol	Chain	Res	Type
2	K	33	HIS
2	K	109	ASN
3	L	210	GLN
3	L	318	ASN
1	M	60	ASN
2	N	32	GLN
2	N	33	HIS
2	N	56	GLN
3	O	200	GLN
3	O	225	ASN
3	O	318	ASN
2	Q	32	GLN
2	Q	33	HIS
2	Q	56	GLN
3	R	225	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	H	202	5,6	4,4,4	0.33	0	6,6,6	0.25	0
4	SO4	N	201	6	4,4,4	0.32	0	6,6,6	0.11	0
4	SO4	Q	202	6	4,4,4	0.33	0	6,6,6	0.10	0
4	SO4	B	201	6	4,4,4	0.36	0	6,6,6	0.16	0
4	SO4	D	201	-	4,4,4	0.32	0	6,6,6	0.11	0
4	SO4	J	201	-	4,4,4	0.25	0	6,6,6	0.10	0
4	SO4	K	202	6	4,4,4	0.33	0	6,6,6	0.09	0
4	SO4	Q	201	-	4,4,4	0.25	0	6,6,6	0.15	0
4	SO4	R	402	-	4,4,4	0.30	0	6,6,6	0.15	0
4	SO4	A	201	-	4,4,4	0.28	0	6,6,6	0.14	0
4	SO4	G	202	-	4,4,4	0.31	0	6,6,6	0.17	0
4	SO4	A	202	-	4,4,4	0.33	0	6,6,6	0.07	0
4	SO4	O	403	-	4,4,4	0.33	0	6,6,6	0.08	0
4	SO4	E	201	-	4,4,4	0.31	0	6,6,6	0.14	0
4	SO4	E	203	5,6	4,4,4	0.33	0	6,6,6	0.33	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 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	O	403	SO4	2	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%)	-1.13	0 100 100	50, 71, 100, 109	0
1	D	131/135 (97%)	-1.12	0 100 100	43, 69, 105, 125	0
1	G	131/135 (97%)	-1.15	0 100 100	44, 69, 103, 128	0
1	J	131/135 (97%)	-1.16	0 100 100	47, 71, 99, 110	0
1	M	131/135 (97%)	-0.95	0 100 100	48, 82, 132, 163	0
1	P	131/135 (97%)	-0.96	0 100 100	49, 81, 128, 153	0
2	B	122/124 (98%)	-0.87	0 100 100	54, 93, 135, 148	0
2	E	122/124 (98%)	-1.01	0 100 100	47, 76, 107, 133	0
2	H	122/124 (98%)	-0.99	0 100 100	50, 75, 110, 134	0
2	K	122/124 (98%)	-0.82	0 100 100	56, 89, 136, 153	0
2	N	122/124 (98%)	-1.00	0 100 100	52, 78, 113, 132	0
2	Q	122/124 (98%)	-0.95	0 100 100	51, 77, 109, 138	0
3	C	191/217 (88%)	-1.12	0 100 100	47, 77, 104, 148	0
3	F	191/217 (88%)	-1.15	0 100 100	46, 70, 107, 118	0
3	I	191/217 (88%)	-1.12	0 100 100	45, 70, 105, 132	0
3	L	191/217 (88%)	-1.13	0 100 100	52, 76, 106, 161	0
3	O	191/217 (88%)	-1.15	0 100 100	44, 67, 98, 122	0
3	R	191/217 (88%)	-1.19	0 100 100	43, 67, 99, 127	0
All	All	2664/2856 (93%)	-1.07	0 100 100	43, 74, 112, 163	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 ⓘ

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	0.12	200,203,216,238	0
7	CL	G	201	1/1	0.95	0.09	75,75,75,75	0
7	CL	L	403	1/1	0.95	0.10	77,77,77,77	0
4	SO4	N	201	5/5	0.96	0.08	97,100,116,125	0
7	CL	C	403	1/1	0.96	0.05	82,82,82,82	0
4	SO4	E	201	5/5	0.97	0.06	90,92,108,116	0
4	SO4	O	403	5/5	0.97	0.04	114,115,137,141	0
7	CL	H	201	1/1	0.97	0.13	77,77,77,77	0
7	CL	K	201	1/1	0.97	0.03	92,92,92,92	0
4	SO4	Q	201	5/5	0.97	0.06	69,82,97,107	0
7	CL	R	403	1/1	0.97	0.03	75,75,75,75	0
6	NA	O	402	1/1	0.98	0.05	65,65,65,65	0
4	SO4	K	202	5/5	0.98	0.04	82,82,110,119	0
7	CL	E	202	1/1	0.98	0.08	73,73,73,73	0
4	SO4	B	201	5/5	0.98	0.05	80,82,126,135	0
4	SO4	A	201	5/5	0.98	0.08	79,80,88,101	0
4	SO4	J	201	5/5	0.98	0.07	71,79,89,104	0
4	SO4	Q	202	5/5	0.98	0.06	100,104,114,126	0
7	CL	M	201	1/1	0.98	0.04	71,71,71,71	0
6	NA	I	402	1/1	0.98	0.04	50,50,50,50	0
6	NA	L	402	1/1	0.99	0.03	46,46,46,46	0
4	SO4	H	202	5/5	0.99	0.03	75,89,108,116	0
6	NA	R	404	1/1	0.99	0.05	72,72,72,72	0
4	SO4	D	201	5/5	0.99	0.03	59,70,84,93	0
4	SO4	E	203	5/5	0.99	0.04	63,76,107,112	0
4	SO4	R	402	5/5	0.99	0.04	64,81,101,104	5
5	SR	L	401	1/1	0.99	0.02	103,103,103,103	0
7	CL	J	202	1/1	0.99	0.09	87,87,87,87	0
5	SR	O	401	1/1	0.99	0.03	143,143,143,143	0
6	NA	C	402	1/1	0.99	0.03	46,46,46,46	0
6	NA	F	402	1/1	0.99	0.04	41,41,41,41	0
4	SO4	G	202	5/5	0.99	0.03	65,65,78,100	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
8	NH4	D	202	1/1	0.99	0.03	43,43,43,43	0
5	SR	I	401	1/1	1.00	0.02	97,97,97,97	0
5	SR	C	401	1/1	1.00	0.02	104,104,104,104	0
5	SR	F	401	1/1	1.00	0.02	95,95,95,95	0
5	SR	R	401	1/1	1.00	0.03	138,138,138,138	0

6.5 Other polymers [i](#)

There are no such residues in this entry.