



Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 11:53 am BST

PDB ID : 1N4R
Title : Protein Geranylgeranyltransferase type-I Complexed with a Geranylgeranylated KKKSKTKCVIL Peptide Product
Authors : Taylor, J.S.; Reid, T.S.; Casey, P.J.; Beese, L.S.
Deposited on : 2002-11-01
Resolution : 2.80 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

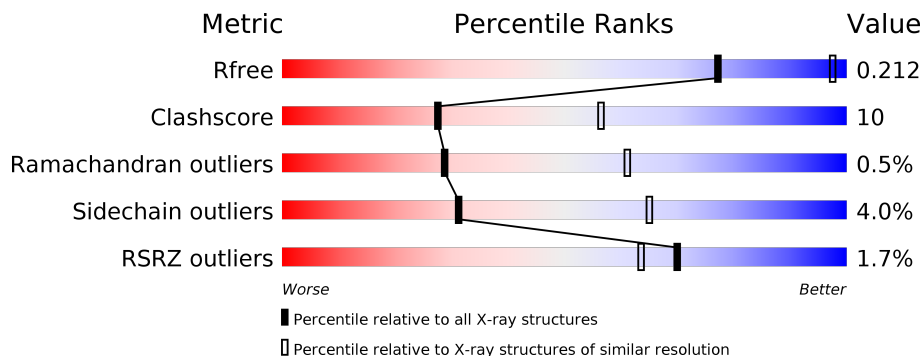
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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}	130704	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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 on 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	377	 2% 64% 19% 17%
1	C	377	 2% 65% 17% 17%
1	E	377	 2% 62% 20% 17%
1	G	377	 2% 65% 17% 17%
1	I	377	 2% 62% 20% 17%
1	K	377	 65% 17% 17%

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Mol	Chain	Length	Quality of chain
2	B	377	
2	D	377	
2	F	377	
2	H	377	
2	J	377	
2	L	377	
3	M	11	
3	N	11	
3	O	11	
3	P	11	
3	Q	11	
3	R	11	

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
7	CL	G	804	-	-	X	-

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 33129 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 Protein farnesyltransferase/geranylgeranyltransferase type-1 subunit alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	314	Total	C	N	O	S	0	0	0
			2614	1673	458	478	5			
1	C	314	Total	C	N	O	S	0	0	0
			2648	1691	462	490	5			
1	E	314	Total	C	N	O	S	0	0	0
			2630	1682	460	483	5			
1	G	314	Total	C	N	O	S	0	0	0
			2632	1683	459	485	5			
1	I	314	Total	C	N	O	S	0	0	0
			2654	1693	461	495	5			
1	K	314	Total	C	N	O	S	0	0	0
			2667	1700	466	496	5			

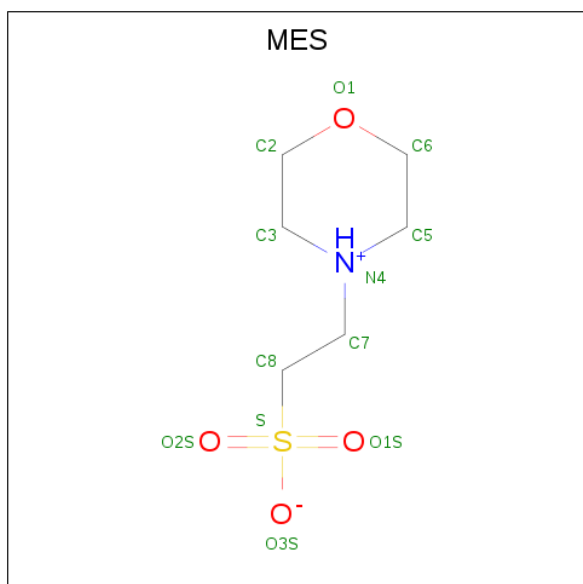
- Molecule 2 is a protein called Geranylgeranyl transferase type-1 subunit beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	346	Total	C	N	O	S	0	0	0
			2689	1702	467	496	24			
2	D	346	Total	C	N	O	S	0	0	0
			2693	1704	467	498	24			
2	F	346	Total	C	N	O	S	0	0	0
			2708	1711	471	502	24			
2	H	346	Total	C	N	O	S	0	0	0
			2670	1692	458	496	24			
2	J	346	Total	C	N	O	S	0	0	0
			2709	1711	471	503	24			
2	L	346	Total	C	N	O	S	0	0	0
			2714	1715	471	504	24			

- Molecule 3 is a protein called Fusion protein consisting of transforming protein p21b and Ras related protein Rap-2b.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	M	6	Total 46	C 30	N 7	O 8	S 1	0	0	0
3	N	6	Total 46	C 30	N 7	O 8	S 1	0	0	0
3	O	6	Total 46	C 30	N 7	O 8	S 1	0	0	0
3	P	6	Total 46	C 30	N 7	O 8	S 1	0	0	0
3	Q	6	Total 46	C 30	N 7	O 8	S 1	0	0	0
3	R	6	Total 46	C 30	N 7	O 8	S 1	0	0	0

- Molecule 4 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (three-letter code: MES) (formula: C₆H₁₃NO₄S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	S		
4	A	1	Total 12	C 6	N 1	O 4	S 1	0	0
4	C	1	Total 12	C 6	N 1	O 4	S 1	0	0
4	E	1	Total 12	C 6	N 1	O 4	S 1	0	0
4	G	1	Total 12	C 6	N 1	O 4	S 1	0	0
4	I	1	Total 12	C 6	N 1	O 4	S 1	0	0

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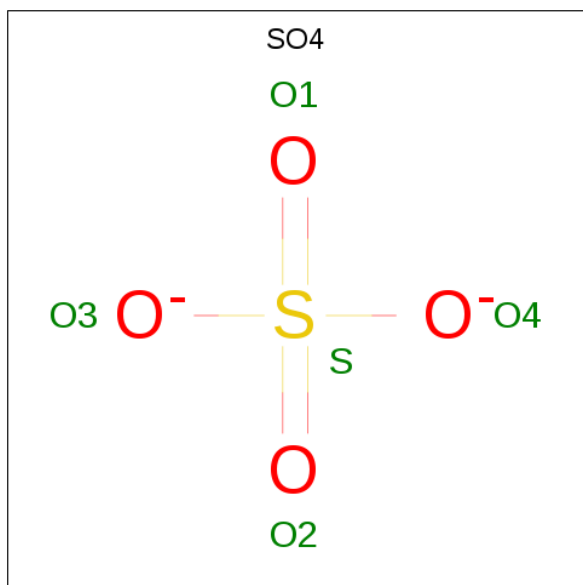
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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	S		
4	K	1	12	6	1	4	1	0	0

- Molecule 5 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	J	1	Total	Zn	0	0
			1	1		
5	D	1	Total	Zn	0	0
			1	1		
5	H	1	Total	Zn	0	0
			1	1		
5	B	1	Total	Zn	0	0
			1	1		
5	L	1	Total	Zn	0	0
			1	1		
5	F	1	Total	Zn	0	0
			1	1		

- Molecule 6 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	B	1	Total	O	S	0	0
			5	4	1		
6	D	1	Total	O	S	0	0
			5	4	1		

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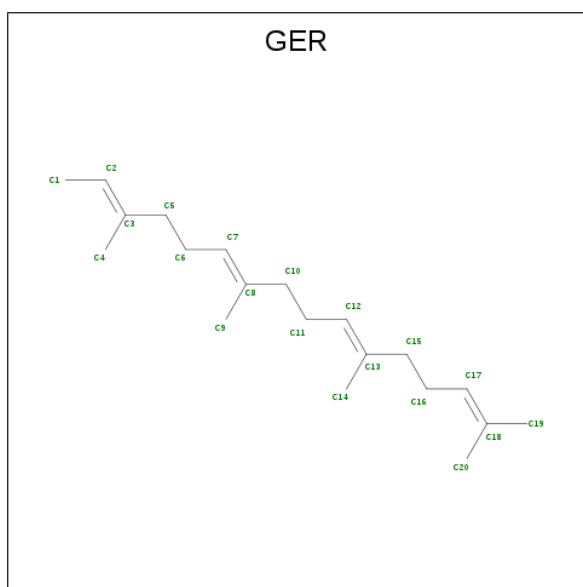
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	F	1	Total	O	S	0	0
			5	4	1		
6	H	1	Total	O	S	0	0
			5	4	1		
6	J	1	Total	O	S	0	0
			5	4	1		
6	L	1	Total	O	S	0	0
			5	4	1		

- Molecule 7 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	G	1	Total	Cl	0	0
			1	1		
7	J	1	Total	Cl	0	0
			1	1		
7	D	1	Total	Cl	0	0
			1	1		
7	K	1	Total	Cl	0	0
			1	1		
7	H	1	Total	Cl	0	0
			1	1		
7	C	1	Total	Cl	0	0
			1	1		
7	F	1	Total	Cl	0	0
			1	1		

- Molecule 8 is GERAN-8-YL GERAN (three-letter code: GER) (formula: C₂₀H₃₄).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	M	1	Total C 20 20	0	0
8	N	1	Total C 20 20	0	0
8	O	1	Total C 20 20	0	0
8	P	1	Total C 20 20	0	0
8	Q	1	Total C 20 20	0	0
8	R	1	Total C 20 20	0	0

- Molecule 9 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	A	45	Total O 45 45	0	0
9	B	35	Total O 35 35	0	0
9	C	49	Total O 49 49	0	0
9	D	45	Total O 45 45	0	0
9	E	40	Total O 40 40	0	0
9	F	51	Total O 51 51	0	0

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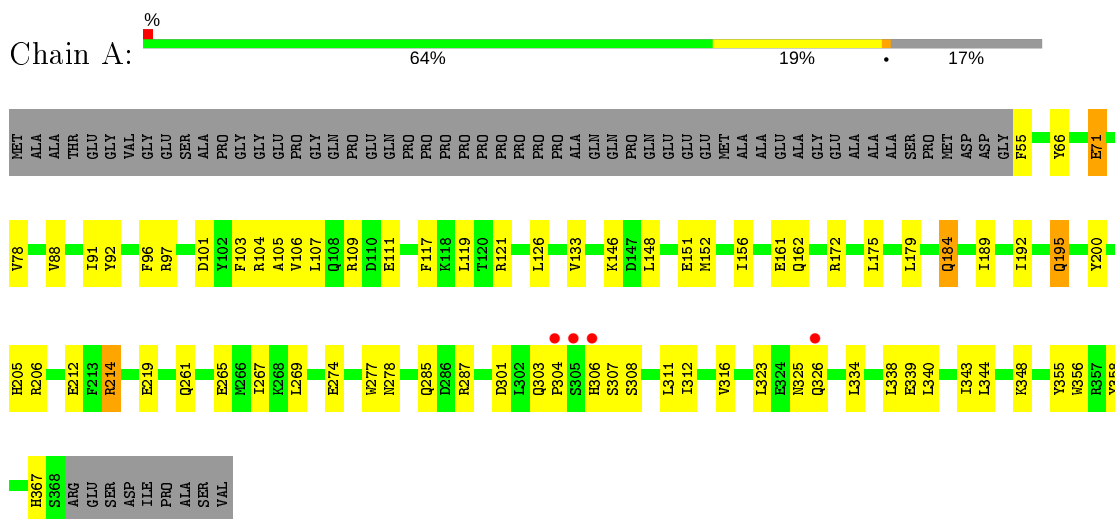
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	G	40	Total 40	O 40	0	0
9	H	23	Total 23	O 23	0	0
9	I	52	Total 52	O 52	0	0
9	J	36	Total 36	O 36	0	0
9	K	101	Total 101	O 101	0	0
9	L	60	Total 60	O 60	0	0
9	M	1	Total 1	O 1	0	0
9	N	3	Total 3	O 3	0	0
9	O	1	Total 1	O 1	0	0
9	Q	3	Total 3	O 3	0	0
9	R	5	Total 5	O 5	0	0

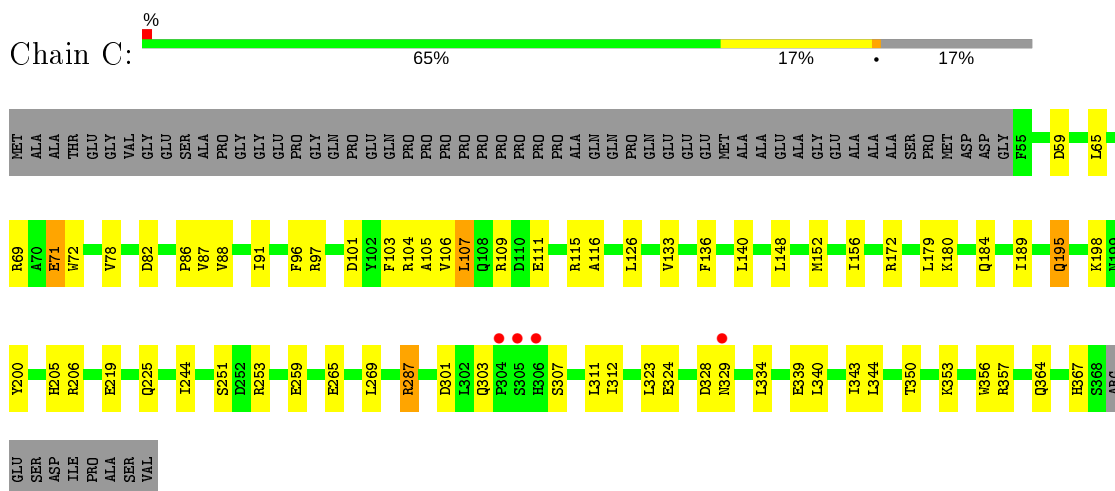
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA 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: Protein farnesyltransferase/geranylgeranyltransferase type-1 subunit alpha

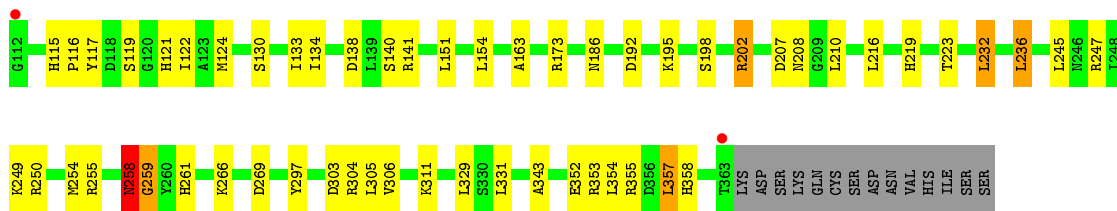


- Molecule 1: Protein farnesyltransferase/geranylgeranyltransferase type-1 subunit alpha

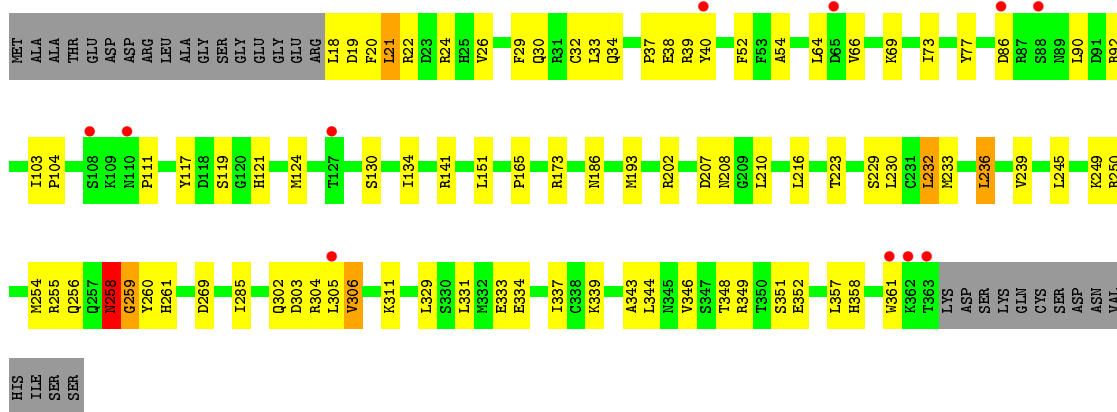


- Molecule 1: Protein farnesyltransferase/geranylgeranyltransferase type-1 subunit alpha

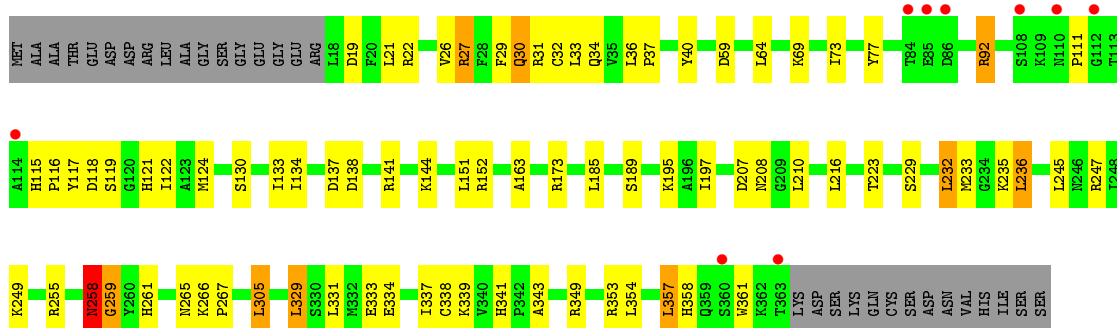




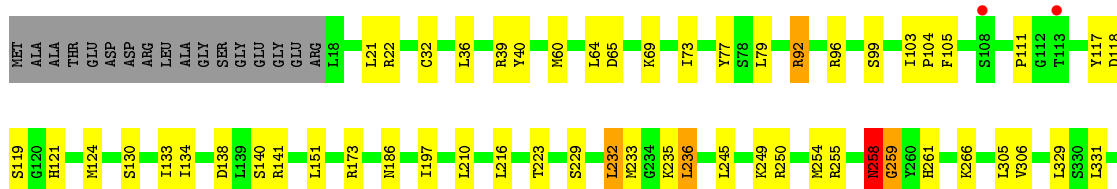
• Molecule 2: Geranylgeranyl transferase type-1 subunit beta

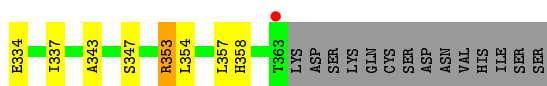


• Molecule 2: Geranylgeranyl transferase type-1 subunit beta

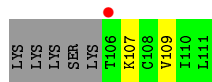
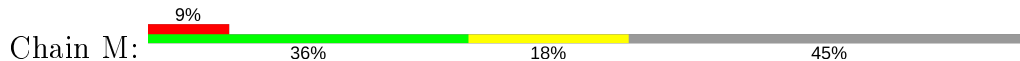


• Molecule 2: Geranylgeranyl transferase type-1 subunit beta

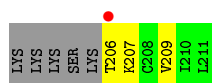
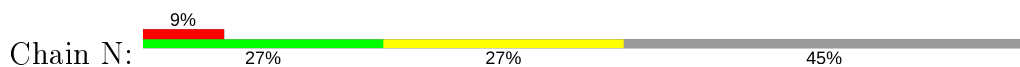




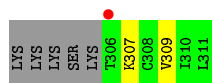
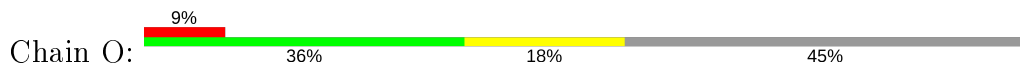
- Molecule 3: Fusion protein consisting of transforming protein p21b and Ras related protein Rap-2b



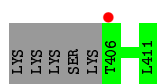
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- Molecule 3: Fusion protein consisting of transforming protein p21b and Ras related protein Rap-2b



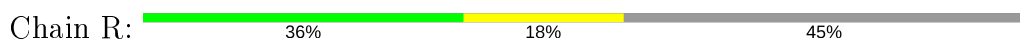
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- Molecule 3: Fusion protein consisting of transforming protein p21b and Ras related protein Rap-2b



- Molecule 3: Fusion protein consisting of transforming protein p21b and Ras related protein Rap-2b



LYS	T606
LYS	K607
LYS	G608
LYS	V609
SER	I610
LYS	L611

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	272.07Å 268.80Å 185.31Å 90.00° 131.55° 90.00°	Depositor
Resolution (Å)	29.86 – 2.80 29.86 – 2.60	Depositor EDS
% Data completeness (in resolution range)	99.1 (29.86-2.80) 95.3 (29.86-2.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.45 (at 2.61Å)	Xtrriage
Refinement program	CNS	Depositor
R, R_{free}	0.200 , 0.218 0.194 , 0.212	Depositor DCC
R_{free} test set	12102 reflections (4.19%)	wwPDB-VP
Wilson B-factor (Å ²)	51.9	Xtrriage
Anisotropy	0.040	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 49.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	0.085 for -h-2*1,-k,l	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	33129	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.61% 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: GER, ZN, SO4, MES, CL

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.33	0/2680	0.52	0/3649
1	C	0.36	0/2714	0.53	0/3690
1	E	0.35	0/2696	0.53	0/3668
1	G	0.36	0/2698	0.53	0/3670
1	I	0.36	0/2720	0.54	0/3698
1	K	0.39	0/2733	0.56	0/3713
2	B	0.36	0/2750	0.60	2/3720 (0.1%)
2	D	0.37	0/2754	0.59	2/3725 (0.1%)
2	F	0.38	0/2769	0.60	2/3743 (0.1%)
2	H	0.35	0/2730	0.58	2/3696 (0.1%)
2	J	0.36	0/2770	0.59	2/3745 (0.1%)
2	L	0.39	0/2775	0.61	2/3750 (0.1%)
3	M	0.54	0/45	0.50	0/58
3	N	0.52	0/45	0.51	0/58
3	O	0.54	0/45	0.53	0/58
3	P	0.50	0/45	0.51	0/58
3	Q	0.55	0/45	0.52	0/58
3	R	0.56	0/45	0.53	0/58
All	All	0.37	0/33059	0.57	12/44815 (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
2	B	0	1
2	F	0	1
All	All	0	2

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	J	259	GLY	N-CA-C	-5.88	98.40	113.10
2	F	259	GLY	N-CA-C	-5.80	98.61	113.10
2	L	259	GLY	N-CA-C	-5.79	98.62	113.10
2	H	259	GLY	N-CA-C	-5.75	98.73	113.10
2	D	259	GLY	N-CA-C	-5.65	98.97	113.10
2	B	258	ASN	N-CA-C	-5.56	95.99	111.00
2	B	259	GLY	N-CA-C	-5.53	99.28	113.10
2	J	258	ASN	N-CA-C	-5.46	96.26	111.00
2	L	258	ASN	N-CA-C	-5.42	96.38	111.00
2	D	258	ASN	N-CA-C	-5.38	96.48	111.00
2	F	258	ASN	N-CA-C	-5.38	96.48	111.00
2	H	258	ASN	N-CA-C	-5.34	96.57	111.00

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	297	TYR	Sidechain
2	F	297	TYR	Sidechain

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	2614	0	2499	58	0
1	C	2648	0	2549	50	0
1	E	2630	0	2524	69	0
1	G	2632	0	2527	57	0
1	I	2654	0	2551	54	0
1	K	2667	0	2577	54	0
2	B	2689	0	2585	55	0
2	D	2693	0	2589	64	0
2	F	2708	0	2613	48	0
2	H	2670	0	2551	66	0
2	J	2709	0	2610	46	0
2	L	2714	0	2621	39	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	M	46	0	54	1	0
3	N	46	0	54	2	0
3	O	46	0	54	1	0
3	P	46	0	54	0	0
3	Q	46	0	54	1	0
3	R	46	0	54	1	0
4	A	12	0	13	3	0
4	C	12	0	13	1	0
4	E	12	0	13	2	0
4	G	12	0	13	2	0
4	I	12	0	13	2	0
4	K	12	0	13	1	0
5	B	1	0	0	0	0
5	D	1	0	0	0	0
5	F	1	0	0	0	0
5	H	1	0	0	0	0
5	J	1	0	0	0	0
5	L	1	0	0	0	0
6	B	5	0	0	0	0
6	D	5	0	0	0	0
6	F	5	0	0	0	0
6	H	5	0	0	0	0
6	J	5	0	0	0	0
6	L	5	0	0	0	0
7	C	1	0	0	1	0
7	D	1	0	0	0	0
7	F	1	0	0	0	0
7	G	1	0	0	2	0
7	H	1	0	0	0	0
7	J	1	0	0	0	0
7	K	1	0	0	1	0
8	M	20	0	32	1	0
8	N	20	0	32	1	0
8	O	20	0	32	1	0
8	P	20	0	32	1	0
8	Q	20	0	32	1	0
8	R	20	0	32	2	0
9	A	45	0	0	3	0
9	B	35	0	0	1	0
9	C	49	0	0	2	0
9	D	45	0	0	2	0
9	E	40	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
9	F	51	0	0	2	0
9	G	40	0	0	4	0
9	H	23	0	0	0	0
9	I	52	0	0	0	0
9	J	36	0	0	2	0
9	K	101	0	0	4	0
9	L	60	0	0	0	0
9	M	1	0	0	0	0
9	N	3	0	0	0	0
9	O	1	0	0	0	0
9	Q	3	0	0	0	0
9	R	5	0	0	0	0
All	All	33129	0	31390	639	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 (639) 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:E:156:ILE:HG12	1:E:172:ARG:HH12	1.12	1.14
1:K:156:ILE:HG12	1:K:172:ARG:HH12	0.99	1.14
1:A:156:ILE:HG12	1:A:172:ARG:HH12	1.12	1.12
1:I:156:ILE:HG12	1:I:172:ARG:HH12	1.15	1.11
1:C:156:ILE:HG12	1:C:172:ARG:HH12	1.16	1.09
1:G:156:ILE:HG12	1:G:172:ARG:HH12	1.26	0.97
1:K:156:ILE:HG12	1:K:172:ARG:NH1	1.83	0.94
7:G:804:CL:CL	9:G:913:HOH:O	2.23	0.92
1:E:81:ASN:N	1:E:81:ASN:HD22	1.70	0.89
1:C:353:LYS:HE2	1:C:357:ARG:HH12	1.38	0.86
1:E:81:ASN:HD22	1:E:81:ASN:H	1.25	0.84
2:H:21:LEU:HD13	2:H:302:GLN:HE22	1.42	0.82
2:H:21:LEU:HD22	2:H:302:GLN:HE21	1.41	0.82
1:E:353:LYS:HE3	1:E:357:ARG:HH22	1.44	0.82
2:D:69:LYS:O	2:D:73:ILE:HG13	1.82	0.79
2:H:348:THR:O	2:H:352:GLU:HG2	1.83	0.79
1:A:156:ILE:HG12	1:A:172:ARG:NH1	1.94	0.78
2:B:92:ARG:NH1	2:B:119:SER:HB3	2.00	0.76
2:B:39:ARG:HB3	2:B:39:ARG:HH11	1.51	0.76
2:F:69:LYS:O	2:F:73:ILE:HG13	1.86	0.76
1:C:353:LYS:HE2	1:C:357:ARG:NH1	2.02	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:343:ILE:HG22	1:G:348:LYS:HG3	1.68	0.75
1:K:107:LEU:HD22	2:L:117:TYR:CD2	2.21	0.75
1:A:152:MET:O	1:A:156:ILE:HG13	1.86	0.75
2:H:229:SER:O	2:H:233:MET:HG3	1.86	0.75
2:J:92:ARG:HH11	2:J:119:SER:HB3	1.52	0.74
1:G:152:MET:O	1:G:156:ILE:HG13	1.88	0.74
1:A:214:ARG:HG2	1:G:180:LYS:HB2	1.71	0.72
2:F:37:PRO:HD2	2:F:40:TYR:CD1	2.23	0.72
1:K:156:ILE:CG1	1:K:172:ARG:HH12	1.92	0.72
1:E:156:ILE:CG1	1:E:172:ARG:HH12	1.99	0.71
1:I:152:MET:O	1:I:156:ILE:HG13	1.90	0.71
7:K:807:CL:CL	9:K:986:HOH:O	2.46	0.71
2:B:69:LYS:O	2:B:73:ILE:HG13	1.89	0.71
1:I:91:ILE:HD12	1:I:91:ILE:O	1.90	0.70
2:J:92:ARG:NH1	2:J:119:SER:HB3	2.05	0.70
1:K:152:MET:O	1:K:156:ILE:HG13	1.91	0.70
1:E:152:MET:O	1:E:156:ILE:HG13	1.91	0.70
1:C:152:MET:O	1:C:156:ILE:HG13	1.90	0.70
1:E:353:LYS:HE2	1:E:357:ARG:HH12	1.56	0.70
1:E:156:ILE:HG12	1:E:172:ARG:NH1	1.97	0.69
1:I:107:LEU:HD22	2:J:117:TYR:CD2	2.27	0.69
2:H:21:LEU:HD13	2:H:302:GLN:NE2	2.07	0.68
2:H:69:LYS:O	2:H:73:ILE:HG13	1.92	0.68
7:G:804:CL:CL	9:G:932:HOH:O	2.47	0.68
1:A:339:GLU:O	1:A:343:ILE:HG13	1.93	0.68
1:E:105:ALA:O	1:E:109:ARG:HG3	1.94	0.68
7:C:801:CL:CL	9:C:922:HOH:O	2.47	0.68
1:A:91:ILE:O	1:A:91:ILE:HD12	1.94	0.67
2:B:186:ASN:HB2	2:B:358:HIS:CE1	2.29	0.67
1:C:189:ILE:HD11	1:C:205:HIS:HD2	1.60	0.67
1:C:323:LEU:HD13	1:C:367:HIS:HD2	1.60	0.67
1:A:340:LEU:HD23	1:A:343:ILE:HD12	1.77	0.67
1:E:81:ASN:N	1:E:81:ASN:ND2	2.42	0.67
1:E:91:ILE:O	1:E:91:ILE:HD12	1.95	0.67
2:H:349:ARG:O	2:H:352:GLU:HB2	1.94	0.67
2:F:37:PRO:HD2	2:F:40:TYR:CE1	2.30	0.65
1:K:91:ILE:O	1:K:91:ILE:HD12	1.97	0.65
2:B:92:ARG:HH11	2:B:119:SER:HB3	1.61	0.65
1:G:91:ILE:HD12	1:G:91:ILE:O	1.96	0.65
1:E:189:ILE:HD11	1:E:205:HIS:HD2	1.62	0.65
1:G:334:LEU:HD22	1:G:367:HIS:O	1.96	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:255:VAL:HG13	1:E:258:ARG:NH2	2.12	0.65
1:G:97:ARG:HG2	1:G:101:ASP:OD2	1.96	0.65
1:G:189:ILE:HD11	1:G:205:HIS:HD2	1.62	0.64
1:C:91:ILE:HD12	1:C:91:ILE:O	1.98	0.64
2:F:30:GLN:O	2:F:34:GLN:HG3	1.97	0.64
1:I:323:LEU:HB3	1:I:367:HIS:CD2	2.32	0.64
1:A:323:LEU:HB3	1:A:367:HIS:CD2	2.33	0.64
1:I:334:LEU:HD22	1:I:367:HIS:O	1.97	0.64
2:D:37:PRO:HD2	2:D:40:TYR:CE1	2.34	0.63
1:G:252:ASP:OD2	1:G:255:VAL:HG23	1.98	0.62
2:B:24:ARG:HD3	2:B:27:ARG:HH12	1.62	0.62
1:I:148:LEU:HB2	1:I:179:LEU:HD21	1.82	0.62
2:J:22:ARG:O	2:J:26:VAL:HG23	1.99	0.62
2:H:245:LEU:O	2:H:249:LYS:HG3	1.98	0.62
2:H:21:LEU:HD22	2:H:302:GLN:NE2	2.13	0.62
1:K:189:ILE:HD11	1:K:205:HIS:HD2	1.64	0.62
1:A:189:ILE:HD11	1:A:205:HIS:HD2	1.65	0.61
2:D:133:ILE:HD13	2:D:354:LEU:HD13	1.81	0.61
1:G:148:LEU:HB2	1:G:179:LEU:HD21	1.82	0.61
2:B:37:PRO:HD2	2:B:40:TYR:CE1	2.35	0.61
1:C:339:GLU:O	1:C:343:ILE:HG13	2.00	0.61
1:G:312:ILE:HG23	1:G:340:LEU:HD22	1.83	0.61
2:B:245:LEU:O	2:B:249:LYS:HG3	2.01	0.61
1:C:148:LEU:HB2	1:C:179:LEU:HD21	1.81	0.61
1:E:320:GLU:OE2	1:E:363:LEU:HD21	2.01	0.61
2:H:193:MET:HG3	2:H:233:MET:CE	2.31	0.61
2:H:26:VAL:O	2:H:30:GLN:HG3	2.01	0.61
2:J:358:HIS:O	2:J:361:TRP:HB2	2.01	0.61
2:J:232:LEU:HD13	2:J:343:ALA:HB1	1.82	0.60
2:D:229:SER:O	2:D:233:MET:HG3	2.01	0.60
1:E:69:ARG:HB3	1:E:71:GLU:OE1	2.01	0.60
1:E:148:LEU:HB2	1:E:179:LEU:HD21	1.81	0.60
1:G:97:ARG:HH11	1:G:97:ARG:HB3	1.67	0.60
1:A:148:LEU:HB2	1:A:179:LEU:HD21	1.82	0.60
1:C:323:LEU:HD13	1:C:367:HIS:CD2	2.35	0.60
2:H:193:MET:HG3	2:H:233:MET:HE1	1.83	0.60
2:F:26:VAL:O	2:F:30:GLN:HG3	2.01	0.60
1:I:100:TYR:O	1:I:104:ARG:HG3	2.00	0.60
2:L:69:LYS:O	2:L:73:ILE:HG13	2.02	0.60
2:B:37:PRO:HD2	2:B:40:TYR:CD1	2.36	0.60
1:E:117:PHE:CE2	1:E:146:LYS:HE2	2.37	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:214:ARG:O	1:E:214:ARG:HG3	2.01	0.60
1:I:189:ILE:HD11	1:I:205:HIS:HD2	1.66	0.60
2:B:353:ARG:HH11	2:B:353:ARG:HG2	1.65	0.60
1:G:330:LYS:HE2	1:G:367:HIS:HB3	1.84	0.59
1:I:207:GLN:HG2	1:I:242:PHE:CE2	2.37	0.59
2:H:232:LEU:HD13	2:H:343:ALA:HB1	1.83	0.59
2:L:245:LEU:O	2:L:249:LYS:HG3	2.03	0.59
1:A:97:ARG:HG2	1:A:101:ASP:OD2	2.01	0.59
2:J:69:LYS:O	2:J:73:ILE:HG13	2.02	0.59
2:D:353:ARG:NH1	2:D:357:LEU:HG	2.16	0.59
2:J:92:ARG:NH1	2:J:118:ASP:O	2.33	0.59
1:E:78:VAL:O	1:E:104:ARG:HD2	2.02	0.59
1:G:156:ILE:HG12	1:G:172:ARG:NH1	2.08	0.59
2:F:303:ASP:OD1	2:F:306:VAL:HG13	2.03	0.58
2:H:18:LEU:N	2:H:18:LEU:HD22	2.17	0.58
2:H:303:ASP:OD1	2:H:306:VAL:HG13	2.03	0.58
1:G:101:ASP:HA	1:G:104:ARG:HH11	1.68	0.58
1:E:67:ARG:NH2	1:E:94:GLU:OE1	2.36	0.58
1:I:105:ALA:O	1:I:109:ARG:HG3	2.03	0.58
1:I:261:GLN:O	1:I:265:GLU:HG2	2.04	0.58
1:C:328:ASP:O	1:C:329:ASN:HB2	2.03	0.58
1:G:339:GLU:O	1:G:343:ILE:HG13	2.03	0.58
1:E:312:ILE:HG23	1:E:340:LEU:HD22	1.85	0.58
2:F:133:ILE:HD13	2:F:354:LEU:HD13	1.86	0.58
1:K:148:LEU:HB2	1:K:179:LEU:HD21	1.86	0.58
2:B:29:PHE:O	2:B:33:LEU:HD22	2.03	0.57
2:D:173:ARG:HG2	8:N:1208:GER:H111	1.85	0.57
1:G:261:GLN:O	1:G:265:GLU:HG2	2.04	0.57
2:F:138:ASP:HA	2:F:357:LEU:HD11	1.85	0.57
1:A:334:LEU:HD22	1:A:367:HIS:O	2.03	0.57
1:E:339:GLU:O	1:E:343:ILE:HG13	2.05	0.57
1:G:58:LEU:HD12	1:G:125:GLU:OE2	2.04	0.57
2:H:92:ARG:HD2	2:H:119:SER:HB3	1.85	0.57
1:K:334:LEU:HD22	1:K:367:HIS:O	2.04	0.57
2:F:173:ARG:HG2	8:O:1308:GER:H111	1.86	0.57
2:L:39:ARG:HG3	2:L:40:TYR:CE1	2.40	0.57
2:H:92:ARG:HG2	2:H:165:PRO:HG3	1.86	0.57
2:J:258:ASN:OD1	2:J:259:GLY:N	2.32	0.57
2:H:210:LEU:HB2	2:H:223:THR:HA	1.87	0.57
1:A:78:VAL:O	1:A:104:ARG:HD2	2.04	0.57
1:I:88:VAL:HG12	2:J:32:CYS:O	2.05	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:312:ILE:HG23	1:C:340:LEU:HD22	1.86	0.56
2:D:37:PRO:HD2	2:D:40:TYR:CD1	2.40	0.56
2:F:198:SER:O	2:F:202:ARG:HG3	2.05	0.56
2:D:296:ASN:HD22	2:D:296:ASN:C	2.07	0.56
2:D:333:GLU:HA	9:D:835:HOH:O	2.06	0.56
1:K:301:ASP:O	1:K:304:PRO:HD2	2.06	0.56
1:C:334:LEU:HD22	1:C:367:HIS:O	2.05	0.56
2:F:130:SER:O	2:F:134:ILE:HG13	2.06	0.56
1:G:101:ASP:HA	1:G:104:ARG:NH1	2.21	0.56
2:H:193:MET:HE2	2:H:233:MET:HB3	1.86	0.56
2:H:52:PHE:HE1	2:H:130:SER:HG	1.51	0.56
2:B:258:ASN:OD1	2:B:259:GLY:N	2.38	0.56
1:G:107:LEU:HD22	2:H:117:TYR:CD2	2.41	0.56
1:I:200:TYR:HB3	4:I:905:MES:H32	1.88	0.56
2:J:173:ARG:HG2	8:Q:1508:GER:H111	1.88	0.56
2:J:210:LEU:HB2	2:J:223:THR:HA	1.87	0.56
2:D:303:ASP:OD1	2:D:306:VAL:HG13	2.05	0.56
1:E:121:ARG:HG3	1:E:121:ARG:NH1	2.19	0.56
2:J:195:LYS:NZ	9:J:840:HOH:O	2.38	0.56
2:L:210:LEU:HB2	2:L:223:THR:HA	1.87	0.56
2:H:202:ARG:HG3	2:H:202:ARG:HH11	1.72	0.55
2:L:133:ILE:HD13	2:L:354:LEU:HD13	1.88	0.55
2:B:26:VAL:O	2:B:30:GLN:HG3	2.06	0.55
2:B:173:ARG:HG2	8:M:1108:GER:H111	1.86	0.55
2:D:26:VAL:O	2:D:30:GLN:HG3	2.06	0.55
1:C:303:GLN:O	1:C:307:SER:HB2	2.07	0.55
2:L:186:ASN:HB2	2:L:358:HIS:NE2	2.21	0.55
1:A:107:LEU:HD22	2:B:117:TYR:CD2	2.41	0.55
2:B:39:ARG:HH11	2:B:39:ARG:CB	2.19	0.55
2:F:210:LEU:HB2	2:F:223:THR:HA	1.88	0.55
2:L:232:LEU:HD13	2:L:343:ALA:HB1	1.88	0.55
2:B:210:LEU:HB2	2:B:223:THR:HA	1.88	0.55
2:D:210:LEU:HB2	2:D:223:THR:HA	1.88	0.55
1:A:303:GLN:O	1:A:307:SER:HB2	2.07	0.55
2:D:202:ARG:HG3	2:D:202:ARG:HH11	1.71	0.55
1:E:107:LEU:HD22	2:F:117:TYR:CD2	2.42	0.55
2:H:258:ASN:OD1	2:H:259:GLY:N	2.36	0.55
1:I:156:ILE:CG1	1:I:172:ARG:HH12	2.05	0.55
2:H:173:ARG:HG2	8:P:1408:GER:H111	1.89	0.55
2:D:30:GLN:O	2:D:34:GLN:HG3	2.06	0.54
1:E:334:LEU:HD22	1:E:367:HIS:O	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:149:GLN:NE2	1:K:179:LEU:HD13	2.23	0.54
2:D:121:HIS:HB3	2:D:124:MET:HG2	1.90	0.54
1:G:100:TYR:O	1:G:104:ARG:HG3	2.06	0.54
1:K:198:LYS:HD3	2:L:266:LYS:HD3	1.89	0.54
2:D:258:ASN:OD1	2:D:259:GLY:N	2.33	0.54
2:D:31:ARG:HH22	2:D:306:VAL:HG23	1.72	0.54
2:J:133:ILE:HD13	2:J:354:LEU:HD13	1.89	0.54
1:A:261:GLN:O	1:A:265:GLU:HG2	2.07	0.54
2:F:64:LEU:HD11	2:F:134:ILE:HG22	1.91	0.53
1:I:65:LEU:O	1:I:69:ARG:HG3	2.08	0.53
1:K:156:ILE:HD11	1:K:172:ARG:HH22	1.73	0.53
1:I:69:ARG:HB3	1:I:71:GLU:OE1	2.08	0.53
1:C:156:ILE:HG12	1:C:172:ARG:NH1	2.02	0.53
2:D:103:ILE:HG23	2:D:104:PRO:HD2	1.91	0.53
1:G:252:ASP:OD2	1:G:254:ALA:HB3	2.09	0.53
1:I:106:VAL:HG13	1:I:111:GLU:HB3	1.90	0.53
1:I:156:ILE:HG12	1:I:172:ARG:NH1	2.01	0.53
1:A:301:ASP:O	1:A:304:PRO:HD2	2.08	0.53
1:E:81:ASN:H	1:E:81:ASN:ND2	2.00	0.53
2:F:186:ASN:HB2	2:F:358:HIS:NE2	2.24	0.53
1:E:255:VAL:HG13	1:E:258:ARG:HH21	1.72	0.53
1:C:350:THR:O	1:C:353:LYS:HB2	2.08	0.53
2:D:19:ASP:OD2	2:D:19:ASP:N	2.42	0.53
1:E:58:LEU:HD22	1:E:95:LYS:HD3	1.90	0.53
2:L:186:ASN:HB2	2:L:358:HIS:CE1	2.44	0.53
2:D:20:PHE:CZ	2:D:337:ILE:HD11	2.44	0.53
1:K:323:LEU:HB3	1:K:367:HIS:CD2	2.44	0.53
1:E:121:ARG:HH11	1:E:121:ARG:HG3	1.74	0.53
1:C:311:LEU:HD23	1:C:311:LEU:C	2.30	0.52
2:D:138:ASP:OD1	2:D:140:SER:HB3	2.08	0.52
2:H:130:SER:O	2:H:134:ILE:HG13	2.09	0.52
2:J:121:HIS:HB3	2:J:124:MET:HG2	1.91	0.52
1:K:88:VAL:HG12	2:L:32:CYS:O	2.10	0.52
2:F:121:HIS:HB3	2:F:124:MET:HG2	1.92	0.52
2:F:232:LEU:HD13	2:F:343:ALA:HB1	1.92	0.52
1:G:303:GLN:O	1:G:307:SER:HB2	2.10	0.52
1:K:311:LEU:HD23	1:K:311:LEU:C	2.29	0.52
2:L:334:GLU:HB3	2:L:337:ILE:HD12	1.92	0.52
1:G:334:LEU:O	1:G:338:LEU:HG	2.09	0.52
1:G:88:VAL:HG12	2:H:32:CYS:O	2.09	0.52
1:A:92:TYR:O	1:A:97:ARG:NH2	2.42	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:18:LEU:HD12	2:D:304:ARG:NH1	2.25	0.52
1:G:65:LEU:O	1:G:69:ARG:HG3	2.09	0.52
2:H:21:LEU:HD23	2:H:24:ARG:HD2	1.90	0.52
1:C:69:ARG:HB3	1:C:71:GLU:OE1	2.09	0.52
2:L:121:HIS:HB3	2:L:124:MET:HG2	1.92	0.52
2:L:173:ARG:HG2	8:R:1608:GER:H111	1.91	0.52
1:K:312:ILE:HG23	1:K:340:LEU:HD22	1.92	0.52
1:E:353:LYS:HE3	1:E:357:ARG:NH2	2.20	0.52
2:J:77:TYR:CE1	2:J:141:ARG:HB2	2.45	0.51
1:A:274:GLU:HG2	1:A:278:ASN:HD21	1.74	0.51
1:E:92:TYR:O	1:E:97:ARG:NH2	2.43	0.51
1:I:311:LEU:C	1:I:311:LEU:HD23	2.31	0.51
2:J:130:SER:O	2:J:134:ILE:HG13	2.10	0.51
1:K:303:GLN:O	1:K:307:SER:HB2	2.11	0.51
1:C:78:VAL:O	1:C:104:ARG:HD2	2.09	0.51
2:H:357:LEU:HD21	2:H:361:TRP:CZ2	2.45	0.51
2:J:334:GLU:HB3	2:J:337:ILE:HD12	1.90	0.51
1:A:312:ILE:HG23	1:A:340:LEU:HD22	1.93	0.51
2:F:186:ASN:HB2	2:F:358:HIS:CE1	2.45	0.51
1:I:200:TYR:CB	4:I:905:MES:H32	2.40	0.51
2:J:138:ASP:HA	2:J:357:LEU:HD11	1.92	0.51
1:I:301:ASP:O	1:I:304:PRO:HD2	2.11	0.51
1:C:180:LYS:HB2	1:E:214:ARG:HG2	1.92	0.51
1:G:156:ILE:CG1	1:G:172:ARG:HH12	2.12	0.51
2:H:334:GLU:HB3	2:H:337:ILE:HD12	1.93	0.51
1:K:106:VAL:HG13	1:K:111:GLU:HB3	1.93	0.51
1:A:214:ARG:O	1:A:214:ARG:HG3	2.11	0.51
1:C:88:VAL:HG12	2:D:32:CYS:O	2.10	0.51
2:F:103:ILE:HG23	2:F:104:PRO:HD2	1.92	0.51
1:G:344:LEU:HA	1:G:348:LYS:HB2	1.92	0.51
1:I:312:ILE:HG23	1:I:340:LEU:HD22	1.92	0.51
1:A:184:GLN:HG3	9:A:907:HOH:O	2.11	0.51
1:K:82:ASP:HB2	1:K:86:PRO:HB3	1.93	0.51
2:B:133:ILE:HD13	2:B:354:LEU:HD13	1.91	0.51
1:C:97:ARG:HG2	1:C:101:ASP:OD2	2.11	0.51
2:D:39:ARG:HG3	2:D:40:TYR:CE1	2.46	0.51
1:E:106:VAL:HG13	1:E:111:GLU:HB3	1.93	0.51
1:E:93:SER:N	2:F:38:GLU:OE1	2.37	0.51
2:H:121:HIS:HB3	2:H:124:MET:HG2	1.93	0.51
9:C:927:HOH:O	1:K:339:GLU:HA	2.10	0.51
1:C:107:LEU:HD22	2:D:117:TYR:CD2	2.45	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:229:SER:O	2:B:233:MET:HG3	2.10	0.50
1:E:101:ASP:HA	1:E:104:ARG:HH11	1.77	0.50
1:E:265:GLU:O	1:E:269:LEU:HD13	2.11	0.50
2:B:22:ARG:HG2	2:B:22:ARG:HH11	1.76	0.50
2:F:352:GLU:O	2:F:355:ARG:HB3	2.11	0.50
2:B:353:ARG:HG2	2:B:353:ARG:NH1	2.24	0.50
1:I:339:GLU:O	1:I:343:ILE:HG13	2.12	0.50
2:L:79:LEU:O	2:L:96:ARG:HG3	2.11	0.50
1:A:103:PHE:CZ	1:A:133:VAL:HG22	2.47	0.50
1:A:88:VAL:HG12	2:B:32:CYS:O	2.11	0.50
2:B:207:ASP:O	2:B:208:ASN:HB2	2.12	0.50
1:G:318:ILE:HG22	1:G:322:MET:HE2	1.93	0.50
1:K:65:LEU:O	1:K:69:ARG:HG3	2.11	0.50
2:B:121:HIS:HB3	2:B:124:MET:HG2	1.94	0.50
1:K:311:LEU:HD23	1:K:311:LEU:O	2.12	0.50
2:D:338:CYS:HB3	9:D:830:HOH:O	2.10	0.50
1:C:105:ALA:O	1:C:109:ARG:HG3	2.11	0.50
1:K:339:GLU:O	1:K:343:ILE:HG13	2.11	0.50
2:L:236:LEU:HD22	2:L:245:LEU:HD21	1.94	0.50
2:B:303:ASP:OD1	2:B:306:VAL:HG13	2.12	0.49
1:A:104:ARG:NH2	9:A:917:HOH:O	2.45	0.49
1:A:106:VAL:HG13	1:A:111:GLU:HB3	1.94	0.49
2:D:295:ARG:CZ	2:D:299:LEU:HD11	2.42	0.49
1:E:173:ARG:NH1	9:E:921:HOH:O	2.41	0.49
2:F:29:PHE:O	2:F:33:LEU:HD22	2.12	0.49
1:G:105:ALA:O	1:G:109:ARG:HG3	2.12	0.49
1:G:69:ARG:HB3	1:G:71:GLU:OE1	2.12	0.49
1:I:350:THR:O	1:I:353:LYS:HB3	2.12	0.49
1:E:274:GLU:HG3	1:E:310:TYR:CE2	2.48	0.49
1:G:311:LEU:HD23	1:G:311:LEU:C	2.32	0.49
1:G:106:VAL:HG13	1:G:111:GLU:HB3	1.93	0.49
1:G:200:TYR:HB3	4:G:904:MES:H32	1.95	0.49
1:I:330:LYS:HE2	1:I:367:HIS:HB3	1.94	0.49
1:K:330:LYS:HE2	1:K:367:HIS:HB3	1.94	0.49
2:J:338:CYS:SG	2:J:349:ARG:NH2	2.85	0.49
1:E:103:PHE:CZ	1:E:133:VAL:HG22	2.47	0.49
1:E:198:LYS:HD3	2:F:266:LYS:HD3	1.94	0.49
1:G:340:LEU:HD23	1:G:343:ILE:HD12	1.94	0.49
1:C:148:LEU:CB	1:C:179:LEU:HD21	2.43	0.49
2:F:77:TYR:CE1	2:F:141:ARG:HB2	2.48	0.49
2:D:133:ILE:HG22	2:D:350:THR:HG23	1.93	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:138:ASP:OD1	2:F:140:SER:HB3	2.12	0.49
2:L:353:ARG:NH1	2:L:357:LEU:HG	2.28	0.49
1:A:344:LEU:HD13	1:A:356:TRP:CE2	2.48	0.49
1:E:100:TYR:O	1:E:104:ARG:HG3	2.12	0.49
2:F:99:SER:HB2	9:F:847:HOH:O	2.12	0.49
2:D:197:ILE:HD11	2:D:235:LYS:HD3	1.94	0.48
1:E:58:LEU:HD23	1:E:63:TYR:CE2	2.47	0.48
1:K:97:ARG:HG2	1:K:101:ASP:OD2	2.13	0.48
1:K:214:ARG:HH11	1:K:214:ARG:HG3	1.76	0.48
1:C:200:TYR:HB3	4:C:902:MES:H32	1.95	0.48
2:F:207:ASP:O	2:F:208:ASN:HB2	2.12	0.48
2:H:64:LEU:HD11	2:H:134:ILE:HG22	1.94	0.48
1:K:265:GLU:O	1:K:269:LEU:HD13	2.13	0.48
1:G:148:LEU:CB	1:G:179:LEU:HD21	2.43	0.48
2:H:22:ARG:HG2	2:H:22:ARG:HH11	1.77	0.48
1:I:148:LEU:CB	1:I:179:LEU:HD21	2.43	0.48
1:A:161:GLU:HG3	1:A:162:GLN:HG3	1.96	0.48
1:A:214:ARG:CG	1:A:214:ARG:O	2.61	0.48
1:A:285:GLN:NE2	2:B:247:ARG:NH1	2.61	0.48
1:A:334:LEU:O	1:A:338:LEU:HG	2.13	0.48
2:D:130:SER:O	2:D:134:ILE:HG13	2.12	0.48
2:D:64:LEU:HD11	2:D:134:ILE:HG22	1.94	0.48
1:E:353:LYS:CE	1:E:357:ARG:HH12	2.25	0.48
1:C:82:ASP:HB2	1:C:86:PRO:HB3	1.95	0.48
2:D:236:LEU:HD22	2:D:245:LEU:HD21	1.94	0.48
2:D:296:ASN:ND2	2:D:296:ASN:C	2.67	0.48
1:G:344:LEU:HD13	1:G:356:TRP:CE2	2.48	0.48
1:I:60:SER:OG	1:I:61:PRO:HD2	2.14	0.48
2:J:27:ARG:HH22	2:J:30:GLN:NE2	2.10	0.48
2:B:130:SER:O	2:B:134:ILE:HG13	2.12	0.48
2:B:258:ASN:CG	2:B:259:GLY:H	2.16	0.48
2:D:115:HIS:ND1	2:D:116:PRO:HD2	2.28	0.48
1:E:303:GLN:O	1:E:307:SER:HB2	2.13	0.48
1:E:82:ASP:HB2	1:E:86:PRO:HB3	1.95	0.48
2:F:208:ASN:ND2	2:F:247:ARG:HB3	2.29	0.48
2:H:269:ASP:OD2	2:H:311:LYS:HE3	2.14	0.48
2:L:92:ARG:NH1	2:L:118:ASP:O	2.47	0.48
1:E:207:GLN:HG2	1:E:242:PHE:CE2	2.49	0.48
2:H:186:ASN:HB2	2:H:358:HIS:NE2	2.28	0.48
2:L:103:ILE:HG23	2:L:104:PRO:HD2	1.94	0.48
1:A:117:PHE:CE2	1:A:146:LYS:HE2	2.49	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:189:ILE:HD11	1:A:205:HIS:CD2	2.49	0.48
1:I:265:GLU:HA	1:I:265:GLU:OE2	2.13	0.48
2:J:77:TYR:CZ	2:J:141:ARG:HB2	2.49	0.48
2:L:77:TYR:CZ	2:L:141:ARG:HB2	2.49	0.48
1:C:106:VAL:HG13	1:C:111:GLU:HB3	1.95	0.47
2:D:22:ARG:O	2:D:26:VAL:HG23	2.14	0.47
1:I:353:LYS:HG2	1:I:354:GLU:N	2.29	0.47
1:K:303:GLN:HB3	1:K:304:PRO:HD3	1.95	0.47
1:K:67:ARG:HD2	9:K:965:HOH:O	2.14	0.47
2:B:269:ASP:OD2	2:B:311:LYS:HE3	2.14	0.47
1:E:148:LEU:CB	1:E:179:LEU:HD21	2.44	0.47
1:E:353:LYS:CE	1:E:357:ARG:HH22	2.22	0.47
2:L:60:MET:HE2	2:L:347:SER:HB3	1.97	0.47
2:D:269:ASP:OD2	2:D:311:LYS:HE3	2.14	0.47
1:G:156:ILE:HD11	1:G:172:ARG:HH22	1.78	0.47
1:A:121:ARG:HD3	9:A:934:HOH:O	2.14	0.47
1:E:281:LYS:HD3	9:E:923:HOH:O	2.14	0.47
1:A:265:GLU:O	1:A:269:LEU:HD13	2.14	0.47
2:D:77:TYR:CZ	2:D:141:ARG:HB2	2.50	0.47
1:E:88:VAL:HG12	2:F:32:CYS:O	2.15	0.47
2:D:232:LEU:HD13	2:D:343:ALA:HB1	1.97	0.47
2:H:339:LYS:O	2:H:348:THR:HG23	2.14	0.47
1:K:82:ASP:OD2	2:L:99:SER:OG	2.33	0.47
1:C:198:LYS:HD3	2:D:266:LYS:HD3	1.96	0.47
1:E:301:ASP:O	1:E:304:PRO:HD2	2.15	0.47
1:I:232:ARG:HD3	2:J:265:ASN:ND2	2.28	0.47
2:B:64:LEU:HD11	2:B:134:ILE:HG22	1.96	0.47
2:B:37:PRO:HB2	2:B:39:ARG:HG2	1.97	0.47
2:H:77:TYR:CZ	2:H:141:ARG:HB2	2.50	0.47
1:G:79:PRO:HA	1:G:101:ASP:OD1	2.15	0.47
2:F:64:LEU:HD23	2:F:64:LEU:HA	1.78	0.46
1:K:189:ILE:HD11	1:K:205:HIS:CD2	2.48	0.46
2:B:236:LEU:HD22	2:B:245:LEU:HD21	1.97	0.46
2:B:357:LEU:HD22	2:B:361:TRP:CZ2	2.50	0.46
2:D:312:TRP:HB3	2:D:313:PRO:HD2	1.97	0.46
2:F:250:ARG:O	2:F:254:MET:HG2	2.14	0.46
2:J:267:PRO:HG2	9:J:836:HOH:O	2.15	0.46
2:J:30:GLN:HB3	2:J:30:GLN:HE21	1.62	0.46
1:K:200:TYR:HB3	4:K:906:MES:H32	1.97	0.46
2:F:77:TYR:CZ	2:F:141:ARG:HB2	2.51	0.46
2:H:22:ARG:O	2:H:26:VAL:HG23	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:151:GLU:HG3	1:A:175:LEU:HD11	1.98	0.46
2:D:86:ASP:N	2:D:86:ASP:OD2	2.45	0.46
2:F:236:LEU:HD22	2:F:245:LEU:HD21	1.98	0.46
1:E:91:ILE:HD11	2:F:38:GLU:H	1.81	0.46
2:H:103:ILE:HG23	2:H:104:PRO:HD2	1.97	0.46
1:E:121:ARG:HH11	1:E:121:ARG:CG	2.29	0.46
1:E:296:LEU:HD22	1:E:322:MET:CE	2.46	0.46
2:F:22:ARG:O	2:F:26:VAL:HG23	2.16	0.46
2:J:333:GLU:OE1	2:J:339:LYS:HE3	2.15	0.46
2:J:37:PRO:HD2	2:J:40:TYR:CE1	2.51	0.46
1:C:189:ILE:HD11	1:C:205:HIS:CD2	2.46	0.46
2:J:152:ARG:HD3	2:J:189:SER:O	2.15	0.46
1:K:214:ARG:NH1	1:K:214:ARG:HG3	2.31	0.46
1:A:91:ILE:HD11	2:B:38:GLU:H	1.81	0.46
1:E:200:TYR:HB3	4:E:903:MES:H32	1.98	0.46
1:K:83:GLY:HA3	2:L:105:PHE:CD1	2.51	0.46
2:B:138:ASP:OD1	2:B:140:SER:HB3	2.16	0.45
2:F:192:ASP:CG	2:F:195:LYS:HG3	2.36	0.45
4:E:903:MES:H51	2:F:219:HIS:NE2	2.30	0.45
1:I:256:LEU:HD22	1:I:287:ARG:NH2	2.32	0.45
2:J:207:ASP:O	2:J:208:ASN:HB2	2.16	0.45
2:J:59:ASP:OD2	2:J:349:ARG:NH1	2.50	0.45
1:A:355:TYR:O	1:A:358:TYR:HB3	2.17	0.45
1:E:296:LEU:HD22	1:E:322:MET:HE3	1.97	0.45
2:H:18:LEU:N	2:H:18:LEU:CD2	2.80	0.45
2:H:20:PHE:CE2	2:H:337:ILE:HD11	2.52	0.45
2:J:245:LEU:O	2:J:249:LYS:HG3	2.17	0.45
2:B:92:ARG:NH1	2:B:118:ASP:O	2.50	0.45
1:C:340:LEU:HD23	1:C:343:ILE:HD12	1.98	0.45
2:D:354:LEU:HD11	2:D:358:HIS:HE2	1.80	0.45
1:I:344:LEU:HD13	1:I:356:TRP:CE2	2.52	0.45
2:J:197:ILE:HD11	2:J:235:LYS:HD3	1.99	0.45
1:G:219:GLU:OE1	1:G:219:GLU:HA	2.17	0.45
2:H:22:ARG:NH1	2:H:22:ARG:HG2	2.31	0.45
1:I:189:ILE:HD11	1:I:205:HIS:CD2	2.50	0.45
1:C:96:PHE:CE1	1:C:126:LEU:HB3	2.52	0.45
2:D:236:LEU:CD2	2:D:245:LEU:HD21	2.47	0.45
2:H:37:PRO:HD2	2:H:40:TYR:CD1	2.52	0.45
2:B:22:ARG:NH1	2:B:22:ARG:HG2	2.31	0.45
1:C:344:LEU:HD13	1:C:356:TRP:CE2	2.52	0.45
1:E:318:ILE:HG22	1:E:322:MET:CE	2.47	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:154:LEU:HD11	9:F:827:HOH:O	2.16	0.45
1:I:346:LYS:O	1:I:347:GLU:HG2	2.16	0.45
2:J:19:ASP:OD2	2:J:19:ASP:N	2.48	0.45
2:B:336:GLY:HA2	2:J:305:LEU:CD1	2.47	0.45
2:B:250:ARG:O	2:B:254:MET:HG2	2.17	0.45
2:H:30:GLN:O	2:H:34:GLN:HG3	2.17	0.45
1:I:355:TYR:O	1:I:358:TYR:HB3	2.16	0.45
2:L:22:ARG:HH11	2:L:22:ARG:HG2	1.81	0.45
2:B:21:LEU:CD1	2:B:21:LEU:N	2.80	0.44
1:C:103:PHE:CZ	1:C:133:VAL:HG22	2.52	0.44
1:C:265:GLU:O	1:C:269:LEU:HD13	2.17	0.44
1:G:265:GLU:O	1:G:269:LEU:HD13	2.17	0.44
2:D:92:ARG:NH1	2:D:118:ASP:O	2.50	0.44
2:J:229:SER:O	2:J:233:MET:HG3	2.18	0.44
2:J:236:LEU:HD22	2:J:245:LEU:HD21	1.99	0.44
2:D:245:LEU:O	2:D:249:LYS:HG3	2.17	0.44
2:D:250:ARG:O	2:D:254:MET:HG2	2.18	0.44
1:C:91:ILE:HD11	2:D:38:GLU:H	1.82	0.44
2:F:245:LEU:O	2:F:249:LYS:HG3	2.16	0.44
1:A:121:ARG:NH1	1:A:121:ARG:HG3	2.32	0.44
2:B:59:ASP:OD2	2:B:349:ARG:NH1	2.49	0.44
2:D:92:ARG:HB3	2:D:119:SER:CB	2.48	0.44
2:H:29:PHE:O	2:H:33:LEU:HD22	2.18	0.44
1:K:325:ASN:O	1:K:326:GLN:C	2.55	0.44
2:L:130:SER:O	2:L:134:ILE:HG13	2.17	0.44
2:D:193:MET:HG3	2:D:233:MET:CE	2.47	0.44
1:E:214:ARG:O	1:E:214:ARG:CG	2.64	0.44
2:H:21:LEU:CD1	2:H:21:LEU:N	2.80	0.44
2:H:236:LEU:HD22	2:H:245:LEU:HD21	1.98	0.44
2:H:357:LEU:CD2	2:H:361:TRP:CZ2	3.01	0.44
1:I:147:ASP:OD1	1:I:150:GLU:N	2.48	0.44
1:I:173:ARG:O	1:I:177:GLU:HG3	2.16	0.44
2:B:103:ILE:HG23	2:B:104:PRO:HD2	1.98	0.44
1:C:353:LYS:CE	1:C:357:ARG:NH1	2.78	0.44
2:D:357:LEU:HD22	2:D:361:TRP:CZ2	2.53	0.44
1:E:156:ILE:HD11	1:E:172:ARG:HH22	1.83	0.44
1:I:96:PHE:CE1	1:I:126:LEU:HB3	2.53	0.44
1:I:151:GLU:HG3	1:I:175:LEU:HD11	1.98	0.44
1:I:303:GLN:O	1:I:307:SER:HB2	2.18	0.44
1:K:65:LEU:HD12	1:K:67:ARG:NH1	2.33	0.44
1:G:318:ILE:HG22	1:G:322:MET:CE	2.47	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:97:ARG:NH1	1:G:97:ARG:HB3	2.32	0.44
1:K:79:PRO:HA	1:K:101:ASP:OD1	2.18	0.44
1:C:219:GLU:OE1	1:C:219:GLU:HA	2.18	0.44
2:H:348:THR:HA	2:H:351:SER:OG	2.17	0.44
1:K:96:PHE:CE1	1:K:126:LEU:HB3	2.53	0.44
2:L:250:ARG:O	2:L:254:MET:HG2	2.18	0.44
1:C:195:GLN:H	1:C:195:GLN:HG2	1.58	0.43
1:E:357:ARG:HH11	1:E:357:ARG:HG3	1.83	0.43
1:I:103:PHE:CZ	1:I:133:VAL:HG22	2.53	0.43
1:I:265:GLU:O	1:I:269:LEU:HD13	2.19	0.43
4:A:901:MES:H51	2:B:219:HIS:NE2	2.34	0.43
2:B:77:TYR:CZ	2:B:141:ARG:HB2	2.53	0.43
1:C:65:LEU:O	1:C:69:ARG:HG3	2.18	0.43
1:G:267:ILE:HD13	1:G:277:TRP:CE2	2.53	0.43
1:I:198:LYS:HD3	2:J:266:LYS:HD3	1.99	0.43
1:C:253:ARG:HD3	1:I:56:LEU:HD22	2.00	0.43
1:C:287:ARG:H	1:C:287:ARG:HG2	1.62	0.43
2:D:77:TYR:CE1	2:D:141:ARG:HB2	2.53	0.43
1:E:96:PHE:CE1	1:E:126:LEU:HB3	2.53	0.43
1:I:287:ARG:H	1:I:287:ARG:HG2	1.56	0.43
1:E:294:ASN:O	1:E:298:GLN:HG3	2.19	0.43
1:G:286:ASP:HB2	9:G:941:HOH:O	2.17	0.43
1:I:325:ASN:O	1:I:326:GLN:C	2.57	0.43
1:A:311:LEU:HD23	1:A:311:LEU:C	2.39	0.43
2:J:77:TYR:HE2	2:J:137:ASP:OD2	2.01	0.43
2:F:92:ARG:HB3	2:F:119:SER:CB	2.48	0.43
1:G:189:ILE:HD11	1:G:205:HIS:CD2	2.47	0.43
1:K:149:GLN:HE22	1:K:179:LEU:HD13	1.84	0.43
1:K:261:GLN:O	1:K:265:GLU:HG2	2.18	0.43
2:F:36:LEU:HA	2:F:37:PRO:HD3	1.86	0.43
2:H:24:ARG:NE	2:H:304:ARG:O	2.52	0.43
2:H:33:LEU:CD2	2:H:54:ALA:HB1	2.49	0.43
2:L:77:TYR:CE1	2:L:141:ARG:HB2	2.54	0.43
3:N:206:THR:HB	3:N:207:LYS:H	1.68	0.43
2:B:267:PRO:HG2	9:B:843:HOH:O	2.18	0.43
2:D:354:LEU:HD11	2:D:358:HIS:NE2	2.33	0.43
2:H:19:ASP:O	2:H:21:LEU:HD13	2.19	0.43
2:H:258:ASN:CG	2:H:259:GLY:H	2.21	0.43
1:A:212:GLU:O	1:G:180:LYS:HE3	2.19	0.42
2:H:344:LEU:HB3	2:H:346:VAL:HG22	2.00	0.42
1:A:200:TYR:HB3	4:A:901:MES:H32	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:72:TRP:CZ2	1:I:115:ARG:HB2	2.53	0.42
2:J:115:HIS:HA	2:J:116:PRO:HD3	1.91	0.42
2:J:29:PHE:O	2:J:33:LEU:HD22	2.18	0.42
1:K:219:GLU:OE1	1:K:219:GLU:HA	2.19	0.42
1:K:78:VAL:O	1:K:104:ARG:HD2	2.18	0.42
2:L:229:SER:O	2:L:233:MET:HG3	2.19	0.42
1:E:318:ILE:HG22	1:E:322:MET:HE3	2.02	0.42
1:G:91:ILE:HD11	2:H:38:GLU:H	1.83	0.42
2:H:90:LEU:HD23	2:H:90:LEU:HA	1.91	0.42
1:C:136:PHE:CE2	1:C:140:LEU:HD11	2.54	0.42
1:G:323:LEU:HB3	1:G:367:HIS:CD2	2.54	0.42
2:H:33:LEU:HD22	2:H:54:ALA:HB1	2.00	0.42
1:I:149:GLN:NE2	1:I:179:LEU:HD13	2.34	0.42
2:F:122:ILE:HG22	2:F:163:ALA:HA	2.01	0.42
2:J:144:LYS:HG2	2:J:185:LEU:HD22	2.01	0.42
2:L:258:ASN:OD1	2:L:259:GLY:N	2.45	0.42
1:A:156:ILE:HD11	1:A:172:ARG:HH22	1.84	0.42
1:A:267:ILE:HD13	1:A:277:TRP:CE2	2.55	0.42
2:B:39:ARG:HG3	2:B:40:TYR:CE1	2.54	0.42
2:D:68:ASN:OD1	2:D:70:ASP:HB2	2.19	0.42
2:D:92:ARG:HH11	2:D:119:SER:HB3	1.85	0.42
1:G:189:ILE:HG21	1:G:206:ARG:HB2	2.02	0.42
1:I:106:VAL:HG11	1:I:116:ALA:CB	2.50	0.42
2:L:173:ARG:HD2	8:R:1608:GER:H142	2.02	0.42
1:C:244:ILE:HD11	1:C:259:GLU:OE1	2.20	0.42
2:D:37:PRO:O	2:D:39:ARG:N	2.53	0.42
1:E:189:ILE:HD11	1:E:205:HIS:CD2	2.47	0.42
1:E:287:ARG:H	1:E:287:ARG:HG2	1.48	0.42
2:H:249:LYS:HB3	2:H:285:ILE:HD13	2.01	0.42
2:J:329:LEU:HD12	2:J:329:LEU:HA	1.90	0.42
1:A:96:PHE:CE1	1:A:126:LEU:HB3	2.55	0.42
1:A:325:ASN:O	1:A:326:GLN:C	2.58	0.42
1:G:96:PHE:CE1	1:G:126:LEU:HB3	2.54	0.42
2:H:256:GLN:HB2	2:H:260:TYR:CE2	2.55	0.42
2:B:86:ASP:N	2:B:86:ASP:OD2	2.47	0.42
1:E:97:ARG:HG2	1:E:101:ASP:OD2	2.20	0.42
1:G:289:LEU:HG	9:G:936:HOH:O	2.19	0.42
2:H:86:ASP:N	2:H:86:ASP:OD2	2.48	0.42
2:B:64:LEU:HD23	2:B:64:LEU:HA	1.81	0.42
2:H:207:ASP:O	2:H:208:ASN:HB2	2.20	0.42
2:H:250:ARG:O	2:H:254:MET:HG2	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:40:TYR:CD1	2:H:40:TYR:N	2.88	0.42
3:M:107:LYS:HE3	3:M:109:VAL:HG22	2.02	0.42
3:N:207:LYS:HE3	3:N:209:VAL:HG22	2.02	0.42
3:R:607:LYS:HE3	3:R:609:VAL:HG22	2.01	0.42
1:A:348:LYS:HD3	1:A:348:LYS:HA	1.79	0.41
1:A:71:GLU:CD	1:A:71:GLU:H	2.15	0.41
2:B:160:SER:HB3	2:B:199:TYR:CE1	2.56	0.41
2:F:258:ASN:CG	2:F:259:GLY:H	2.22	0.41
2:H:37:PRO:HD2	2:H:40:TYR:CE1	2.55	0.41
1:I:82:ASP:HB2	1:I:86:PRO:HB3	2.01	0.41
1:K:353:LYS:HB2	1:K:353:LYS:HE3	1.70	0.41
2:L:92:ARG:HB3	2:L:119:SER:CB	2.49	0.41
1:A:312:ILE:O	1:A:316:VAL:HG23	2.20	0.41
2:B:256:GLN:HB2	2:B:260:TYR:CE2	2.55	0.41
2:D:255:ARG:CZ	2:D:264:PRO:HG3	2.50	0.41
2:H:186:ASN:HB2	2:H:358:HIS:CE1	2.55	0.41
2:J:36:LEU:HA	2:J:37:PRO:HD3	1.91	0.41
1:A:105:ALA:O	1:A:109:ARG:HG3	2.21	0.41
1:C:87:VAL:CG1	2:D:33:LEU:HD12	2.51	0.41
2:J:122:ILE:HG22	2:J:163:ALA:HA	2.01	0.41
1:K:142:ARG:HA	1:K:142:ARG:HD3	1.83	0.41
1:K:361:ARG:HG3	9:K:947:HOH:O	2.20	0.41
1:A:189:ILE:HG21	1:A:206:ARG:HB2	2.03	0.41
1:A:66:TYR:CE1	1:A:119:LEU:HD13	2.55	0.41
1:C:156:ILE:HD11	1:C:172:ARG:HH22	1.85	0.41
2:F:37:PRO:C	2:F:39:ARG:H	2.24	0.41
2:B:186:ASN:HB2	2:B:358:HIS:NE2	2.34	0.41
2:D:92:ARG:NH1	2:D:119:SER:HB3	2.36	0.41
2:F:37:PRO:C	2:F:39:ARG:N	2.73	0.41
1:K:101:ASP:HA	1:K:104:ARG:HH11	1.86	0.41
1:A:148:LEU:CB	1:A:179:LEU:HD21	2.47	0.41
1:A:308:SER:O	1:A:312:ILE:HG12	2.21	0.41
2:B:115:HIS:HA	2:B:116:PRO:HD3	1.87	0.41
2:D:19:ASP:O	2:D:21:LEU:HD13	2.20	0.41
2:H:37:PRO:C	2:H:39:ARG:N	2.74	0.41
1:I:106:VAL:HG11	1:I:116:ALA:HB1	2.03	0.41
1:K:225:GLN:NE2	1:K:229:GLU:OE2	2.52	0.41
2:L:22:ARG:HG2	2:L:22:ARG:NH1	2.35	0.41
4:A:901:MES:H82	4:A:901:MES:H31	1.92	0.41
1:C:106:VAL:HG11	1:C:116:ALA:HB1	2.03	0.41
1:E:223:VAL:HG11	1:E:240:ARG:HB2	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:235:LYS:O	2:D:239:VAL:HG23	2.21	0.41
1:G:287:ARG:H	1:G:287:ARG:HG2	1.56	0.41
1:G:355:TYR:O	1:G:358:TYR:HB3	2.21	0.41
3:Q:507:LYS:HE3	3:Q:509:VAL:HG22	2.03	0.41
1:A:192:ILE:O	1:A:195:GLN:HG3	2.21	0.41
1:A:219:GLU:HA	1:A:219:GLU:OE1	2.21	0.41
1:G:106:VAL:HG11	1:G:116:ALA:HB1	2.03	0.41
4:G:904:MES:H31	4:G:904:MES:H82	1.94	0.41
1:I:285:GLN:NE2	2:J:247:ARG:NH1	2.68	0.41
1:K:148:LEU:CB	1:K:179:LEU:HD21	2.49	0.41
2:L:236:LEU:CD2	2:L:245:LEU:HD21	2.51	0.41
1:C:189:ILE:HG21	1:C:206:ARG:HB2	2.03	0.41
1:C:251:SER:HA	1:C:287:ARG:HH12	1.86	0.41
2:H:230:LEU:HD22	2:H:239:VAL:HG21	2.03	0.41
1:I:101:ASP:HA	1:I:104:ARG:HH11	1.86	0.41
1:K:91:ILE:HG13	2:L:36:LEU:O	2.21	0.41
2:L:197:ILE:HD11	2:L:235:LYS:HD3	2.02	0.41
2:L:64:LEU:HD23	2:L:64:LEU:HA	1.86	0.41
2:B:232:LEU:HD13	2:B:343:ALA:HB1	2.03	0.40
2:B:36:LEU:HA	2:B:37:PRO:HD3	1.83	0.40
1:E:136:PHE:CE2	1:E:140:LEU:HD11	2.56	0.40
2:F:115:HIS:HA	2:F:116:PRO:HD3	1.92	0.40
1:G:301:ASP:O	1:G:304:PRO:HD2	2.20	0.40
2:J:64:LEU:HA	2:J:64:LEU:HD23	1.88	0.40
1:K:117:PHE:CE2	1:K:146:LYS:HE2	2.56	0.40
2:D:202:ARG:HH11	2:D:202:ARG:CG	2.34	0.40
1:E:267:ILE:HD13	1:E:277:TRP:CE2	2.56	0.40
2:L:133:ILE:CD1	2:L:354:LEU:HD13	2.51	0.40
3:O:307:LYS:HE3	3:O:309:VAL:HG22	2.02	0.40
1:A:274:GLU:HG2	1:A:278:ASN:ND2	2.36	0.40
1:C:72:TRP:CZ2	1:C:115:ARG:HB2	2.55	0.40
2:D:348:THR:HA	2:D:351:SER:OG	2.21	0.40
2:D:79:LEU:O	2:D:96:ARG:HG3	2.21	0.40
2:F:269:ASP:OD2	2:F:311:LYS:HE3	2.21	0.40
2:B:236:LEU:CD2	2:B:245:LEU:HD21	2.51	0.40
1:G:325:ASN:O	1:G:326:GLN:C	2.59	0.40
1:K:192:ILE:O	1:K:195:GLN:HG3	2.21	0.40
2:L:306:VAL:HG23	2:L:306:VAL:O	2.20	0.40
1:A:156:ILE:CD1	1:A:172:ARG:HH22	2.35	0.40
2:D:292:GLU:HG3	9:K:989:HOH:O	2.22	0.40
2:F:19:ASP:OD2	2:F:304:ARG:NH2	2.54	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:287:ARG:HG2	1:K:287:ARG:H	1.60	0.40
1:K:60:SER:OG	1:K:61:PRO:HD2	2.21	0.40
2:L:138:ASP:OD1	2:L:140:SER:HB3	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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	312/377 (83%)	288 (92%)	23 (7%)	1 (0%)	41	72
1	C	312/377 (83%)	292 (94%)	20 (6%)	0	100	100
1	E	312/377 (83%)	291 (93%)	21 (7%)	0	100	100
1	G	312/377 (83%)	287 (92%)	24 (8%)	1 (0%)	41	72
1	I	312/377 (83%)	290 (93%)	22 (7%)	0	100	100
1	K	312/377 (83%)	292 (94%)	19 (6%)	1 (0%)	41	72
2	B	344/377 (91%)	324 (94%)	18 (5%)	2 (1%)	25	56
2	D	344/377 (91%)	326 (95%)	14 (4%)	4 (1%)	13	39
2	F	344/377 (91%)	323 (94%)	19 (6%)	2 (1%)	25	56
2	H	344/377 (91%)	318 (92%)	22 (6%)	4 (1%)	13	39
2	J	344/377 (91%)	324 (94%)	17 (5%)	3 (1%)	17	46
2	L	344/377 (91%)	325 (94%)	17 (5%)	2 (1%)	25	56
3	M	4/11 (36%)	4 (100%)	0	0	100	100
3	N	4/11 (36%)	4 (100%)	0	0	100	100
3	O	4/11 (36%)	4 (100%)	0	0	100	100
3	P	4/11 (36%)	4 (100%)	0	0	100	100
3	Q	4/11 (36%)	4 (100%)	0	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	R	4/11 (36%)	4 (100%)	0	0	100	100
All	All	3960/4590 (86%)	3704 (94%)	236 (6%)	20 (0%)	29	61

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	111	PRO
2	B	258	ASN
2	D	111	PRO
2	D	258	ASN
2	F	111	PRO
2	F	258	ASN
2	H	111	PRO
2	H	258	ASN
2	J	111	PRO
2	J	258	ASN
1	K	306	HIS
2	L	111	PRO
2	L	258	ASN
2	D	38	GLU
2	H	333	GLU
2	J	34	GLN
1	A	306	HIS
2	D	333	GLU
1	G	306	HIS
2	H	66	VAL

5.3.2 Protein sidechains [i](#)

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	276/338 (82%)	270 (98%)	6 (2%)	52	83
1	C	285/338 (84%)	275 (96%)	10 (4%)	36	70
1	E	280/338 (83%)	269 (96%)	11 (4%)	32	66

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	G	281/338 (83%)	276 (98%)	5 (2%)	59	86
1	I	287/338 (85%)	278 (97%)	9 (3%)	40	74
1	K	290/338 (86%)	279 (96%)	11 (4%)	33	67
2	B	286/326 (88%)	273 (96%)	13 (4%)	27	60
2	D	287/326 (88%)	270 (94%)	17 (6%)	19	49
2	F	291/326 (89%)	276 (95%)	15 (5%)	23	55
2	H	282/326 (86%)	271 (96%)	11 (4%)	32	66
2	J	291/326 (89%)	274 (94%)	17 (6%)	20	50
2	L	292/326 (90%)	279 (96%)	13 (4%)	27	60
3	M	6/11 (54%)	6 (100%)	0	100	100
3	N	6/11 (54%)	6 (100%)	0	100	100
3	O	6/11 (54%)	6 (100%)	0	100	100
3	P	6/11 (54%)	6 (100%)	0	100	100
3	Q	6/11 (54%)	6 (100%)	0	100	100
3	R	6/11 (54%)	6 (100%)	0	100	100
All	All	3464/4050 (86%)	3326 (96%)	138 (4%)	31	65

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

Mol	Chain	Res	Type
1	A	55	PHE
1	A	71	GLU
1	A	184	GLN
1	A	195	GLN
1	A	214	ARG
1	A	287	ARG
2	B	21	LEU
2	B	33	LEU
2	B	39	ARG
2	B	151	LEU
2	B	216	LEU
2	B	232	LEU
2	B	236	LEU
2	B	255	ARG
2	B	261	HIS
2	B	305	LEU
2	B	329	LEU

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Mol	Chain	Res	Type
2	B	331	LEU
2	B	357	LEU
1	C	59	ASP
1	C	71	GLU
1	C	107	LEU
1	C	184	GLN
1	C	195	GLN
1	C	225	GLN
1	C	287	ARG
1	C	301	ASP
1	C	324	GLU
1	C	364	GLN
2	D	21	LEU
2	D	39	ARG
2	D	65	ASP
2	D	151	LEU
2	D	216	LEU
2	D	232	LEU
2	D	236	LEU
2	D	255	ARG
2	D	258	ASN
2	D	261	HIS
2	D	296	ASN
2	D	305	LEU
2	D	306	VAL
2	D	329	LEU
2	D	331	LEU
2	D	353	ARG
2	D	357	LEU
1	E	55	PHE
1	E	71	GLU
1	E	81	ASN
1	E	121	ARG
1	E	184	GLN
1	E	194	ASN
1	E	214	ARG
1	E	287	ARG
1	E	320	GLU
1	E	324	GLU
1	E	364	GLN
2	F	21	LEU
2	F	24	ARG

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Mol	Chain	Res	Type
2	F	151	LEU
2	F	202	ARG
2	F	216	LEU
2	F	232	LEU
2	F	236	LEU
2	F	255	ARG
2	F	258	ASN
2	F	261	HIS
2	F	305	LEU
2	F	329	LEU
2	F	331	LEU
2	F	353	ARG
2	F	357	LEU
1	G	71	GLU
1	G	107	LEU
1	G	184	GLN
1	G	287	ARG
1	G	324	GLU
2	H	21	LEU
2	H	151	LEU
2	H	216	LEU
2	H	232	LEU
2	H	236	LEU
2	H	255	ARG
2	H	261	HIS
2	H	305	LEU
2	H	306	VAL
2	H	329	LEU
2	H	331	LEU
1	I	55	PHE
1	I	67	ARG
1	I	71	GLU
1	I	107	LEU
1	I	184	GLN
1	I	194	ASN
1	I	287	ARG
1	I	324	GLU
1	I	353	LYS
2	J	21	LEU
2	J	27	ARG
2	J	30	GLN
2	J	31	ARG

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Mol	Chain	Res	Type
2	J	92	ARG
2	J	151	LEU
2	J	216	LEU
2	J	232	LEU
2	J	236	LEU
2	J	255	ARG
2	J	261	HIS
2	J	305	LEU
2	J	329	LEU
2	J	331	LEU
2	J	341	HIS
2	J	353	ARG
2	J	357	LEU
1	K	55	PHE
1	K	59	ASP
1	K	71	GLU
1	K	107	LEU
1	K	121	ARG
1	K	142	ARG
1	K	194	ASN
1	K	195	GLN
1	K	287	ARG
1	K	298	GLN
1	K	324	GLU
2	L	21	LEU
2	L	65	ASP
2	L	92	ARG
2	L	151	LEU
2	L	216	LEU
2	L	232	LEU
2	L	236	LEU
2	L	255	ARG
2	L	261	HIS
2	L	305	LEU
2	L	329	LEU
2	L	331	LEU
2	L	353	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (53) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	184	GLN

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Mol	Chain	Res	Type
1	A	278	ASN
1	A	285	GLN
1	A	367	HIS
2	B	208	ASN
2	B	265	ASN
1	C	81	ASN
1	C	108	GLN
1	C	170	HIS
1	C	184	GLN
1	C	195	GLN
1	C	201	HIS
1	C	225	GLN
1	C	298	GLN
1	C	364	GLN
1	C	367	HIS
2	D	246	ASN
2	D	265	ASN
2	D	296	ASN
1	E	81	ASN
1	E	89	GLN
1	E	184	GLN
1	E	367	HIS
2	F	30	GLN
2	F	246	ASN
2	F	265	ASN
1	G	80	GLN
1	G	81	ASN
1	G	89	GLN
1	G	162	GLN
1	G	184	GLN
1	G	297	ASN
2	H	246	ASN
2	H	265	ASN
2	H	302	GLN
1	I	81	ASN
1	I	89	GLN
1	I	149	GLN
1	I	184	GLN
1	I	195	GLN
1	I	285	GLN
1	I	364	GLN
2	J	30	GLN

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Mol	Chain	Res	Type
2	J	34	GLN
2	J	208	ASN
2	J	246	ASN
2	J	265	ASN
1	K	149	GLN
1	K	184	GLN
1	K	297	ASN
1	K	298	GLN
2	L	246	ASN
2	L	265	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 31 ligands modelled in this entry, 13 are monoatomic - leaving 18 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
6	SO4	H	813	-	4,4,4	0.20	0	6,6,6	0.17	0
4	MES	E	903	-	12,12,12	8.84	8 (66%)	14,16,16	2.76	5 (35%)
6	SO4	B	810	-	4,4,4	0.22	0	6,6,6	0.16	0
8	GER	M	1108	3	19,19,19	0.69	0	22,22,22	0.69	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	MES	K	906	-	12,12,12	8.86	8 (66%)	14,16,16	2.85	6 (42%)
4	MES	A	901	-	12,12,12	8.95	8 (66%)	14,16,16	2.74	5 (35%)
8	GER	P	1408	3	19,19,19	0.59	0	22,22,22	0.70	0
4	MES	I	905	-	12,12,12	8.74	8 (66%)	14,16,16	2.88	6 (42%)
8	GER	O	1308	3	19,19,19	0.66	0	22,22,22	0.71	0
4	MES	G	904	-	12,12,12	8.81	8 (66%)	14,16,16	2.77	5 (35%)
6	SO4	L	815	-	4,4,4	0.17	0	6,6,6	0.25	0
8	GER	R	1608	3	19,19,19	0.64	0	22,22,22	0.71	0
8	GER	N	1208	3	19,19,19	0.64	0	22,22,22	0.69	0
6	SO4	D	811	-	4,4,4	0.22	0	6,6,6	0.18	0
6	SO4	J	814	-	4,4,4	0.21	0	6,6,6	0.21	0
6	SO4	F	812	-	4,4,4	0.26	0	6,6,6	0.19	0
8	GER	Q	1508	3	19,19,19	0.65	0	22,22,22	0.72	0
4	MES	C	902	-	12,12,12	8.81	8 (66%)	14,16,16	2.73	5 (35%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	MES	E	903	-	-	3/6/14/14	0/1/1/1
8	GER	M	1108	3	-	2/20/20/20	-
8	GER	O	1308	3	-	2/20/20/20	-
4	MES	K	906	-	-	3/6/14/14	0/1/1/1
4	MES	A	901	-	-	3/6/14/14	0/1/1/1
8	GER	P	1408	3	-	2/20/20/20	-
4	MES	I	905	-	-	3/6/14/14	0/1/1/1
4	MES	G	904	-	-	3/6/14/14	0/1/1/1
8	GER	R	1608	3	-	2/20/20/20	-
8	GER	N	1208	3	-	2/20/20/20	-
8	GER	Q	1508	3	-	2/20/20/20	-
4	MES	C	902	-	-	3/6/14/14	0/1/1/1

All (48) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	K	906	MES	C8-S	-23.49	1.44	1.77

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	901	MES	C8-S	-23.47	1.44	1.77
4	G	904	MES	C8-S	-23.00	1.44	1.77
4	E	903	MES	C8-S	-22.96	1.44	1.77
4	C	902	MES	C8-S	-22.91	1.44	1.77
4	I	905	MES	C8-S	-22.79	1.45	1.77
4	E	903	MES	O2S-S	12.07	1.80	1.45
4	C	902	MES	O2S-S	11.96	1.80	1.45
4	A	901	MES	O2S-S	11.95	1.80	1.45
4	C	902	MES	O1S-S	11.84	1.80	1.45
4	E	903	MES	O1S-S	11.82	1.79	1.45
4	G	904	MES	O2S-S	11.81	1.79	1.45
4	I	905	MES	O2S-S	11.77	1.79	1.45
4	G	904	MES	O1S-S	11.72	1.79	1.45
4	A	901	MES	O1S-S	11.72	1.79	1.45
4	I	905	MES	O1S-S	11.64	1.79	1.45
4	K	906	MES	O2S-S	11.61	1.79	1.45
4	K	906	MES	O1S-S	11.49	1.79	1.45
4	A	901	MES	O3S-S	8.97	1.79	1.47
4	G	904	MES	O3S-S	8.88	1.78	1.47
4	E	903	MES	O3S-S	8.87	1.78	1.47
4	C	902	MES	O3S-S	8.78	1.78	1.47
4	I	905	MES	O3S-S	8.76	1.78	1.47
4	K	906	MES	O3S-S	8.60	1.78	1.47
4	K	906	MES	C7-C8	-4.89	1.39	1.52
4	A	901	MES	C7-C8	-4.79	1.39	1.52
4	G	904	MES	C7-C8	-4.78	1.39	1.52
4	I	905	MES	C7-C8	-4.74	1.39	1.52
4	E	903	MES	C7-C8	-4.70	1.39	1.52
4	C	902	MES	C7-C8	-4.63	1.40	1.52
4	A	901	MES	C3-C2	-2.76	1.39	1.50
4	K	906	MES	C3-C2	-2.73	1.39	1.50
4	C	902	MES	C3-C2	-2.71	1.39	1.50
4	G	904	MES	C3-C2	-2.71	1.39	1.50
4	E	903	MES	C3-C2	-2.67	1.40	1.50
4	I	905	MES	C3-C2	-2.62	1.40	1.50
4	A	901	MES	C5-C6	-2.57	1.40	1.50
4	I	905	MES	C5-C6	-2.57	1.40	1.50
4	E	903	MES	C5-C6	-2.56	1.40	1.50
4	C	902	MES	C5-C6	-2.54	1.40	1.50
4	K	906	MES	C5-C6	-2.54	1.40	1.50
4	G	904	MES	C5-C6	-2.54	1.40	1.50
4	A	901	MES	C7-N4	-2.31	1.42	1.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	G	904	MES	C7-N4	-2.21	1.42	1.47
4	C	902	MES	C7-N4	-2.18	1.42	1.47
4	K	906	MES	C7-N4	-2.17	1.42	1.47
4	E	903	MES	C7-N4	-2.14	1.42	1.47
4	I	905	MES	C7-N4	-2.08	1.42	1.47

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	G	904	MES	O3S-S-C8	5.85	115.23	105.77
4	I	905	MES	O3S-S-C8	5.76	115.08	105.77
4	E	903	MES	O3S-S-C8	5.72	115.02	105.77
4	A	901	MES	O3S-S-C8	5.67	114.94	105.77
4	K	906	MES	O3S-S-C8	5.64	114.89	105.77
4	C	902	MES	O3S-S-C8	5.42	114.53	105.77
4	A	901	MES	O1S-S-C8	5.27	113.26	106.92
4	I	905	MES	O1S-S-C8	5.21	113.19	106.92
4	K	906	MES	O2S-S-C8	5.08	113.03	106.92
4	C	902	MES	O1S-S-C8	5.01	112.94	106.92
4	E	903	MES	O1S-S-C8	5.00	112.94	106.92
4	K	906	MES	O1S-S-C8	4.96	112.88	106.92
4	G	904	MES	O1S-S-C8	4.82	112.72	106.92
4	I	905	MES	O2S-S-C8	4.78	112.67	106.92
4	C	902	MES	O2S-S-C8	4.55	112.39	106.92
4	G	904	MES	O2S-S-C8	4.42	112.24	106.92
4	E	903	MES	O2S-S-C8	4.29	112.08	106.92
4	A	901	MES	O2S-S-C8	4.08	111.83	106.92
4	I	905	MES	O3S-S-O2S	-3.22	103.41	111.27
4	K	906	MES	O3S-S-O2S	-3.12	103.64	111.27
4	G	904	MES	O3S-S-O2S	-3.08	103.74	111.27
4	C	902	MES	O3S-S-O2S	-3.07	103.78	111.27
4	E	903	MES	O3S-S-O2S	-2.98	103.98	111.27
4	A	901	MES	O2S-S-O1S	-2.70	104.59	113.95
4	A	901	MES	O3S-S-O2S	-2.67	104.76	111.27
4	G	904	MES	O2S-S-O1S	-2.56	105.09	113.95
4	K	906	MES	O2S-S-O1S	-2.55	105.11	113.95
4	E	903	MES	O2S-S-O1S	-2.55	105.13	113.95
4	I	905	MES	O2S-S-O1S	-2.55	105.13	113.95
4	C	902	MES	O2S-S-O1S	-2.48	105.35	113.95
4	I	905	MES	O1-C2-C3	-2.03	107.32	111.80
4	K	906	MES	O3S-S-O1S	-2.01	106.36	111.27

There are no chirality outliers.

All (30) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	E	903	MES	C7-C8-S-O1S
4	E	903	MES	C7-C8-S-O3S
4	C	902	MES	C7-C8-S-O1S
4	C	902	MES	C7-C8-S-O3S
4	I	905	MES	C7-C8-S-O1S
4	I	905	MES	C7-C8-S-O3S
4	G	904	MES	C7-C8-S-O1S
4	K	906	MES	C7-C8-S-O1S
4	K	906	MES	C7-C8-S-O3S
4	A	901	MES	C7-C8-S-O1S
4	A	901	MES	C7-C8-S-O2S
4	A	901	MES	C7-C8-S-O3S
4	G	904	MES	C7-C8-S-O3S
4	E	903	MES	C7-C8-S-O2S
4	C	902	MES	C7-C8-S-O2S
4	I	905	MES	C7-C8-S-O2S
4	G	904	MES	C7-C8-S-O2S
4	K	906	MES	C7-C8-S-O2S
8	O	1308	GER	C14-C13-C15-C16
8	R	1608	GER	C14-C13-C15-C16
8	M	1108	GER	C14-C13-C15-C16
8	Q	1508	GER	C14-C13-C15-C16
8	P	1408	GER	C14-C13-C15-C16
8	N	1208	GER	C14-C13-C15-C16
8	M	1108	GER	C10-C11-C12-C13
8	O	1308	GER	C10-C11-C12-C13
8	R	1608	GER	C10-C11-C12-C13
8	P	1408	GER	C10-C11-C12-C13
8	N	1208	GER	C10-C11-C12-C13
8	Q	1508	GER	C10-C11-C12-C13

There are no ring outliers.

12 monomers are involved in 18 short contacts:

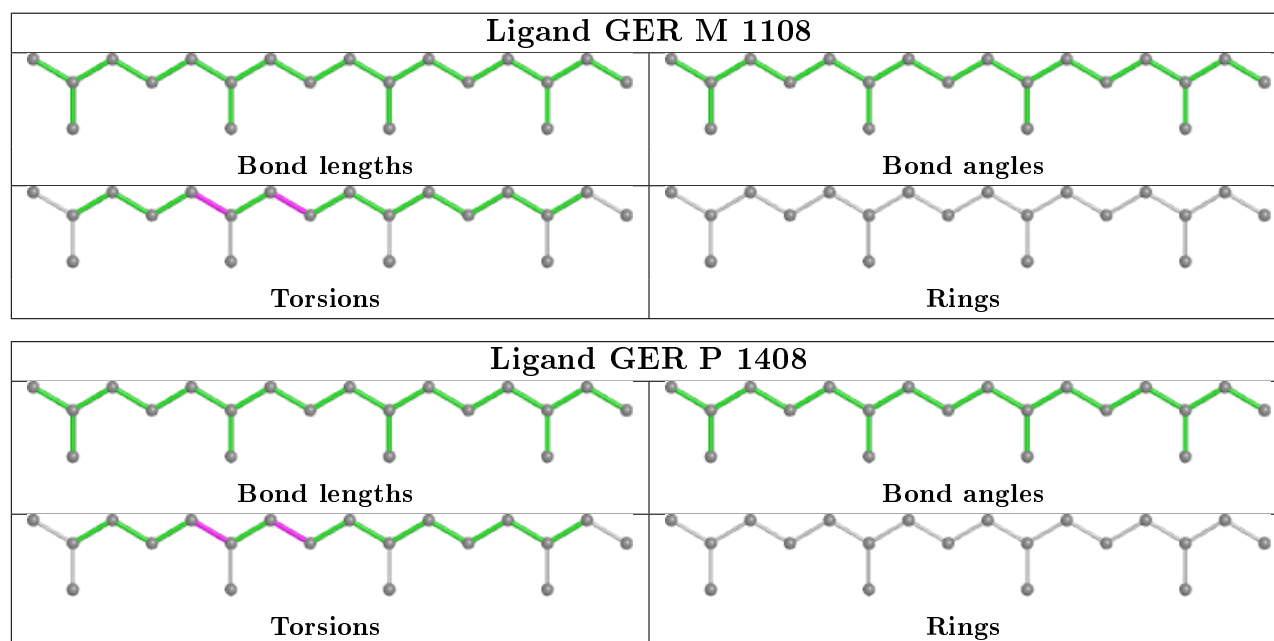
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	E	903	MES	2	0
8	M	1108	GER	1	0
4	K	906	MES	1	0
4	A	901	MES	3	0
8	P	1408	GER	1	0

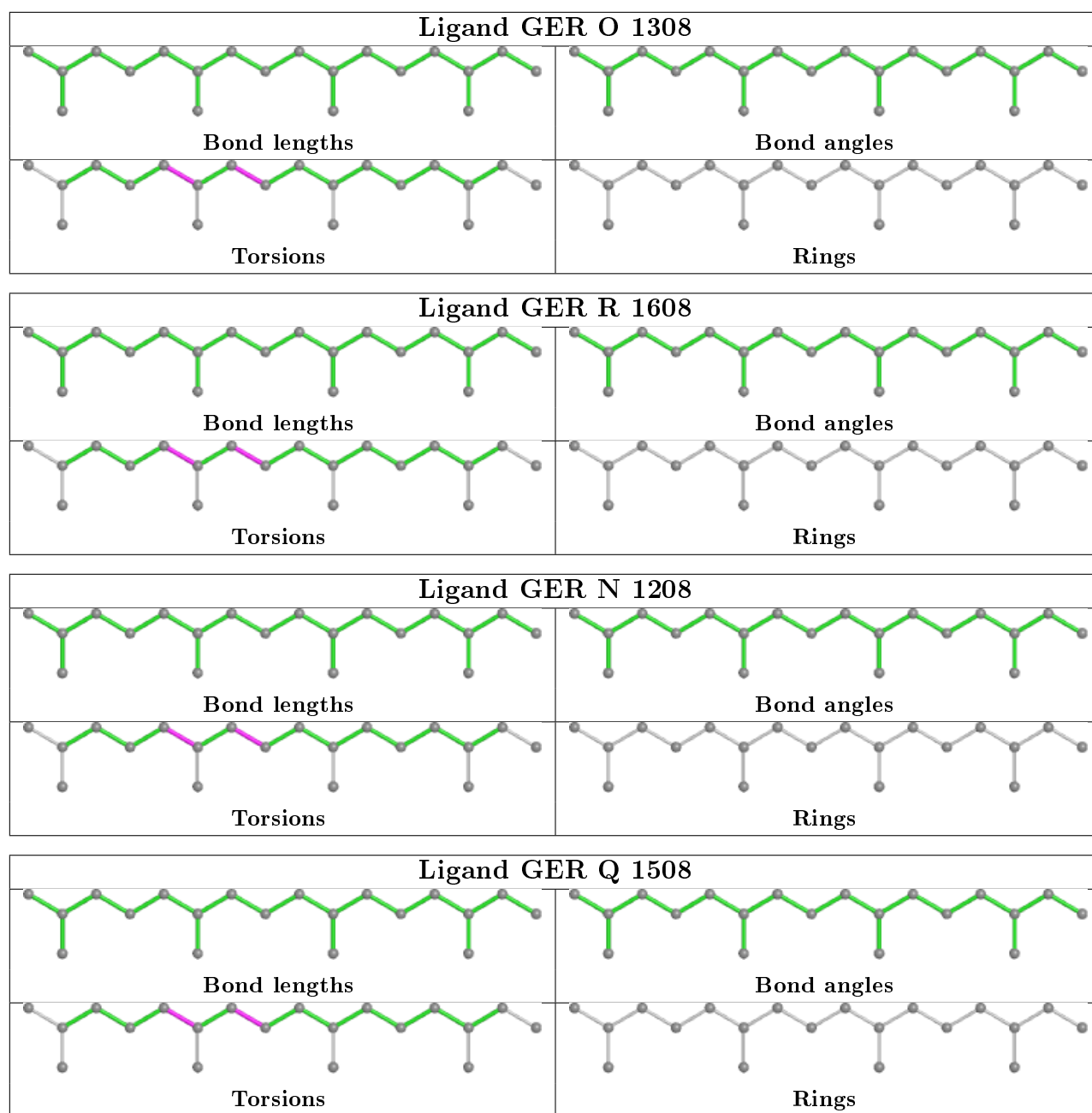
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	I	905	MES	2	0
8	O	1308	GER	1	0
4	G	904	MES	2	0
8	R	1608	GER	2	0
8	N	1208	GER	1	0
8	Q	1508	GER	1	0
4	C	902	MES	1	0

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





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data 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	314/377 (83%)	-0.36	4 (1%) 77 72	36, 57, 91, 105	0
1	C	314/377 (83%)	-0.39	4 (1%) 77 72	35, 55, 81, 99	0
1	E	314/377 (83%)	-0.36	3 (0%) 82 77	34, 57, 83, 103	0
1	G	314/377 (83%)	-0.32	6 (1%) 66 59	38, 58, 85, 104	0
1	I	314/377 (83%)	-0.43	5 (1%) 72 66	32, 52, 80, 91	0
1	K	314/377 (83%)	-0.57	1 (0%) 94 93	25, 43, 67, 82	0
2	B	346/377 (91%)	-0.34	6 (1%) 70 63	36, 52, 78, 100	0
2	D	346/377 (91%)	-0.30	6 (1%) 70 63	34, 48, 78, 92	0
2	F	346/377 (91%)	-0.31	5 (1%) 75 70	32, 48, 77, 103	0
2	H	346/377 (91%)	-0.02	11 (3%) 47 37	37, 64, 92, 111	0
2	J	346/377 (91%)	-0.30	9 (2%) 56 46	31, 50, 79, 102	0
2	L	346/377 (91%)	-0.41	3 (0%) 84 80	27, 42, 66, 88	0
3	M	6/11 (54%)	0.17	1 (16%) 1 1	52, 58, 82, 89	0
3	N	6/11 (54%)	0.18	1 (16%) 1 1	56, 61, 79, 85	0
3	O	6/11 (54%)	0.15	1 (16%) 1 1	52, 59, 78, 84	0
3	P	6/11 (54%)	0.42	1 (16%) 1 1	64, 68, 89, 92	0
3	Q	6/11 (54%)	0.10	0 100 100	51, 56, 72, 79	0
3	R	6/11 (54%)	0.03	0 100 100	51, 57, 71, 76	0
All	All	3996/4590 (87%)	-0.33	67 (1%) 70 63	25, 52, 83, 111	0

All (67) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	H	363	THR	5.0
2	H	108	SER	4.8
2	B	363	THR	4.1

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Mol	Chain	Res	Type	RSRZ
2	J	363	THR	3.9
1	C	306	HIS	3.8
2	F	88	SER	3.8
1	G	305	SER	3.7
2	J	108	SER	3.7
1	E	306	HIS	3.7
1	E	91	ILE	3.5
1	G	306	HIS	3.5
2	D	363	THR	3.4
2	H	361	TRP	3.1
1	I	306	HIS	3.0
1	G	55	PHE	3.0
1	C	305	SER	3.0
2	J	360	SER	2.9
1	I	305	SER	2.9
2	B	86	ASP	2.9
2	H	88	SER	2.9
2	L	363	THR	2.8
2	F	40	TYR	2.7
3	M	106	THR	2.7
2	D	108	SER	2.7
2	F	363	THR	2.7
2	H	127	THR	2.7
2	L	113	THR	2.7
3	N	206	THR	2.7
1	G	326	GLN	2.6
2	F	112	GLY	2.6
1	A	304	PRO	2.6
2	D	37	PRO	2.6
3	P	406	THR	2.6
2	D	88	SER	2.6
1	A	306	HIS	2.5
1	C	329	ASN	2.5
3	O	306	THR	2.4
2	H	65	ASP	2.4
2	L	108	SER	2.4
1	E	304	PRO	2.4
2	H	362	LYS	2.4
2	F	86	ASP	2.3
1	K	306	HIS	2.3
2	J	85	GLU	2.3
2	H	40	TYR	2.3

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Mol	Chain	Res	Type	RSRZ
2	J	110	ASN	2.2
1	A	305	SER	2.2
1	I	85	SER	2.2
2	J	112	GLY	2.2
1	I	328	ASP	2.2
2	H	86	ASP	2.2
2	D	361	TRP	2.2
2	B	84	THR	2.2
2	H	305	LEU	2.2
1	C	304	PRO	2.2
2	J	84	THR	2.2
2	J	114	ALA	2.1
2	J	86	ASP	2.1
1	G	91	ILE	2.1
1	G	304	PRO	2.1
1	I	84	PRO	2.1
2	B	362	LYS	2.1
2	H	110	ASN	2.1
1	A	326	GLN	2.0
2	D	86	ASP	2.0
2	B	40	TYR	2.0
2	B	112	GLY	2.0

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 [i](#)

There are no carbohydrates 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(\AA^2)	Q<0.9
8	GER	P	1408	20/20	0.89	0.31	61,66,73,73	0

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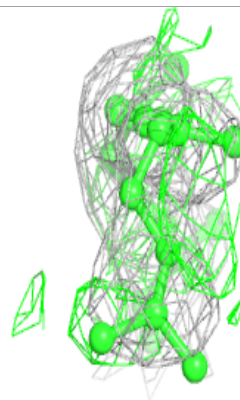
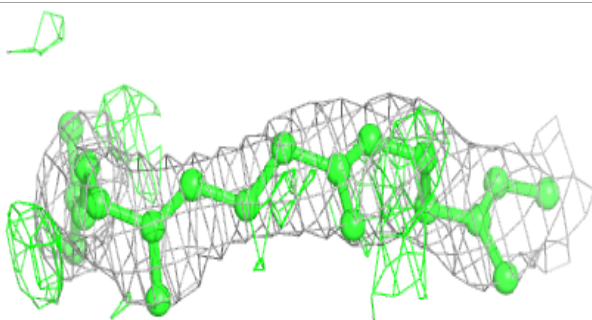
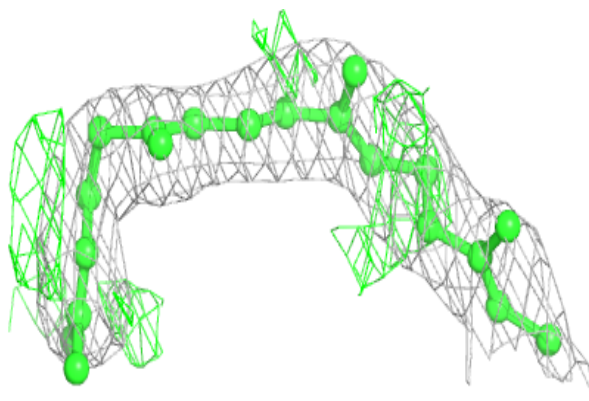
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	MES	G	904	12/12	0.89	0.34	97,101,104,104	0
4	MES	A	901	12/12	0.90	0.33	96,103,106,106	0
4	MES	I	905	12/12	0.91	0.29	78,84,88,89	0
4	MES	C	902	12/12	0.92	0.30	86,93,95,96	0
4	MES	E	903	12/12	0.93	0.26	84,90,94,94	0
8	GER	M	1108	20/20	0.93	0.29	48,52,61,61	0
8	GER	N	1208	20/20	0.93	0.29	49,53,58,59	0
8	GER	R	1608	20/20	0.93	0.34	47,50,54,54	0
4	MES	K	906	12/12	0.93	0.25	65,77,83,84	0
8	GER	O	1308	20/20	0.93	0.26	48,51,58,60	0
8	GER	Q	1508	20/20	0.94	0.30	47,50,55,56	0
7	CL	J	806	1/1	0.96	0.06	58,58,58,58	0
5	ZN	H	378	1/1	0.96	0.04	87,87,87,87	0
7	CL	H	805	1/1	0.96	0.08	60,60,60,60	0
6	SO4	D	811	5/5	0.97	0.15	66,66,67,67	0
5	ZN	J	378	1/1	0.97	0.05	68,68,68,68	0
7	CL	F	803	1/1	0.98	0.05	51,51,51,51	0
6	SO4	F	812	5/5	0.98	0.14	67,67,68,69	0
6	SO4	B	810	5/5	0.98	0.14	67,68,68,69	0
7	CL	K	807	1/1	0.98	0.09	44,44,44,44	0
5	ZN	L	378	1/1	0.98	0.05	58,58,58,58	0
5	ZN	F	378	1/1	0.99	0.05	78,78,78,78	0
7	CL	C	801	1/1	0.99	0.16	53,53,53,53	0
6	SO4	L	815	5/5	0.99	0.16	54,54,56,56	0
6	SO4	H	813	5/5	0.99	0.11	68,69,69,70	0
5	ZN	D	378	1/1	0.99	0.05	66,66,66,66	0
6	SO4	J	814	5/5	0.99	0.13	52,52,54,55	0
7	CL	D	802	1/1	0.99	0.06	45,45,45,45	0
7	CL	G	804	1/1	0.99	0.11	52,52,52,52	0
5	ZN	B	378	1/1	0.99	0.04	77,77,77,77	0

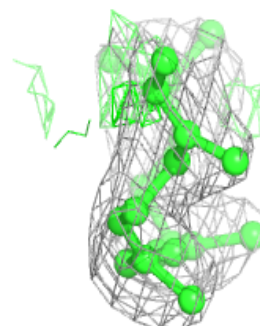
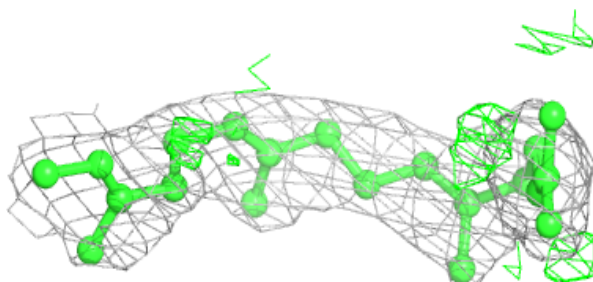
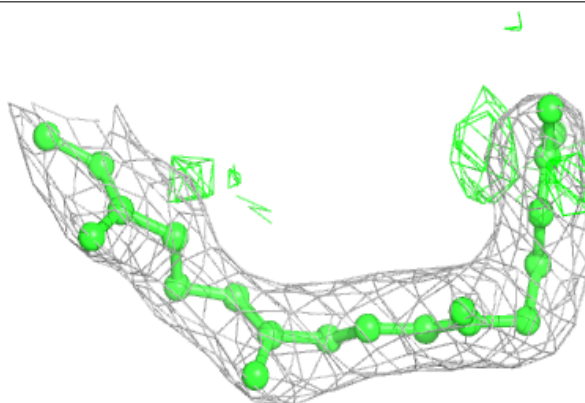
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around GER P 1408:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

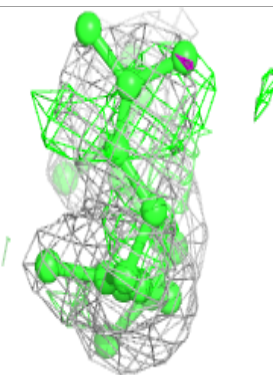
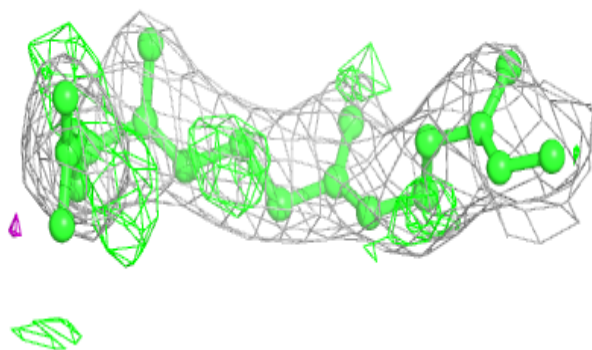
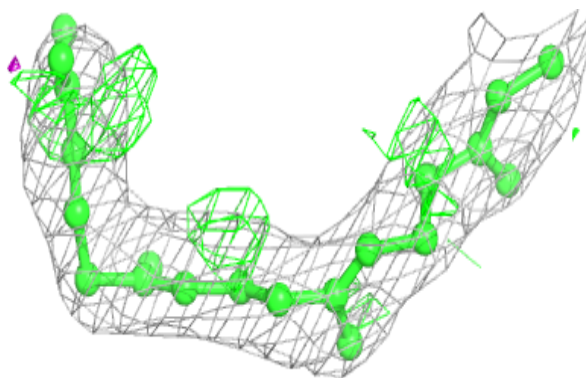
**Electron density around GER M 1108:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

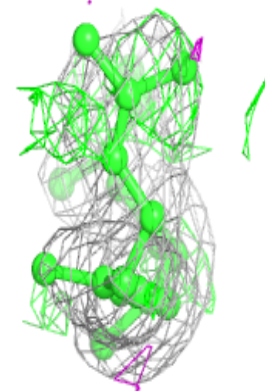
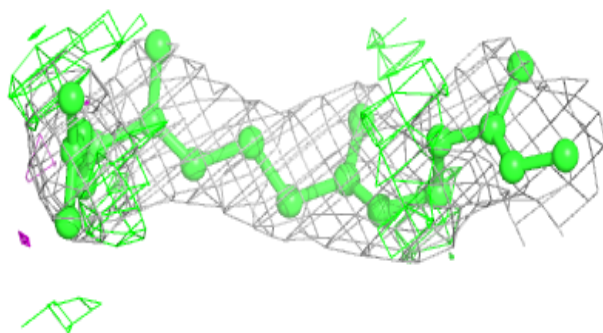
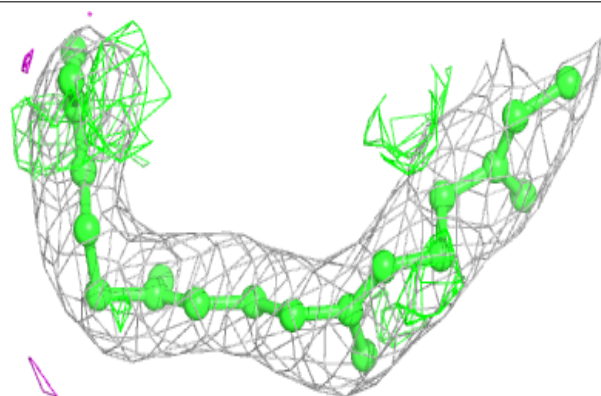


Electron density around GER N 1208:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

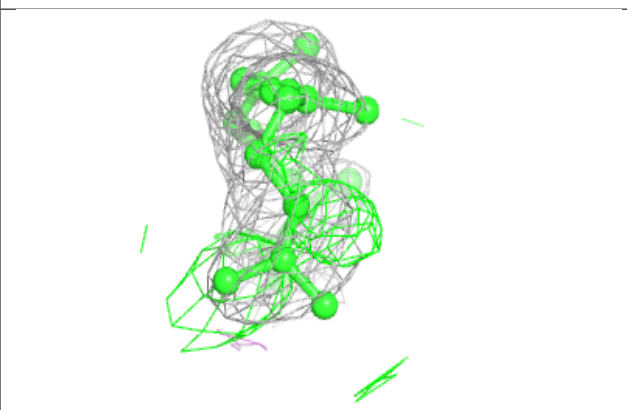
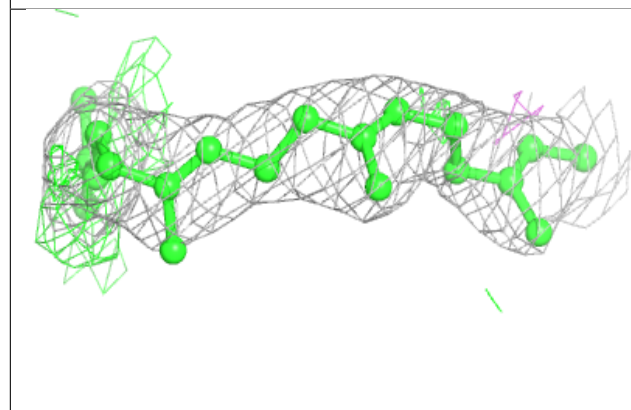
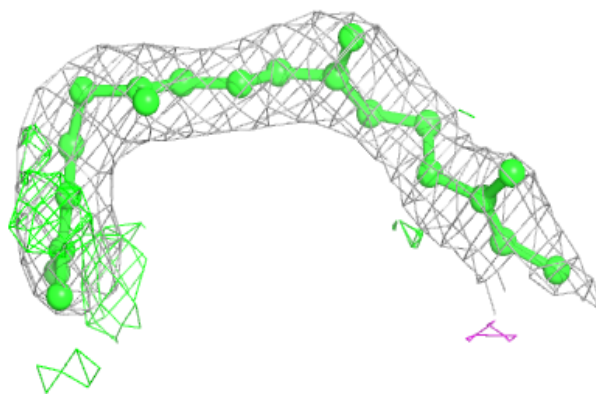
**Electron density around GER R 1608:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

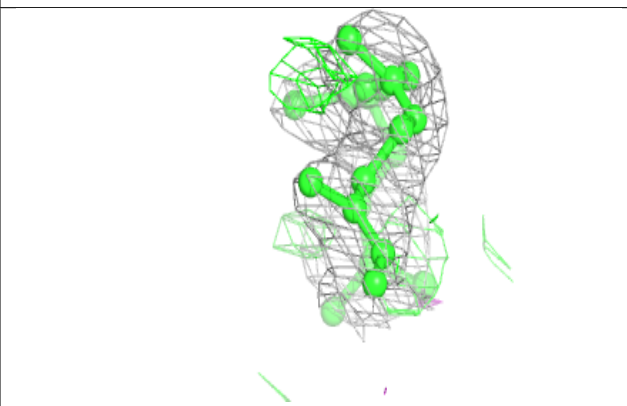
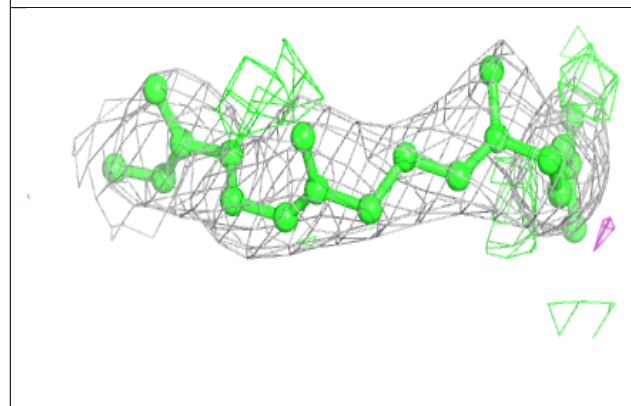
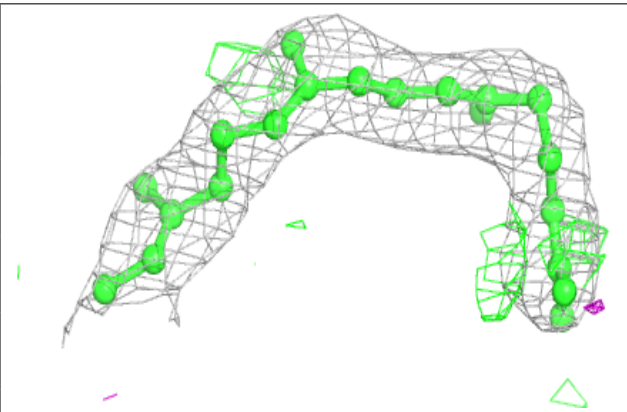


Electron density around GER O 1308:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around GER Q 1508:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.