



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 12:35 PM UTC

PDB ID : 3MP3 / pdb_00003mp3
Title : Crystal Structure of Human Lyase in complex with inhibitor HG-CoA
Authors : Fu, Z.; Runquist, J.A.; Montgomery, C.; Mizioro, H.M.; Kim, J.-J.P.
Deposited on : 2010-04-24
Resolution : 2.40 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
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
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

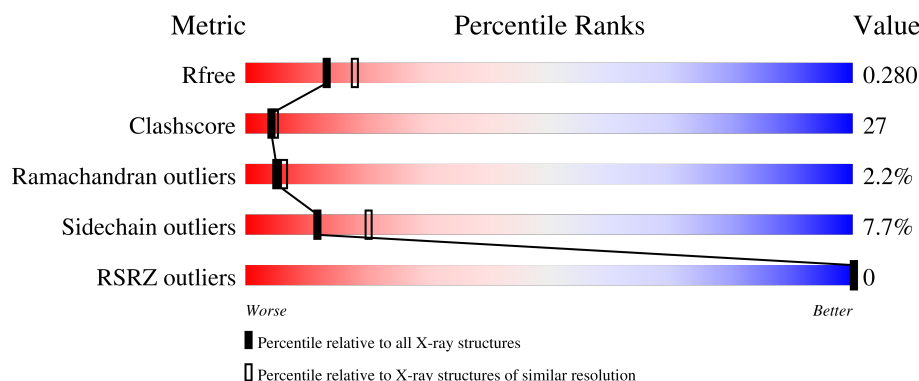
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	298	 62% 32% 5% ..
1	B	298	 59% 35% 5% .
1	C	298	 51% 43% . ..
1	D	298	 44% 45% 7% .
1	E	298	 58% 34% 7% .

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Mol	Chain	Length	Quality of chain
1	F	298	43% 49% 7% •

2 Entry composition [i](#)

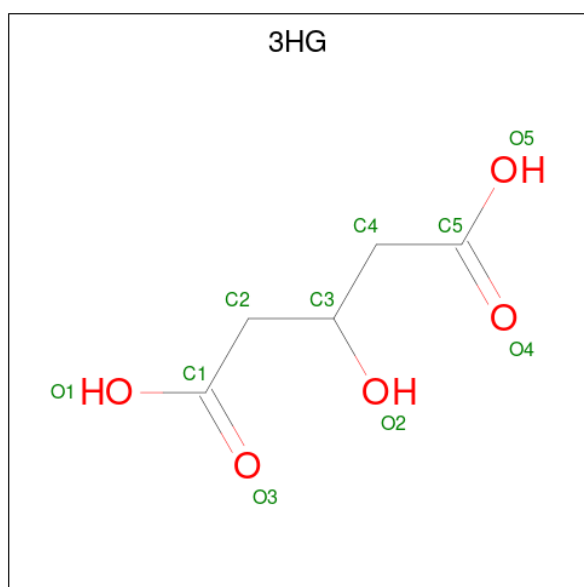
There are 5 unique types of molecules in this entry. The entry contains 13597 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Hydroxymethylglutaryl-CoA lyase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	296	Total	C	N	O	S	0	0	0
			2189	1390	365	417	17			
1	B	296	Total	C	N	O	S	0	0	0
			2189	1390	365	417	17			
1	C	296	Total	C	N	O	S	0	0	0
			2189	1390	365	417	17			
1	D	289	Total	C	N	O	S	0	0	0
			2142	1361	357	408	16			
1	E	296	Total	C	N	O	S	0	0	0
			2189	1390	365	417	17			
1	F	296	Total	C	N	O	S	0	0	0
			2189	1390	365	417	17			

- Molecule 2 is 3-HYDROXYPENTANEDIOIC ACID (CCD ID: 3HG) (formula: $C_5H_8O_5$).

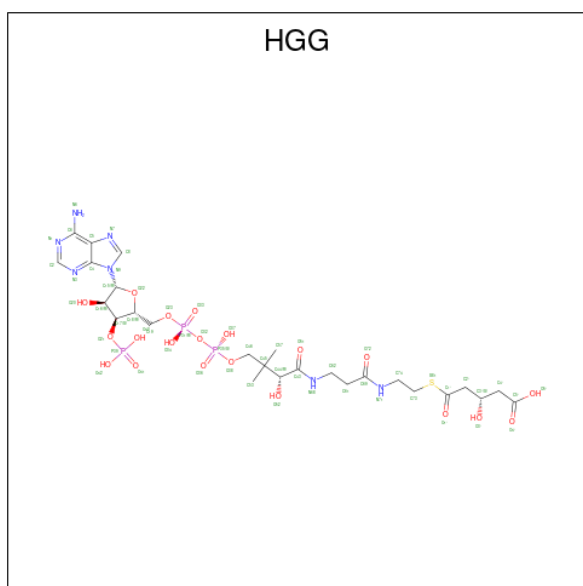


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			10	5	5		

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mg	0	0
			1	1		
3	B	1	Total	Mg	0	0
			1	1		
3	C	1	Total	Mg	0	0
			1	1		
3	D	1	Total	Mg	0	0
			1	1		
3	E	1	Total	Mg	0	0
			1	1		
3	F	1	Total	Mg	0	0
			1	1		

- Molecule 4 is (3R,5S,9R,21S)-1-[(2R,3S,4R,5R)-5-(6-amino-9H-purin-9-yl)-4-hydroxy-3-(phosphonooxy)tetrahydrofuran-2-yl]-3,5,9,21-tetrahydroxy-8,8-dimethyl-10,14,19-trioxo-2,4,6-trioxa-18-thia-11,15-diaza-3,5-diphosphatricosan-23-oic acid 3,5-dioxide (CCD ID: HGG) (formula: C₂₆H₄₂N₇O₂₀P₃S).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
4	B	1	Total	C	N	O	P	S	0	0
			57	26	7	20	3	1		

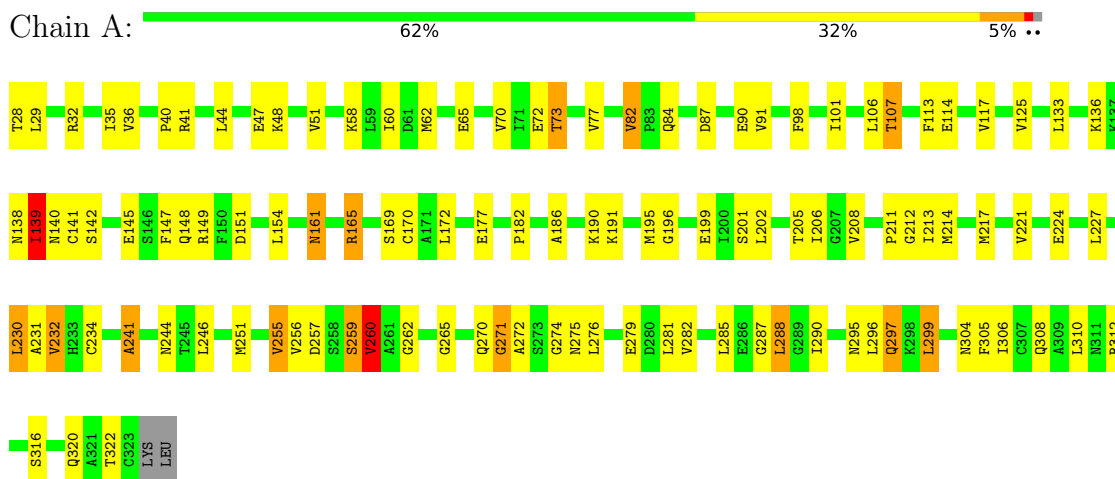
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	100	Total 100	O 100	0	0
5	B	70	Total 70	O 70	0	0
5	C	70	Total 70	O 70	0	0
5	D	65	Total 65	O 65	0	0
5	E	74	Total 74	O 74	0	0
5	F	58	Total 58	O 58	0	0

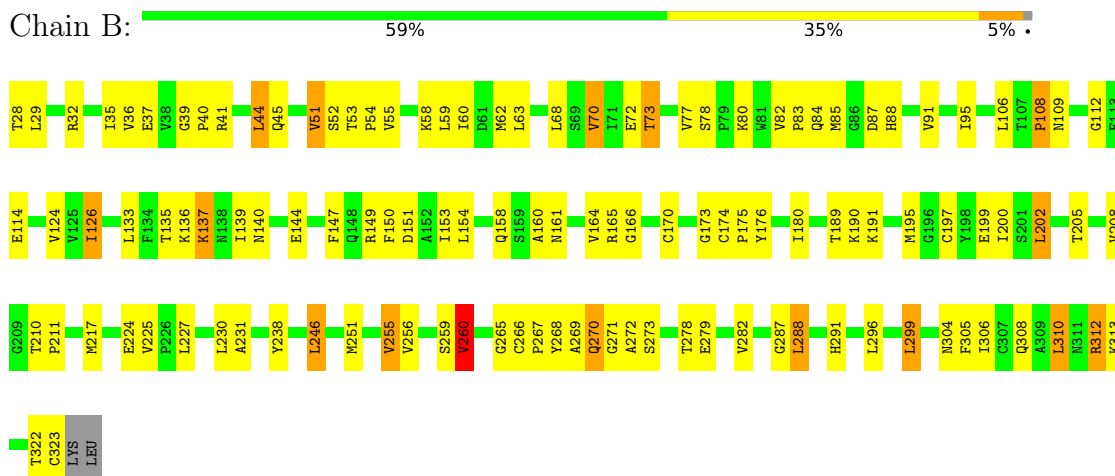
3 Residue-property plots

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: Hydroxymethylglutaryl-CoA lyase

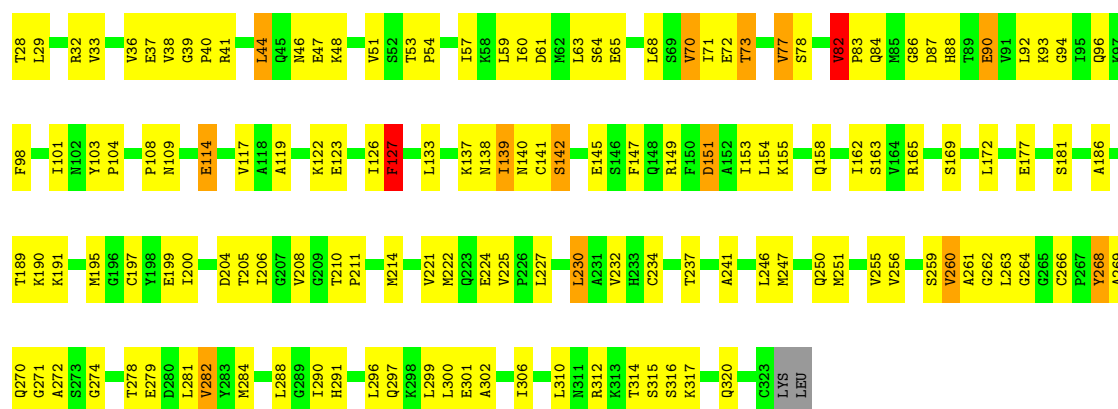


• Molecule 1: Hydroxymethylglutaryl-CoA lyase

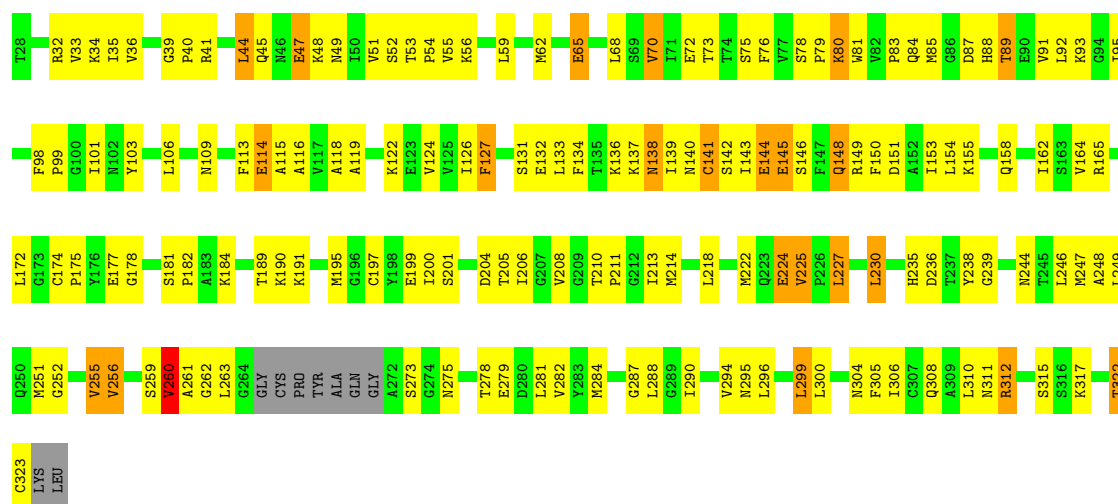


• Molecule 1: Hydroxymethylglutaryl-CoA lyase

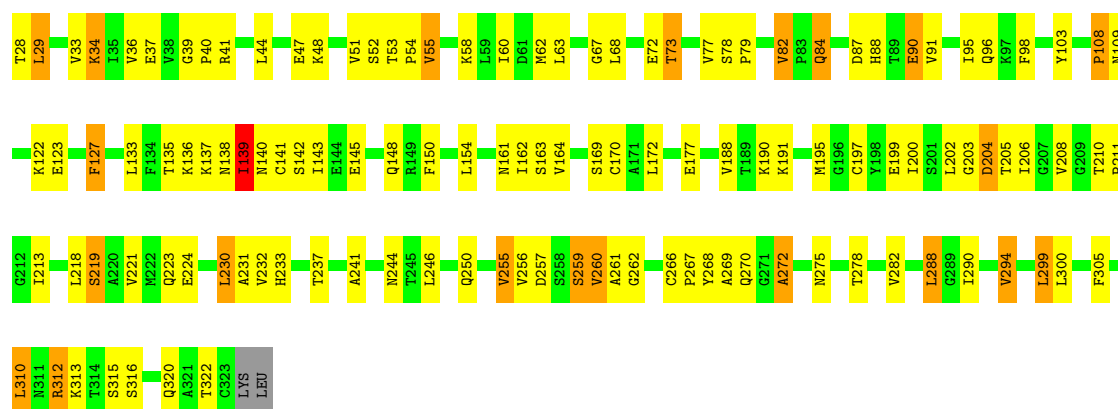




• Molecule 1: Hydroxymethylglutaryl-CoA lyase

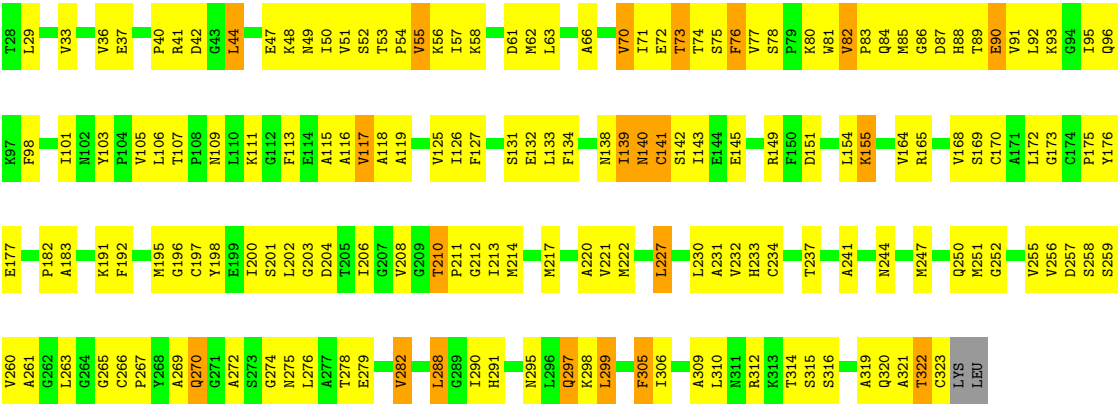


• Molecule 1: Hydroxymethylglutaryl-CoA lyase



• Molecule 1: Hydroxymethylglutaryl-CoA lyase

Chain F: 43% 49% 7% .



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	197.37Å 117.00Å 86.87Å 90.00° 112.31° 90.00°	Depositor
Resolution (Å)	29.70 – 2.40 29.70 – 2.40	Depositor EDS
% Data completeness (in resolution range)	93.3 (29.70-2.40) 93.3 (29.70-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.78 (at 2.39Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.212 , 0.281 0.212 , 0.280	Depositor DCC
R_{free} test set	6749 reflections (10.10%)	wwPDB-VP
Wilson B-factor (Å ²)	22.8	Xtriage
Anisotropy	0.045	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 23.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	13597	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.71% 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: HGG, MG, 3HG

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	0.51	0/2224	0.98	8/3011 (0.3%)
1	B	0.48	0/2224	1.01	9/3011 (0.3%)
1	C	0.48	0/2224	1.00	7/3011 (0.2%)
1	D	0.45	0/2174	0.95	6/2941 (0.2%)
1	E	0.47	0/2224	0.98	7/3011 (0.2%)
1	F	0.42	0/2224	0.93	10/3011 (0.3%)
All	All	0.47	0/13294	0.97	47/17996 (0.3%)

There are no bond length outliers.

All (47) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	146	SER	N-CA-C	-8.62	102.89	113.41
1	A	73	THR	N-CA-C	8.55	122.20	111.69
1	E	73	THR	N-CA-C	8.52	121.80	111.40
1	A	29	LEU	N-CA-C	7.58	118.32	109.60
1	B	73	THR	N-CA-C	6.97	118.77	111.03
1	B	88	HIS	N-CA-C	6.33	118.99	111.33
1	A	29	LEU	CA-C-N	6.25	126.16	119.85
1	A	29	LEU	C-N-CA	6.25	126.16	119.85
1	E	294	VAL	N-CA-C	6.03	117.27	108.53
1	A	262	GLY	N-CA-C	-5.98	105.87	115.08
1	F	227	LEU	N-CA-C	5.96	117.78	111.28
1	A	107	THR	CA-C-N	5.94	125.64	119.05
1	A	107	THR	C-N-CA	5.94	125.64	119.05
1	B	126	ILE	N-CA-C	-5.89	103.63	110.05
1	E	88	HIS	N-CA-C	5.80	118.38	111.71
1	E	204	ASP	N-CA-C	-5.77	100.96	109.62
1	E	29	LEU	N-CA-C	5.77	118.04	110.08
1	E	67	GLY	N-CA-C	5.77	121.34	114.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	127	PHE	N-CA-C	5.75	117.76	108.34
1	F	241	ALA	N-CA-C	5.57	117.81	111.02
1	F	66	ALA	N-CA-C	-5.54	106.03	112.89
1	E	241	ALA	N-CA-C	5.52	117.75	111.02
1	A	241	ALA	N-CA-C	5.50	117.70	111.11
1	F	204	ASP	N-CA-C	-5.49	98.98	108.52
1	C	82	VAL	CA-C-N	5.44	125.06	119.56
1	C	82	VAL	C-N-CA	5.44	125.06	119.56
1	C	73	THR	N-CA-C	5.41	118.00	111.40
1	D	127	PHE	N-CA-C	5.36	116.82	107.49
1	B	202	LEU	CA-C-N	-5.33	118.50	122.18
1	B	202	LEU	C-N-CA	-5.33	118.50	122.18
1	C	282	VAL	N-CA-C	-5.28	104.49	111.09
1	F	29	LEU	N-CA-C	5.25	116.50	109.83
1	D	225	VAL	N-CA-C	5.22	114.21	108.05
1	D	81	TRP	N-CA-C	-5.18	105.63	111.28
1	F	29	LEU	CA-C-N	5.18	125.08	119.85
1	F	29	LEU	C-N-CA	5.18	125.08	119.85
1	C	241	ALA	N-CA-C	5.15	117.31	111.02
1	B	287	GLY	N-CA-C	-5.13	107.95	114.16
1	D	148	GLN	N-CA-C	-5.12	105.61	111.14
1	F	88	HIS	N-CA-C	5.11	117.24	111.11
1	F	275	ASN	N-CA-C	5.09	117.25	110.53
1	B	227	LEU	N-CA-C	5.08	117.72	111.82
1	C	88	HIS	N-CA-C	5.07	118.24	111.75
1	B	137	LYS	N-CA-C	5.07	118.73	112.54
1	D	227	LEU	N-CA-C	5.06	117.19	111.11
1	B	260	VAL	N-CA-C	5.04	119.82	109.34
1	F	117	VAL	N-CA-C	-5.00	105.67	110.72

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2189	0	2244	96	0
1	B	2189	0	2244	95	0
1	C	2189	0	2244	141	0
1	D	2142	0	2203	144	0
1	E	2189	0	2244	109	0
1	F	2189	0	2244	152	0
2	A	10	0	6	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
3	E	1	0	0	0	0
3	F	1	0	0	0	0
4	B	57	0	37	10	0
5	A	100	0	0	6	0
5	B	70	0	0	3	0
5	C	70	0	0	7	0
5	D	65	0	0	3	0
5	E	74	0	0	4	0
5	F	58	0	0	3	0
All	All	13597	0	13466	721	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 27.

All (721) 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:D:47:GLU:HG2	1:D:310:LEU:HD21	1.46	0.98
1:E:211:PRO:HG2	1:F:288:LEU:HD11	1.44	0.96
1:D:263:LEU:HD11	1:D:306:ILE:HD12	1.48	0.96
1:E:84:GLN:H	1:E:84:GLN:NE2	1.65	0.94
1:E:51:VAL:HB	1:E:55:VAL:HG21	1.48	0.93
1:D:92:LEU:HD23	1:D:119:ALA:HB3	1.51	0.92
1:C:266:CYS:HB3	1:C:269:ALA:HB3	1.52	0.91
1:C:172:LEU:HD11	1:C:214:MET:HE2	1.55	0.89
1:A:190:LYS:HD3	1:A:224:GLU:HB3	1.56	0.87
1:C:48:LYS:H	1:C:48:LYS:HD2	1.40	0.87
1:D:142:SER:HB2	1:D:145:GLU:HB2	1.57	0.86
1:A:48:LYS:N	1:A:48:LYS:HD2	1.90	0.86
1:C:84:GLN:CD	1:C:84:GLN:H	1.85	0.85
1:F:52:SER:OG	1:F:55:VAL:HG12	1.77	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:93:LYS:HE3	1:C:119:ALA:HA	1.57	0.84
1:C:149:ARG:HH11	1:E:161:ASN:HD22	1.21	0.84
1:E:63:LEU:HD21	1:E:260:VAL:HG11	1.60	0.83
1:C:36:VAL:HA	1:C:70:VAL:HG13	1.61	0.83
1:A:48:LYS:HD2	1:A:48:LYS:H	1.40	0.83
1:E:68:LEU:HD13	1:E:294:VAL:HG21	1.61	0.83
1:A:288:LEU:HD11	1:B:211:PRO:HG2	1.60	0.83
1:B:28:THR:HG22	1:B:29:LEU:H	1.44	0.83
1:B:58:LYS:NZ	1:B:62:MET:HE2	1.94	0.82
1:C:47:GLU:HG2	1:C:310:LEU:HD21	1.61	0.82
1:C:28:THR:HG22	1:C:29:LEU:H	1.43	0.82
1:D:52:SER:OG	1:D:55:VAL:HG23	1.80	0.82
1:C:191:LYS:HE2	1:C:195:MET:HE3	1.63	0.81
1:D:59:LEU:HD22	1:D:306:ILE:HG21	1.61	0.81
1:D:190:LYS:HD2	1:D:224:GLU:HB3	1.62	0.81
1:B:265:GLY:O	1:B:267:PRO:HD3	1.81	0.81
1:E:28:THR:HG22	1:E:29:LEU:H	1.46	0.81
1:E:84:GLN:H	1:E:84:GLN:HE21	1.30	0.80
1:B:251:MET:HE2	1:B:251:MET:HA	1.63	0.80
1:E:260:VAL:HG23	1:E:261:ALA:H	1.46	0.80
1:A:36:VAL:HA	1:A:70:VAL:HG13	1.64	0.79
1:E:51:VAL:HG13	1:E:310:LEU:HD13	1.64	0.79
1:C:39:GLY:HA3	1:C:260:VAL:HG23	1.63	0.79
1:C:33:VAL:HG23	1:C:290:ILE:HG21	1.65	0.78
1:F:134:PHE:O	1:F:138:ASN:HB2	1.83	0.77
1:E:52:SER:O	1:E:55:VAL:HG23	1.82	0.77
1:E:141:CYS:HB2	1:E:145:GLU:HB2	1.66	0.77
1:F:50:ILE:HG22	1:F:84:GLN:HG3	1.67	0.77
1:F:138:ASN:HD22	1:F:139:ILE:HG12	1.51	0.76
1:D:138:ASN:ND2	1:D:139:ILE:HG13	2.00	0.76
1:C:60:ILE:HD13	1:C:73:THR:HG23	1.67	0.76
1:D:151:ASP:O	1:D:155:LYS:HB2	1.84	0.76
1:F:278:THR:O	1:F:282:VAL:HG12	1.84	0.76
1:F:279:GLU:HG3	1:F:299:LEU:HD13	1.67	0.76
1:F:306:ILE:O	1:F:310:LEU:HD13	1.85	0.76
1:F:90:GLU:H	1:F:90:GLU:CD	1.93	0.75
1:F:133:LEU:HD13	1:F:177:GLU:HG2	1.68	0.75
1:C:270:GLN:O	1:C:272:ALA:N	2.19	0.75
1:D:109:ASN:HA	1:D:153:ILE:HD11	1.68	0.74
1:E:270:GLN:HG3	1:F:321:ALA:O	1.88	0.74
1:A:47:GLU:HG2	1:A:310:LEU:HD21	1.69	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:58:LYS:HZ3	1:B:62:MET:HE2	1.50	0.74
1:E:154:LEU:CD1	1:E:195:MET:HB3	2.18	0.74
1:B:322:THR:O	1:B:323:CYS:HB2	1.87	0.73
1:C:59:LEU:HD13	1:C:306:ILE:HB	1.70	0.73
1:F:113:PHE:O	1:F:117:VAL:HG23	1.87	0.73
1:E:40:PRO:HG2	1:E:72:GLU:O	1.89	0.73
1:A:138:ASN:C	1:A:139:ILE:HG12	2.12	0.73
1:C:264:GLY:O	1:C:274:GLY:HA3	1.88	0.73
1:F:75:SER:O	1:F:77:VAL:HG12	1.88	0.72
1:B:191:LYS:HG3	1:B:195:MET:HE2	1.71	0.72
1:A:206:ILE:HG13	1:A:208:VAL:HG13	1.72	0.72
1:B:77:VAL:HG11	4:B:399:HGG:H61	1.72	0.72
1:A:147:PHE:CE1	1:A:191:LYS:HG2	2.25	0.71
1:C:40:PRO:HG2	1:C:72:GLU:O	1.90	0.71
1:A:141:CYS:HA	1:A:145:GLU:OE2	1.89	0.71
1:E:316:SER:O	1:E:320:GLN:HG3	1.90	0.71
1:B:133:LEU:HD23	1:B:176:TYR:HB3	1.71	0.71
1:E:237:THR:HG23	1:E:272:ALA:HB3	1.73	0.71
1:C:83:PRO:HD2	1:C:84:GLN:HE22	1.56	0.71
1:D:172:LEU:HD12	1:D:204:ASP:HB2	1.73	0.71
1:E:58:LYS:HD3	1:E:305:PHE:CZ	2.25	0.70
1:F:36:VAL:HG13	1:F:70:VAL:HG22	1.73	0.70
1:A:36:VAL:HG22	1:A:70:VAL:HG11	1.74	0.70
1:F:58:LYS:HD3	1:F:305:PHE:CZ	2.27	0.70
1:F:33:VAL:HG23	1:F:290:ILE:HG21	1.73	0.70
1:B:150:PHE:O	1:B:154:LEU:HD12	1.92	0.70
1:F:51:VAL:HG23	1:F:56:LYS:HE3	1.74	0.70
1:C:284:MET:O	1:C:288:LEU:HB2	1.92	0.70
1:C:262:GLY:HA3	1:C:315:SER:HB2	1.73	0.69
1:C:149:ARG:HH11	1:E:161:ASN:ND2	1.89	0.69
1:B:108:PRO:HG2	1:B:109:ASN:H	1.56	0.69
1:D:59:LEU:HD22	1:D:306:ILE:CG2	2.22	0.69
1:E:91:VAL:O	1:E:95:ILE:HG23	1.93	0.69
1:D:322:THR:O	1:D:323:CYS:HB2	1.91	0.69
1:C:149:ARG:NH1	1:E:161:ASN:HD22	1.91	0.69
1:B:109:ASN:ND2	4:B:399:HGG:H57A	2.06	0.69
1:D:263:LEU:CD1	1:D:306:ILE:HD12	2.23	0.69
1:E:48:LYS:HD3	5:E:668:HOH:O	1.91	0.69
1:A:36:VAL:HA	1:A:70:VAL:CG1	2.23	0.68
1:F:74:THR:H	1:F:105:VAL:HG12	1.59	0.68
1:A:259:SER:HB2	1:A:275:ASN:HB2	1.75	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:37:GLU:HG2	1:C:63:LEU:HD13	1.76	0.68
1:E:53:THR:N	1:E:54:PRO:HD2	2.09	0.68
1:E:278:THR:HG22	1:E:299:LEU:HD11	1.76	0.68
1:B:278:THR:HG22	1:B:299:LEU:HD11	1.76	0.67
1:F:52:SER:O	1:F:55:VAL:HG13	1.92	0.67
1:C:82:VAL:HG23	1:C:84:GLN:OE1	1.94	0.67
1:E:300:LEU:HD13	1:E:322:THR:HG21	1.76	0.67
1:F:36:VAL:HG21	1:F:165:ARG:HH22	1.59	0.67
1:A:251:MET:HA	1:A:251:MET:HE2	1.76	0.67
1:D:222:MET:HE3	1:D:227:LEU:HD13	1.77	0.67
1:F:53:THR:N	1:F:54:PRO:HD2	2.10	0.66
1:C:48:LYS:HD2	1:C:48:LYS:N	2.07	0.66
1:B:59:LEU:HD13	1:B:306:ILE:HB	1.77	0.66
1:C:211:PRO:HG2	1:D:288:LEU:HD21	1.77	0.66
1:B:112:GLY:CA	4:B:399:HGG:N6	2.59	0.66
1:F:155:LYS:N	1:F:155:LYS:HE2	2.11	0.66
1:F:87:ASP:O	1:F:91:VAL:HG23	1.95	0.66
1:D:92:LEU:HD23	1:D:119:ALA:CB	2.26	0.66
1:F:98:PHE:HB2	1:F:101:ILE:HD12	1.77	0.65
1:D:76:PHE:CZ	1:D:116:ALA:HA	2.31	0.65
1:F:55:VAL:HG23	1:F:305:PHE:CE2	2.31	0.65
1:A:297:GLN:CD	1:A:297:GLN:H	2.04	0.65
1:F:297:GLN:HB3	5:F:560:HOH:O	1.97	0.65
1:A:161:ASN:HD21	1:F:149:ARG:HH22	1.45	0.65
1:C:93:LYS:CE	1:C:119:ALA:HA	2.26	0.65
1:D:200:ILE:HD12	1:D:200:ILE:N	2.11	0.65
1:D:191:LYS:O	1:D:195:MET:HB2	1.97	0.64
1:E:37:GLU:HG2	1:E:63:LEU:HD13	1.79	0.64
1:E:154:LEU:HD13	1:E:195:MET:HB3	1.79	0.64
1:F:202:LEU:HD13	1:F:217:MET:SD	2.37	0.64
1:B:144:GLU:OE1	1:B:147:PHE:HD1	1.81	0.64
1:D:62:MET:HE1	1:D:305:PHE:CG	2.33	0.64
1:D:246:LEU:HD13	1:D:246:LEU:C	2.22	0.64
1:F:106:LEU:HD12	1:F:106:LEU:N	2.13	0.64
1:A:58:LYS:HD3	1:A:305:PHE:CZ	2.32	0.64
1:B:112:GLY:HA2	4:B:399:HGG:N6	2.11	0.64
1:F:81:TRP:O	1:F:83:PRO:HD3	1.97	0.64
1:A:40:PRO:HG2	1:A:72:GLU:O	1.97	0.64
1:C:96:GLN:HB2	1:C:98:PHE:HE1	1.62	0.64
1:D:165:ARG:NH2	1:D:199:GLU:OE2	2.31	0.64
1:F:316:SER:O	1:F:320:GLN:HG3	1.97	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:151:ASP:HB3	5:C:573:HOH:O	1.97	0.64
1:C:278:THR:O	1:C:282:VAL:HG13	1.98	0.63
1:C:317:LYS:HA	1:C:320:GLN:OE1	1.98	0.63
1:C:127:PHE:CD1	1:C:127:PHE:N	2.65	0.63
1:E:51:VAL:HB	1:E:55:VAL:CG2	2.28	0.63
1:E:169:SER:HB2	1:E:205:THR:OG1	1.98	0.63
1:F:154:LEU:HD13	1:F:195:MET:HB3	1.79	0.63
1:F:78:SER:HB2	1:F:80:LYS:HG3	1.81	0.63
1:A:172:LEU:HD11	1:A:214:MET:HE2	1.81	0.63
1:A:169:SER:HB2	1:A:205:THR:OG1	1.99	0.62
1:B:304:ASN:O	1:B:308:GLN:HG3	1.99	0.62
1:D:115:ALA:O	1:D:118:ALA:HB3	2.00	0.62
1:F:62:MET:HE1	1:F:305:PHE:CD2	2.33	0.62
1:C:57:ILE:HD13	1:C:94:GLY:HA3	1.82	0.62
1:D:33:VAL:HG23	1:D:290:ILE:HG21	1.82	0.62
1:F:138:ASN:HD22	1:F:139:ILE:H	1.46	0.62
1:B:202:LEU:HD13	1:B:217:MET:SD	2.40	0.62
1:D:246:LEU:HD13	1:D:246:LEU:O	1.99	0.62
1:F:47:GLU:HG2	1:F:310:LEU:HD23	1.82	0.62
1:E:28:THR:HG22	1:E:29:LEU:N	2.15	0.62
1:C:38:VAL:HG12	1:C:72:GLU:HB2	1.81	0.61
1:B:40:PRO:HG2	1:B:72:GLU:O	2.01	0.61
1:B:62:MET:HE1	1:B:305:PHE:CG	2.36	0.61
1:A:316:SER:O	1:A:320:GLN:HG3	2.00	0.61
1:C:154:LEU:HD13	1:C:195:MET:HG2	1.83	0.61
1:F:172:LEU:HD22	1:F:213:ILE:HG22	1.83	0.61
1:B:112:GLY:CA	4:B:399:HGG:HN6A	2.14	0.61
1:D:278:THR:HG22	1:D:299:LEU:HD11	1.82	0.61
1:C:288:LEU:HD13	1:D:211:PRO:HG2	1.83	0.60
1:E:33:VAL:HG23	1:E:290:ILE:HG21	1.81	0.60
1:C:60:ILE:CD1	1:C:73:THR:HG23	2.31	0.60
1:E:41:ARG:C	1:E:41:ARG:HD2	2.26	0.60
1:F:231:ALA:HA	1:F:255:VAL:HG13	1.81	0.60
1:A:58:LYS:HZ3	1:A:62:MET:HE2	1.66	0.60
1:A:295:ASN:OD1	1:A:297:GLN:HG2	2.01	0.60
1:B:151:ASP:HB3	5:B:985:HOH:O	2.01	0.60
1:D:210:THR:HB	1:D:211:PRO:HD2	1.83	0.60
1:F:279:GLU:HA	1:F:282:VAL:HG13	1.83	0.60
1:A:172:LEU:HD22	1:A:213:ILE:HG22	1.84	0.60
1:A:260:VAL:O	1:A:279:GLU:OE1	2.18	0.60
1:B:291:HIS:HB2	5:B:961:HOH:O	2.00	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:77:VAL:HG11	1:C:82:VAL:HG11	1.82	0.60
1:C:263:LEU:HD11	1:C:306:ILE:HG12	1.82	0.60
1:D:62:MET:HA	1:D:65:GLU:OE2	2.01	0.60
1:C:261:ALA:HB1	1:C:314:THR:OG1	2.02	0.60
1:F:133:LEU:HD13	1:F:177:GLU:CG	2.31	0.60
1:D:41:ARG:HD2	1:D:41:ARG:C	2.27	0.60
1:D:59:LEU:HD13	1:D:306:ILE:CG2	2.31	0.60
1:C:133:LEU:HD13	1:C:177:GLU:HG2	1.82	0.60
1:A:82:VAL:HG23	1:A:84:GLN:OE1	2.02	0.60
1:E:288:LEU:HD11	1:F:211:PRO:HG2	1.84	0.59
1:C:151:ASP:O	1:C:155:LYS:HD3	2.01	0.59
1:C:165:ARG:NH2	1:C:199:GLU:OE2	2.32	0.59
1:B:322:THR:O	1:B:323:CYS:CB	2.50	0.59
1:A:58:LYS:NZ	1:A:62:MET:HE2	2.17	0.59
1:C:127:PHE:N	1:C:127:PHE:HD1	2.01	0.59
1:A:138:ASN:O	1:A:139:ILE:HG23	2.03	0.59
1:D:131:SER:HB3	1:D:134:PHE:HB3	1.85	0.59
1:F:89:THR:O	1:F:93:LYS:HG3	2.02	0.59
1:C:28:THR:HG22	1:C:29:LEU:N	2.15	0.59
1:C:282:VAL:CG2	1:C:296:LEU:HD13	2.32	0.59
1:E:133:LEU:HB3	1:E:177:GLU:HG3	1.84	0.59
1:C:41:ARG:C	1:C:41:ARG:HD2	2.28	0.59
1:C:282:VAL:HG21	1:C:296:LEU:HD13	1.85	0.59
1:E:312:ARG:HD3	1:E:313:LYS:O	2.03	0.59
1:C:84:GLN:CD	1:C:84:GLN:N	2.58	0.59
1:D:140:ASN:O	1:D:141:CYS:HB3	2.02	0.59
1:B:53:THR:N	1:B:54:PRO:HD2	2.17	0.59
1:F:227:LEU:HA	1:F:230:LEU:HD23	1.83	0.59
1:D:34:LYS:HD3	1:D:70:VAL:HG11	1.85	0.58
1:D:36:VAL:HG13	1:D:70:VAL:HG22	1.85	0.58
1:D:175:PRO:HD3	1:D:206:ILE:HD13	1.83	0.58
1:F:206:ILE:HG13	1:F:208:VAL:HG22	1.84	0.58
1:E:84:GLN:HG2	5:E:570:HOH:O	2.03	0.58
1:E:221:VAL:HG12	1:E:230:LEU:HD21	1.85	0.58
1:D:154:LEU:O	1:D:158:GLN:HG3	2.02	0.58
1:B:109:ASN:HD22	4:B:399:HGG:H57A	1.67	0.58
1:C:221:VAL:HG12	1:C:230:LEU:HD21	1.85	0.58
1:F:87:ASP:O	1:F:90:GLU:HG2	2.04	0.58
1:A:62:MET:HE1	1:A:305:PHE:CG	2.38	0.58
1:D:227:LEU:HD21	1:D:252:GLY:HA3	1.84	0.57
1:F:210:THR:C	1:F:244:ASN:HD21	2.12	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:106:LEU:H	1:F:106:LEU:CD1	2.17	0.57
1:A:98:PHE:HB2	1:A:101:ILE:HD12	1.86	0.57
1:C:96:GLN:HB2	1:C:98:PHE:CE1	2.39	0.57
1:C:44:LEU:HD11	1:C:306:ILE:HD11	1.87	0.57
1:D:109:ASN:CA	1:D:153:ILE:HD11	2.34	0.57
1:A:141:CYS:HB2	1:A:145:GLU:CB	2.35	0.57
1:B:87:ASP:O	1:B:91:VAL:HG23	2.04	0.57
1:C:28:THR:CG2	1:C:29:LEU:H	2.12	0.57
1:E:260:VAL:HG23	1:E:261:ALA:N	2.19	0.57
1:A:65:GLU:HG3	1:A:98:PHE:CZ	2.40	0.57
1:D:84:GLN:H	1:D:84:GLN:CD	2.13	0.57
1:F:131:SER:O	1:F:143:ILE:HD11	2.05	0.57
1:F:201:SER:HA	1:F:231:ALA:HB3	1.86	0.57
1:C:142:SER:HG	1:C:145:GLU:HG3	1.69	0.57
1:C:297:GLN:O	1:C:301:GLU:HG3	2.04	0.57
1:D:53:THR:N	1:D:54:PRO:HD2	2.19	0.57
1:E:87:ASP:O	1:E:91:VAL:HG23	2.05	0.56
1:B:28:THR:HG22	1:B:29:LEU:N	2.18	0.56
1:C:186:ALA:HA	1:C:221:VAL:HG22	1.87	0.56
1:D:83:PRO:HG2	1:D:84:GLN:NE2	2.20	0.56
1:E:82:VAL:O	1:E:82:VAL:HG22	2.04	0.56
1:B:52:SER:OG	1:B:55:VAL:HG23	2.05	0.56
1:A:77:VAL:HG11	1:A:82:VAL:HG13	1.88	0.56
1:E:231:ALA:CB	1:E:255:VAL:HG22	2.36	0.56
1:F:62:MET:HE1	1:F:305:PHE:CG	2.40	0.56
1:A:77:VAL:CG1	1:A:82:VAL:HG13	2.35	0.56
1:B:190:LYS:HD3	1:B:224:GLU:HB3	1.87	0.56
1:E:191:LYS:O	1:E:195:MET:HG3	2.06	0.56
1:F:142:SER:H	1:F:145:GLU:HB2	1.71	0.56
1:C:32:ARG:HB3	1:C:291:HIS:HB3	1.87	0.56
1:F:200:ILE:HD12	1:F:200:ILE:N	2.21	0.56
1:D:91:VAL:O	1:D:95:ILE:HG23	2.05	0.56
1:E:52:SER:OG	1:E:55:VAL:HG22	2.06	0.56
1:F:106:LEU:HD12	1:F:106:LEU:H	1.70	0.56
1:D:59:LEU:CD2	1:D:306:ILE:HG21	2.33	0.56
1:A:77:VAL:HG11	1:A:82:VAL:CG1	2.36	0.56
1:D:89:THR:O	1:D:93:LYS:HG3	2.06	0.56
1:D:151:ASP:OD1	1:D:155:LYS:HD3	2.06	0.56
1:E:138:ASN:O	1:E:139:ILE:HG23	2.06	0.55
1:A:271:GLY:HA2	5:A:505:HOH:O	2.05	0.55
1:E:63:LEU:HD21	1:E:260:VAL:CG1	2.35	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:83:PRO:HG2	1:C:84:GLN:NE2	2.21	0.55
1:C:234:CYS:HB3	5:C:817:HOH:O	2.06	0.55
1:C:123:GLU:HG3	1:C:163:SER:CB	2.36	0.55
1:F:258:SER:HB2	1:F:276:LEU:O	2.07	0.55
1:C:133:LEU:HD13	1:C:177:GLU:CG	2.36	0.55
1:F:131:SER:HB3	1:F:134:PHE:HB3	1.88	0.55
1:F:74:THR:OG1	1:F:75:SER:N	2.39	0.55
1:E:230:LEU:O	1:E:255:VAL:HG13	2.07	0.55
1:F:85:MET:HE2	1:F:85:MET:HA	1.88	0.54
1:F:261:ALA:HB1	1:F:314:THR:OG1	2.06	0.54
1:C:139:ILE:HD13	1:C:139:ILE:N	2.22	0.54
1:B:35:ILE:CD1	1:B:282:VAL:HG12	2.37	0.54
1:C:77:VAL:HB	1:C:108:PRO:HG3	1.90	0.54
1:C:126:ILE:HD11	1:C:197:CYS:SG	2.47	0.54
1:D:39:GLY:HA3	1:D:260:VAL:HG23	1.89	0.54
1:E:262:GLY:HA3	1:E:315:SER:HB2	1.90	0.54
1:B:191:LYS:HE2	1:B:195:MET:SD	2.47	0.54
1:E:51:VAL:HG13	1:E:310:LEU:CD1	2.35	0.54
1:B:139:ILE:O	1:B:140:ASN:HB3	2.06	0.54
1:D:114:GLU:HG3	5:D:729:HOH:O	2.07	0.54
1:E:36:VAL:HB	1:E:257:ASP:OD2	2.08	0.54
1:E:90:GLU:CD	1:E:90:GLU:H	2.15	0.54
1:E:172:LEU:HD22	1:E:213:ILE:HG22	1.90	0.54
1:C:142:SER:OG	1:C:145:GLU:HG3	2.07	0.53
1:E:84:GLN:HE21	1:E:84:GLN:N	2.03	0.53
1:A:107:THR:HG21	1:A:113:PHE:HA	1.90	0.53
1:B:174:CYS:SG	1:B:175:PRO:HD2	2.49	0.53
1:F:51:VAL:HG22	5:F:770:HOH:O	2.08	0.53
1:D:75:SER:HB3	1:D:85:MET:HG3	1.90	0.53
1:F:37:GLU:HG2	1:F:63:LEU:HD13	1.90	0.53
1:E:135:THR:HG22	1:E:141:CYS:O	2.07	0.53
1:F:76:PHE:CZ	1:F:116:ALA:HA	2.44	0.53
1:A:297:GLN:N	1:A:297:GLN:NE2	2.56	0.53
1:B:91:VAL:O	1:B:95:ILE:HG23	2.08	0.53
1:E:145:GLU:HA	1:E:148:GLN:OE1	2.09	0.53
1:B:126:ILE:O	1:B:166:GLY:HA2	2.07	0.53
1:D:122:LYS:O	1:D:162:ILE:HG23	2.08	0.53
1:D:230:LEU:O	1:D:255:VAL:HG13	2.08	0.53
1:A:87:ASP:O	1:A:91:VAL:HG23	2.09	0.53
1:C:191:LYS:HE2	1:C:195:MET:CE	2.37	0.53
1:C:259:SER:OG	1:C:263:LEU:HB2	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:114:GLU:OE2	1:C:114:GLU:N	2.42	0.53
1:D:279:GLU:HG3	1:D:299:LEU:HD13	1.91	0.53
1:E:206:ILE:HG13	1:E:208:VAL:HG13	1.91	0.53
1:C:47:GLU:HG2	1:C:310:LEU:CD2	2.38	0.52
1:D:136:LYS:O	1:D:140:ASN:HA	2.08	0.52
1:B:189:THR:HG22	1:B:225:VAL:HG21	1.91	0.52
1:B:60:ILE:HD13	1:B:73:THR:HG23	1.91	0.52
1:D:145:GLU:HA	1:D:148:GLN:HG2	1.91	0.52
1:D:236:ASP:OD1	1:D:239:GLY:HA2	2.09	0.52
1:E:58:LYS:O	1:E:62:MET:HG3	2.09	0.52
1:C:247:MET:O	1:C:251:MET:HG2	2.08	0.52
1:F:247:MET:HE3	1:F:250:GLN:HE21	1.72	0.52
1:F:127:PHE:HZ	1:F:139:ILE:HD11	1.73	0.52
1:B:58:LYS:HZ2	1:B:62:MET:HE2	1.71	0.52
1:E:77:VAL:HG11	1:E:82:VAL:HG11	1.92	0.52
1:E:199:GLU:C	1:E:200:ILE:HD12	2.35	0.52
1:F:78:SER:HB2	1:F:80:LYS:CG	2.39	0.52
1:C:133:LEU:HD13	1:C:177:GLU:OE1	2.10	0.52
1:E:259:SER:HB2	1:E:275:ASN:HB2	1.91	0.52
1:F:131:SER:C	1:F:143:ILE:HD11	2.35	0.52
1:C:172:LEU:HD12	1:C:204:ASP:HB2	1.91	0.52
1:F:227:LEU:HD21	1:F:252:GLY:HA3	1.92	0.52
1:A:113:PHE:O	1:A:117:VAL:HG23	2.10	0.52
1:B:231:ALA:CB	1:B:255:VAL:HG22	2.40	0.52
1:B:266:CYS:SG	1:B:269:ALA:HB3	2.50	0.52
1:D:139:ILE:HG23	1:D:149:ARG:HH12	1.75	0.52
1:A:259:SER:HB2	1:A:275:ASN:CB	2.40	0.51
1:C:210:THR:HG21	1:D:287:GLY:HA3	1.91	0.51
1:F:142:SER:HB3	1:F:145:GLU:OE1	2.10	0.51
1:A:141:CYS:HB2	1:A:145:GLU:HB3	1.93	0.51
1:B:260:VAL:O	1:B:279:GLU:OE1	2.28	0.51
1:E:211:PRO:HG2	1:F:288:LEU:CD1	2.29	0.51
1:F:151:ASP:OD1	1:F:155:LYS:HE3	2.10	0.51
1:A:139:ILE:HB	1:A:141:CYS:SG	2.50	0.51
1:A:288:LEU:CD1	1:B:211:PRO:HG2	2.38	0.51
1:C:127:PHE:HD1	1:C:127:PHE:H	1.57	0.51
1:F:91:VAL:O	1:F:95:ILE:HG23	2.10	0.51
1:D:73:THR:OG1	1:D:103:TYR:HB3	2.10	0.51
1:B:147:PHE:CZ	1:B:191:LYS:HG2	2.46	0.51
1:C:122:LYS:NZ	1:C:122:LYS:HB3	2.25	0.51
1:C:117:VAL:HG22	1:C:162:ILE:HD11	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:75:SER:O	1:F:77:VAL:N	2.44	0.51
1:F:198:TYR:HD1	1:F:198:TYR:H	1.59	0.51
1:C:147:PHE:HB3	1:C:195:MET:HE1	1.92	0.51
1:E:48:LYS:HD2	1:E:48:LYS:N	2.25	0.51
1:A:35:ILE:CD1	1:A:282:VAL:HG12	2.41	0.51
1:A:65:GLU:HG3	1:A:98:PHE:HZ	1.75	0.51
1:D:143:ILE:HG23	1:D:144:GLU:H	1.76	0.51
1:A:62:MET:HE3	1:A:305:PHE:CD2	2.46	0.51
1:A:182:PRO:HD2	5:A:575:HOH:O	2.11	0.51
1:E:200:ILE:HD12	1:E:200:ILE:N	2.26	0.51
1:A:133:LEU:HD13	1:A:177:GLU:HG2	1.93	0.51
1:D:189:THR:HG22	1:D:225:VAL:HG21	1.93	0.51
1:E:60:ILE:HD13	1:E:73:THR:HG23	1.93	0.51
1:A:191:LYS:O	1:A:195:MET:HB2	2.10	0.50
1:B:32:ARG:NH1	1:B:291:HIS:CD2	2.80	0.50
1:E:58:LYS:HD3	1:E:305:PHE:CE1	2.47	0.50
1:F:266:CYS:SG	1:F:269:ALA:HB3	2.52	0.50
1:A:297:GLN:CD	1:A:297:GLN:N	2.68	0.50
1:E:123:GLU:HG3	1:E:163:SER:HB2	1.94	0.50
1:A:281:LEU:O	1:A:285:LEU:HG	2.12	0.50
1:B:231:ALA:HB2	1:B:255:VAL:HG22	1.93	0.50
1:E:164:VAL:O	1:E:197:CYS:HA	2.12	0.50
1:D:52:SER:HB2	1:D:54:PRO:HD2	1.94	0.50
1:B:312:ARG:HD3	1:B:313:LYS:O	2.12	0.50
1:C:33:VAL:HG23	1:C:290:ILE:CG2	2.39	0.50
1:C:154:LEU:O	1:C:158:GLN:HG3	2.12	0.49
1:E:266:CYS:HB3	1:E:269:ALA:HB3	1.93	0.49
1:F:251:MET:HA	1:F:251:MET:HE2	1.93	0.49
1:B:173:GLY:HA2	1:B:180:ILE:HG12	1.93	0.49
1:F:33:VAL:HG23	1:F:290:ILE:CG2	2.41	0.49
1:B:39:GLY:HA2	1:B:259:SER:OG	2.12	0.49
1:B:165:ARG:HD2	1:B:199:GLU:OE2	2.13	0.49
1:D:260:VAL:C	1:D:262:GLY:H	2.19	0.49
1:B:39:GLY:HA3	1:B:260:VAL:HG23	1.94	0.49
1:D:133:LEU:HD13	1:D:177:GLU:HG2	1.95	0.49
1:D:150:PHE:O	1:D:154:LEU:HB2	2.11	0.49
1:D:174:CYS:SG	1:D:175:PRO:HD2	2.52	0.49
1:F:109:ASN:OD1	1:F:111:LYS:HB3	2.11	0.49
1:B:37:GLU:HG2	1:B:63:LEU:HD13	1.93	0.49
1:D:51:VAL:O	1:D:56:LYS:NZ	2.45	0.49
1:E:47:GLU:HG2	1:E:310:LEU:HD21	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:259:SER:OG	1:F:263:LEU:HB2	2.12	0.49
1:A:82:VAL:HG22	1:A:82:VAL:O	2.11	0.49
1:C:211:PRO:CG	1:D:288:LEU:HD21	2.43	0.49
1:D:143:ILE:HG23	1:D:144:GLU:N	2.26	0.49
1:D:199:GLU:C	1:D:200:ILE:HD12	2.37	0.49
1:E:154:LEU:HD11	1:E:195:MET:HB3	1.91	0.49
1:A:84:GLN:CD	1:A:84:GLN:H	2.21	0.49
1:D:251:MET:HA	1:D:251:MET:HE2	1.95	0.49
1:E:82:VAL:HA	1:E:84:GLN:HE22	1.78	0.49
1:A:58:LYS:HD3	1:A:305:PHE:CE1	2.48	0.48
1:C:302:ALA:O	1:C:306:ILE:HG22	2.13	0.48
1:F:165:ARG:NH1	1:F:201:SER:HB2	2.28	0.48
1:F:279:GLU:HA	1:F:282:VAL:CG1	2.43	0.48
1:B:124:VAL:CG2	1:B:164:VAL:HG22	2.43	0.48
1:D:214:MET:HE2	1:D:244:ASN:HB3	1.94	0.48
1:E:139:ILE:HD12	1:E:141:CYS:SG	2.53	0.48
1:E:267:PRO:C	1:E:269:ALA:H	2.20	0.48
1:D:208:VAL:HG12	1:D:238:TYR:CE2	2.48	0.48
1:F:269:ALA:O	1:F:270:GLN:O	2.31	0.48
1:B:170:CYS:HA	1:B:173:GLY:O	2.14	0.48
1:D:35:ILE:CD1	1:D:282:VAL:HG12	2.44	0.48
1:E:53:THR:N	1:E:54:PRO:CD	2.76	0.48
1:F:58:LYS:HG2	1:F:62:MET:HE2	1.96	0.48
1:C:138:ASN:HB2	1:C:139:ILE:HD13	1.94	0.48
1:C:247:MET:HE3	1:C:250:GLN:HE21	1.78	0.48
1:A:202:LEU:HD13	1:A:217:MET:SD	2.54	0.48
1:B:133:LEU:HG	1:B:137:LYS:NZ	2.29	0.48
1:D:83:PRO:HG2	1:D:84:GLN:HE22	1.77	0.48
1:B:135:THR:HG23	1:B:139:ILE:HD11	1.95	0.48
1:C:139:ILE:HB	1:C:149:ARG:NH2	2.28	0.48
1:C:172:LEU:HD21	1:C:214:MET:HA	1.95	0.48
1:F:73:THR:OG1	1:F:103:TYR:HB3	2.14	0.48
1:C:73:THR:OG1	1:C:103:TYR:HB3	2.14	0.48
1:C:222:MET:HG2	1:C:227:LEU:HD13	1.95	0.48
1:F:106:LEU:N	1:F:106:LEU:CD1	2.74	0.48
1:A:279:GLU:HG3	1:A:299:LEU:HD13	1.96	0.48
1:A:306:ILE:O	1:A:310:LEU:HB2	2.12	0.48
1:C:98:PHE:HB2	1:C:101:ILE:HD12	1.96	0.48
1:D:205:THR:HA	1:D:235:HIS:CD2	2.49	0.48
1:D:248:ALA:O	1:D:251:MET:HB2	2.14	0.48
1:E:169:SER:O	1:E:170:CYS:HB2	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:51:VAL:HG13	1:B:310:LEU:CD1	2.44	0.47
1:D:133:LEU:HD22	1:D:177:GLU:HG2	1.95	0.47
1:E:288:LEU:HD11	1:F:212:GLY:H	1.78	0.47
1:F:222:MET:HG2	1:F:227:LEU:HD13	1.96	0.47
1:D:126:ILE:HD11	1:D:197:CYS:SG	2.54	0.47
1:D:306:ILE:O	1:D:310:LEU:HB2	2.14	0.47
1:F:265:GLY:N	1:F:274:GLY:HA3	2.29	0.47
1:B:59:LEU:O	1:B:63:LEU:HG	2.14	0.47
1:B:200:ILE:N	1:B:200:ILE:HD12	2.29	0.47
1:C:141:CYS:HB2	1:C:145:GLU:HB2	1.96	0.47
1:C:169:SER:HB2	1:C:205:THR:OG1	2.13	0.47
1:D:85:MET:O	1:D:88:HIS:HD2	1.97	0.47
1:F:232:VAL:HG22	1:F:234:CYS:SG	2.54	0.47
1:D:134:PHE:CD1	1:D:134:PHE:C	2.92	0.47
1:E:136:LYS:NZ	1:E:142:SER:HB3	2.29	0.47
1:E:218:LEU:O	1:E:219:SER:C	2.58	0.47
1:D:132:GLU:HG2	1:D:143:ILE:HB	1.97	0.47
1:F:172:LEU:HD11	1:F:214:MET:HE2	1.96	0.47
1:A:28:THR:HA	5:A:864:HOH:O	2.15	0.47
1:B:270:GLN:O	1:B:272:ALA:N	2.47	0.47
1:D:174:CYS:N	1:D:178:GLY:O	2.46	0.47
1:F:41:ARG:HD2	1:F:41:ARG:C	2.39	0.47
1:F:82:VAL:HG22	1:F:82:VAL:O	2.14	0.47
1:F:211:PRO:N	1:F:244:ASN:HD21	2.13	0.47
1:A:41:ARG:HD2	1:A:41:ARG:C	2.39	0.47
1:A:60:ILE:HD13	1:A:73:THR:HG23	1.96	0.47
1:A:141:CYS:HB2	1:A:145:GLU:HB2	1.95	0.47
1:A:287:GLY:HA3	1:B:210:THR:HG21	1.97	0.47
1:B:52:SER:HG	1:B:55:VAL:HG23	1.80	0.47
1:C:40:PRO:HA	1:C:44:LEU:HD22	1.96	0.47
1:C:190:LYS:HD3	1:C:224:GLU:HB3	1.96	0.47
1:C:306:ILE:HD11	1:C:310:LEU:HD22	1.96	0.47
1:D:33:VAL:CG2	1:D:290:ILE:HG21	2.44	0.47
1:C:206:ILE:HG13	1:C:208:VAL:HG13	1.97	0.47
1:D:247:MET:HE3	1:D:247:MET:O	2.15	0.47
1:D:284:MET:O	1:D:288:LEU:HD23	2.14	0.47
1:F:63:LEU:HB2	1:F:71:ILE:HD13	1.97	0.47
1:F:138:ASN:ND2	1:F:139:ILE:H	2.12	0.47
1:F:170:CYS:HA	1:F:173:GLY:O	2.15	0.47
1:A:231:ALA:CB	1:A:255:VAL:HG22	2.45	0.47
1:C:71:ILE:O	1:C:104:PRO:HD2	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:181:SER:HB2	5:C:1017:HOH:O	2.15	0.47
1:F:310:LEU:N	1:F:310:LEU:HD12	2.30	0.47
1:B:60:ILE:CD1	1:B:73:THR:HG23	2.45	0.47
1:C:48:LYS:H	1:C:48:LYS:CD	2.19	0.47
1:C:77:VAL:HG13	1:C:82:VAL:HG13	1.95	0.47
1:C:133:LEU:N	1:C:133:LEU:HD12	2.29	0.47
1:A:322:THR:HG23	5:A:728:HOH:O	2.14	0.46
1:B:126:ILE:HD11	1:B:197:CYS:SG	2.55	0.46
1:B:282:VAL:CG2	1:B:296:LEU:HD13	2.45	0.46
1:D:259:SER:CB	1:D:275:ASN:HB2	2.45	0.46
1:E:190:LYS:HD3	1:E:224:GLU:HB3	1.96	0.46
1:F:164:VAL:O	1:F:197:CYS:HA	2.15	0.46
1:F:203:GLY:HA2	1:F:233:HIS:O	2.15	0.46
1:A:288:LEU:HB3	1:A:290:ILE:HG13	1.96	0.46
1:C:53:THR:O	1:C:57:ILE:HG13	2.15	0.46
1:C:189:THR:HG22	1:C:225:VAL:HG21	1.98	0.46
1:F:126:ILE:HD11	1:F:197:CYS:SG	2.56	0.46
1:F:247:MET:CE	1:F:250:GLN:HE21	2.29	0.46
1:C:77:VAL:CG1	1:C:82:VAL:HG11	2.46	0.46
1:D:41:ARG:HG3	1:D:106:LEU:HD13	1.98	0.46
1:D:45:GLN:HB2	1:D:85:MET:HE3	1.96	0.46
1:F:90:GLU:CD	1:F:90:GLU:N	2.70	0.46
1:C:77:VAL:CG1	1:C:82:VAL:CG1	2.94	0.46
1:C:200:ILE:N	1:C:200:ILE:HD12	2.30	0.46
1:F:169:SER:O	1:F:170:CYS:HB2	2.14	0.46
1:F:295:ASN:OD1	1:F:298:LYS:HG3	2.16	0.46
1:B:68:LEU:H	1:B:68:LEU:HD22	1.80	0.46
1:C:247:MET:CE	1:C:250:GLN:HE21	2.29	0.46
1:B:199:GLU:C	1:B:200:ILE:HD12	2.41	0.46
1:F:81:TRP:O	1:F:83:PRO:CD	2.63	0.46
1:F:115:ALA:O	1:F:118:ALA:HB3	2.15	0.46
1:F:175:PRO:HD3	1:F:206:ILE:HG21	1.98	0.46
4:B:399:HGG:H19	4:B:399:HGG:O36	2.15	0.46
1:D:122:LYS:C	1:D:162:ILE:HG23	2.41	0.46
1:F:52:SER:O	1:F:55:VAL:CG1	2.62	0.46
1:F:95:ILE:HD12	1:F:103:TYR:CD1	2.51	0.46
1:F:127:PHE:CZ	1:F:139:ILE:HD11	2.50	0.46
1:A:125:VAL:HG22	1:A:165:ARG:HB3	1.98	0.46
1:B:150:PHE:C	1:B:154:LEU:HD12	2.40	0.46
1:C:46:ASN:HB3	5:C:682:HOH:O	2.16	0.46
1:C:53:THR:N	1:C:54:PRO:HD2	2.31	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:172:LEU:HD22	1:D:213:ILE:HG22	1.99	0.45
1:D:256:VAL:CG2	1:D:281:LEU:HD21	2.46	0.45
1:D:263:LEU:HD23	1:D:312:ARG:NH1	2.31	0.45
1:E:51:VAL:CG1	1:E:310:LEU:HD13	2.40	0.45
1:F:92:LEU:HD23	1:F:119:ALA:HB3	1.97	0.45
1:B:58:LYS:HD3	1:B:305:PHE:CZ	2.51	0.45
1:C:36:VAL:HG13	1:C:70:VAL:HG22	1.97	0.45
1:D:144:GLU:O	1:D:148:GLN:HG2	2.15	0.45
1:B:41:ARG:HD3	1:B:106:LEU:HD22	1.98	0.45
1:C:260:VAL:O	1:C:279:GLU:OE1	2.35	0.45
1:C:210:THR:HG23	5:C:818:HOH:O	2.17	0.45
1:C:268:TYR:CD1	1:C:268:TYR:N	2.85	0.45
1:D:40:PRO:HA	1:D:44:LEU:HD22	1.99	0.45
1:E:77:VAL:CG1	1:E:82:VAL:HG11	2.46	0.45
1:F:165:ARG:HH12	1:F:257:ASP:CG	2.24	0.45
1:C:61:ASP:CG	1:C:96:GLN:H	2.24	0.45
1:C:306:ILE:O	1:C:310:LEU:HB2	2.17	0.45
1:D:40:PRO:HG2	1:D:72:GLU:O	2.16	0.45
1:E:172:LEU:HD12	1:E:204:ASP:HB2	1.97	0.45
1:F:77:VAL:HG22	1:F:78:SER:N	2.31	0.45
1:C:117:VAL:HG22	1:C:162:ILE:CD1	2.47	0.45
1:F:132:GLU:CD	1:F:143:ILE:HD12	2.41	0.45
1:F:168:VAL:HG23	1:F:200:ILE:CG2	2.47	0.45
1:A:282:VAL:HG21	1:A:296:LEU:HD13	1.99	0.45
1:C:123:GLU:HG3	1:C:163:SER:HB2	1.98	0.45
1:F:40:PRO:HA	1:F:44:LEU:HD22	1.99	0.45
1:F:230:LEU:N	1:F:230:LEU:HD22	2.31	0.45
1:A:154:LEU:HD13	1:A:195:MET:HG2	1.99	0.45
1:A:227:LEU:HA	1:A:230:LEU:HD22	1.99	0.45
1:C:92:LEU:HD23	1:C:119:ALA:HB3	1.99	0.45
1:D:145:GLU:OE1	1:D:145:GLU:N	2.49	0.44
1:D:282:VAL:CG2	1:D:296:LEU:HD13	2.47	0.44
1:E:96:GLN:HB2	1:E:98:PHE:HE1	1.82	0.44
1:E:135:THR:O	1:E:139:ILE:HG13	2.17	0.44
1:F:266:CYS:HA	1:F:267:PRO:HD3	1.87	0.44
1:A:136:LYS:HE3	1:A:142:SER:HA	2.00	0.44
1:A:154:LEU:HD22	1:A:196:GLY:HA3	1.99	0.44
1:C:247:MET:HE3	1:C:247:MET:HA	1.98	0.44
1:D:52:SER:C	1:D:54:PRO:HD2	2.43	0.44
1:D:246:LEU:C	1:D:246:LEU:CD1	2.88	0.44
1:E:210:THR:C	1:E:244:ASN:HD21	2.25	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:297:GLN:HE21	1:A:297:GLN:CA	2.29	0.44
1:B:154:LEU:O	1:B:158:GLN:HG3	2.17	0.44
1:C:37:GLU:OE2	1:C:260:VAL:N	2.51	0.44
1:F:165:ARG:NH1	1:F:257:ASP:OD1	2.50	0.44
1:A:148:GLN:O	1:A:151:ASP:HB3	2.18	0.44
1:B:144:GLU:OE1	1:B:147:PHE:CD1	2.67	0.44
1:C:37:GLU:CG	1:C:63:LEU:HD13	2.47	0.44
1:D:78:SER:OG	1:D:80:LYS:HG2	2.18	0.44
1:D:87:ASP:O	1:D:91:VAL:HG23	2.18	0.44
1:D:206:ILE:HG13	1:D:208:VAL:HG22	1.98	0.44
1:D:113:PHE:CD2	1:D:124:VAL:HG11	2.53	0.44
1:D:200:ILE:N	1:D:200:ILE:CD1	2.81	0.44
1:E:84:GLN:H	1:E:84:GLN:CD	2.18	0.44
1:F:106:LEU:O	1:F:107:THR:HG23	2.17	0.44
1:A:165:ARG:HH12	1:A:257:ASP:CG	2.26	0.44
1:D:154:LEU:HD23	1:D:164:VAL:HG21	1.99	0.44
1:D:191:LYS:HG3	1:D:195:MET:HE3	2.00	0.44
1:D:259:SER:O	1:D:260:VAL:O	2.36	0.44
1:F:133:LEU:HD22	1:F:177:GLU:HG2	1.98	0.44
1:F:237:THR:HG23	1:F:272:ALA:HB3	2.00	0.44
1:F:306:ILE:O	1:F:309:ALA:HB3	2.17	0.44
1:E:73:THR:OG1	1:E:103:TYR:HB3	2.17	0.44
1:E:108:PRO:HG2	1:E:109:ASN:H	1.83	0.44
1:A:265:GLY:HA2	1:A:274:GLY:CA	2.47	0.43
1:B:44:LEU:HD12	1:B:44:LEU:HA	1.86	0.43
1:D:259:SER:HB3	1:D:275:ASN:HB2	2.00	0.43
1:E:77:VAL:HG13	1:E:82:VAL:HG13	2.00	0.43
1:F:230:LEU:O	1:F:255:VAL:CG1	2.66	0.43
1:F:310:LEU:N	1:F:310:LEU:CD1	2.81	0.43
1:B:208:VAL:HG12	1:B:238:TYR:CE2	2.52	0.43
1:D:140:ASN:HD22	1:D:141:CYS:N	2.16	0.43
1:D:210:THR:CB	1:D:211:PRO:HD2	2.46	0.43
1:E:139:ILE:C	1:E:141:CYS:H	2.25	0.43
1:F:61:ASP:CG	1:F:96:GLN:HE21	2.24	0.43
1:F:133:LEU:HD12	1:F:133:LEU:N	2.33	0.43
1:F:139:ILE:HG22	1:F:149:ARG:HH22	1.83	0.43
1:F:291:HIS:HB2	5:F:608:HOH:O	2.18	0.43
1:F:322:THR:O	1:F:323:CYS:HB2	2.17	0.43
1:A:241:ALA:HB1	1:A:276:LEU:HD13	2.00	0.43
1:C:191:LYS:O	1:C:195:MET:HB2	2.18	0.43
1:D:279:GLU:HG2	1:D:300:LEU:HD23	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:77:VAL:HG13	1:E:82:VAL:CG1	2.48	0.43
1:C:237:THR:O	1:D:317:LYS:HB3	2.18	0.43
1:D:48:LYS:O	1:D:49:ASN:C	2.60	0.43
1:D:165:ARG:NH1	1:D:201:SER:HB2	2.33	0.43
1:A:90:GLU:OE1	1:A:90:GLU:N	2.42	0.43
1:A:106:LEU:HD12	1:A:125:VAL:HB	1.99	0.43
1:A:141:CYS:SG	1:A:149:ARG:NH2	2.92	0.43
1:B:109:ASN:OD1	4:B:399:HGG:N7	2.50	0.43
1:D:32:ARG:HD3	5:D:819:HOH:O	2.18	0.43
1:B:36:VAL:HA	1:B:70:VAL:HG13	2.00	0.43
1:D:278:THR:O	1:D:282:VAL:HG13	2.18	0.43
1:A:51:VAL:CG1	1:A:310:LEU:HD13	2.49	0.43
1:C:36:VAL:HG22	1:C:70:VAL:HG11	2.00	0.43
1:C:114:GLU:OE2	1:C:114:GLU:CA	2.66	0.43
1:A:133:LEU:HD13	1:A:177:GLU:CG	2.49	0.43
1:A:182:PRO:HG3	5:A:791:HOH:O	2.18	0.43
1:B:83:PRO:HG2	1:B:84:GLN:NE2	2.33	0.43
1:C:133:LEU:O	1:C:137:LYS:CG	2.67	0.43
1:F:41:ARG:HB2	1:F:72:GLU:O	2.19	0.43
1:F:41:ARG:HG3	1:F:106:LEU:HD13	2.01	0.43
1:A:147:PHE:HD1	1:A:191:LYS:HZ3	1.66	0.43
1:B:136:LYS:O	1:B:140:ASN:HA	2.19	0.43
1:C:133:LEU:HD13	1:C:177:GLU:CD	2.44	0.43
1:D:80:LYS:HD3	1:D:80:LYS:N	2.34	0.43
1:F:106:LEU:HA	1:F:125:VAL:HB	2.01	0.43
1:F:154:LEU:HD22	1:F:196:GLY:HA3	2.01	0.43
1:B:68:LEU:HD22	1:B:68:LEU:N	2.34	0.43
1:E:267:PRO:HD2	5:E:685:HOH:O	2.17	0.43
1:F:138:ASN:O	1:F:139:ILE:C	2.62	0.43
1:A:212:GLY:H	1:B:288:LEU:CD1	2.32	0.42
1:B:32:ARG:NH2	5:B:649:HOH:O	2.50	0.42
1:D:33:VAL:HG23	1:D:290:ILE:CG2	2.46	0.42
1:E:122:LYS:C	1:E:162:ILE:HG23	2.44	0.42
1:E:270:GLN:C	1:E:272:ALA:H	2.27	0.42
1:F:182:PRO:O	1:F:183:ALA:C	2.62	0.42
1:D:279:GLU:O	1:D:282:VAL:HG22	2.20	0.42
1:E:82:VAL:HG23	1:E:84:GLN:OE1	2.19	0.42
1:E:127:PHE:O	1:E:150:PHE:HZ	2.02	0.42
1:A:169:SER:O	1:A:170:CYS:HB2	2.20	0.42
1:A:211:PRO:HD3	1:A:244:ASN:ND2	2.33	0.42
1:D:143:ILE:HD11	1:D:184:LYS:HZ1	1.84	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:109:ASN:C	1:C:153:ILE:HD11	2.44	0.42
1:F:140:ASN:O	1:F:141:CYS:HB3	2.19	0.42
1:A:139:ILE:O	1:A:141:CYS:N	2.52	0.42
1:B:124:VAL:HG23	1:B:164:VAL:HG22	2.00	0.42
1:C:51:VAL:HG13	1:C:310:LEU:HD13	2.01	0.42
1:D:206:ILE:HG13	1:D:208:VAL:HG13	2.01	0.42
1:A:234:CYS:HB3	5:A:523:HOH:O	2.18	0.42
1:B:78:SER:OG	1:B:80:LYS:HB2	2.20	0.42
1:B:251:MET:HA	1:B:251:MET:CE	2.42	0.42
1:D:218:LEU:HD12	1:D:251:MET:HG3	2.00	0.42
1:F:191:LYS:O	1:F:195:MET:HG3	2.19	0.42
1:C:82:VAL:O	1:C:82:VAL:HG22	2.20	0.42
1:D:322:THR:HG22	1:D:323:CYS:N	2.34	0.42
1:E:219:SER:O	1:E:223:GLN:HG2	2.20	0.42
1:E:145:GLU:HA	1:E:148:GLN:CD	2.45	0.42
1:F:220:ALA:O	1:F:221:VAL:C	2.62	0.42
1:B:53:THR:N	1:B:54:PRO:CD	2.81	0.42
1:D:59:LEU:HD13	1:D:306:ILE:HG23	2.02	0.42
1:D:127:PHE:CE2	1:D:139:ILE:HD11	2.55	0.42
1:D:311:ASN:OD1	1:D:311:ASN:O	2.38	0.42
1:F:261:ALA:O	1:F:315:SER:HB2	2.20	0.42
1:A:165:ARG:NH2	1:A:199:GLU:OE2	2.52	0.42
1:A:232:VAL:HG13	1:A:234:CYS:SG	2.59	0.42
1:C:83:PRO:CD	1:C:84:GLN:HE22	2.28	0.42
1:A:279:GLU:HB2	1:A:316:SER:OG	2.20	0.41
1:B:246:LEU:O	1:B:246:LEU:HD22	2.20	0.41
1:F:57:ILE:HG23	1:F:95:ILE:HA	2.02	0.41
1:C:316:SER:O	1:C:320:GLN:HG3	2.20	0.41
1:B:84:GLN:O	1:B:85:MET:HE2	2.20	0.41
1:C:206:ILE:H	1:C:206:ILE:HG12	1.61	0.41
1:D:143:ILE:CD1	1:D:184:LYS:NZ	2.84	0.41
1:C:210:THR:HB	1:C:211:PRO:HD2	2.02	0.41
1:D:36:VAL:HG13	1:D:70:VAL:CG2	2.49	0.41
1:F:134:PHE:HA	1:F:176:TYR:HD1	1.85	0.41
1:F:196:GLY:O	1:F:197:CYS:C	2.63	0.41
4:B:399:HGG:H57	4:B:399:HGG:O34	2.21	0.41
1:D:78:SER:HA	1:D:79:PRO:HD3	1.90	0.41
1:D:235:HIS:CE1	1:D:275:ASN:OD1	2.74	0.41
1:E:140:ASN:HD22	1:E:140:ASN:N	2.18	0.41
1:F:316:SER:HB3	1:F:319:ALA:CB	2.50	0.41
1:A:265:GLY:N	1:A:274:GLY:HA3	2.36	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:134:PHE:HB2	1:D:174:CYS:SG	2.60	0.41
1:D:260:VAL:O	1:D:262:GLY:N	2.53	0.41
1:E:78:SER:HA	1:E:79:PRO:HD3	1.99	0.41
1:F:133:LEU:HD13	1:F:177:GLU:OE1	2.20	0.41
1:A:304:ASN:O	1:A:308:GLN:HG3	2.21	0.41
1:C:87:ASP:O	1:C:90:GLU:HG2	2.21	0.41
1:C:133:LEU:O	1:C:137:LYS:HG2	2.19	0.41
1:F:53:THR:N	1:F:54:PRO:CD	2.80	0.41
1:D:181:SER:HA	1:D:182:PRO:HD3	1.93	0.41
1:E:68:LEU:H	1:E:68:LEU:HD22	1.85	0.41
1:F:42:ASP:OD1	1:F:233:HIS:NE2	2.53	0.41
1:F:244:ASN:HD22	1:F:244:ASN:HA	1.69	0.41
1:F:316:SER:HB3	1:F:319:ALA:HB3	2.03	0.41
1:A:186:ALA:HA	1:A:221:VAL:HG22	2.02	0.41
1:A:201:SER:HA	1:A:231:ALA:HB3	2.03	0.41
1:B:28:THR:HG22	1:B:29:LEU:HG	2.02	0.41
1:C:61:ASP:HA	1:C:64:SER:HB2	2.03	0.41
1:C:133:LEU:HD12	1:C:133:LEU:H	1.86	0.41
1:C:279:GLU:HG2	1:C:300:LEU:HD23	2.02	0.41
1:D:246:LEU:HA	1:D:249:LEU:HD12	2.02	0.41
1:F:98:PHE:HD1	1:F:103:TYR:OH	2.03	0.41
1:B:45:GLN:HB2	1:B:85:MET:HE3	2.03	0.41
1:D:47:GLU:OE1	1:D:47:GLU:HA	2.21	0.41
1:E:39:GLY:N	1:E:40:PRO:HD2	2.35	0.41
1:E:145:GLU:HG2	5:E:1022:HOH:O	2.20	0.41
1:B:39:GLY:N	1:B:40:PRO:HD2	2.35	0.40
1:C:77:VAL:CG2	1:C:78:SER:N	2.83	0.40
1:C:281:LEU:HD12	1:C:281:LEU:O	2.21	0.40
1:D:141:CYS:SG	1:D:149:ARG:NH2	2.94	0.40
1:E:231:ALA:HB2	1:E:255:VAL:HG22	2.03	0.40
1:F:106:LEU:C	1:F:107:THR:HG23	2.46	0.40
1:F:133:LEU:HD13	1:F:177:GLU:CD	2.45	0.40
1:B:112:GLY:HA3	4:B:399:HGG:HN6A	1.84	0.40
1:E:203:GLY:HA2	1:E:233:HIS:O	2.22	0.40
1:D:98:PHE:CB	1:D:101:ILE:HD12	2.51	0.40
1:D:99:PRO:HD2	5:D:1010:HOH:O	2.20	0.40
1:D:304:ASN:O	1:D:308:GLN:HG3	2.22	0.40
1:B:37:GLU:CG	1:B:63:LEU:HD13	2.51	0.40
1:C:86:GLY:HA2	5:C:554:HOH:O	2.20	0.40
5:C:869:HOH:O	1:E:34:LYS:HE2	2.20	0.40
1:D:214:MET:CE	1:D:244:ASN:HB3	2.51	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:136:LYS:HZ2	1:E:142:SER:HB3	1.84	0.40
1:B:149:ARG:O	1:B:153:ILE:HG13	2.21	0.40
1:C:61:ASP:O	1:C:65:GLU:HG3	2.21	0.40
1:D:282:VAL:HG21	1:D:296:LEU:HD13	2.03	0.40
1:D:294:VAL:HG12	1:D:295:ASN:N	2.36	0.40
1:E:29:LEU:HD23	1:E:250:GLN:HA	2.04	0.40
1:E:133:LEU:O	1:E:137:LYS:HG3	2.21	0.40
1:E:231:ALA:HB1	1:E:255:VAL:HG22	2.00	0.40
1:F:55:VAL:HG23	1:F:305:PHE:HE2	1.81	0.40
1:F:77:VAL:CG2	1:F:78:SER:N	2.84	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	294/298 (99%)	272 (92%)	16 (5%)	6 (2%)	6	7
1	B	294/298 (99%)	272 (92%)	16 (5%)	6 (2%)	6	7
1	C	294/298 (99%)	263 (90%)	27 (9%)	4 (1%)	9	13
1	D	285/298 (96%)	254 (89%)	25 (9%)	6 (2%)	5	7
1	E	294/298 (99%)	273 (93%)	15 (5%)	6 (2%)	6	7
1	F	294/298 (99%)	249 (85%)	35 (12%)	10 (3%)	3	2
All	All	1755/1788 (98%)	1583 (90%)	134 (8%)	38 (2%)	5	6

All (38) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	139	ILE
1	A	140	ASN

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Mol	Chain	Res	Type
1	A	260	VAL
1	A	270	GLN
1	A	272	ALA
1	B	260	VAL
1	B	273	SER
1	C	140	ASN
1	C	271	GLY
1	D	260	VAL
1	E	139	ILE
1	E	260	VAL
1	F	73	THR
1	F	76	PHE
1	F	260	VAL
1	F	270	GLN
1	B	108	PRO
1	B	160	ALA
1	B	271	GLY
1	D	138	ASN
1	D	141	CYS
1	F	48	LYS
1	F	86	GLY
1	F	139	ILE
1	F	140	ASN
1	F	141	CYS
1	A	271	GLY
1	C	260	VAL
1	D	137	LYS
1	E	272	ALA
1	D	261	ALA
1	E	143	ILE
1	B	270	GLN
1	C	142	SER
1	E	268	TYR
1	F	305	PHE
1	D	273	SER
1	E	108	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	238/240 (99%)	220 (92%)	18 (8%)	12	21
1	B	238/240 (99%)	221 (93%)	17 (7%)	13	23
1	C	238/240 (99%)	220 (92%)	18 (8%)	12	21
1	D	234/240 (98%)	215 (92%)	19 (8%)	11	18
1	E	238/240 (99%)	216 (91%)	22 (9%)	8	14
1	F	238/240 (99%)	222 (93%)	16 (7%)	15	26
All	All	1424/1440 (99%)	1314 (92%)	110 (8%)	12	20

All (110) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	32	ARG
1	A	44	LEU
1	A	82	VAL
1	A	114	GLU
1	A	139	ILE
1	A	161	ASN
1	A	165	ARG
1	A	230	LEU
1	A	232	VAL
1	A	246	LEU
1	A	255	VAL
1	A	256	VAL
1	A	259	SER
1	A	260	VAL
1	A	288	LEU
1	A	297	GLN
1	A	299	LEU
1	A	312	ARG
1	B	44	LEU
1	B	51	VAL
1	B	70	VAL
1	B	82	VAL
1	B	114	GLU
1	B	161	ASN
1	B	205	THR
1	B	230	LEU
1	B	246	LEU

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Mol	Chain	Res	Type
1	B	255	VAL
1	B	256	VAL
1	B	260	VAL
1	B	268	TYR
1	B	288	LEU
1	B	299	LEU
1	B	310	LEU
1	B	312	ARG
1	C	44	LEU
1	C	68	LEU
1	C	70	VAL
1	C	77	VAL
1	C	82	VAL
1	C	90	GLU
1	C	114	GLU
1	C	127	PHE
1	C	139	ILE
1	C	151	ASP
1	C	230	LEU
1	C	232	VAL
1	C	246	LEU
1	C	255	VAL
1	C	256	VAL
1	C	268	TYR
1	C	299	LEU
1	C	312	ARG
1	D	44	LEU
1	D	47	GLU
1	D	65	GLU
1	D	68	LEU
1	D	70	VAL
1	D	80	LYS
1	D	89	THR
1	D	114	GLU
1	D	144	GLU
1	D	145	GLU
1	D	224	GLU
1	D	230	LEU
1	D	255	VAL
1	D	256	VAL
1	D	260	VAL
1	D	299	LEU

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Mol	Chain	Res	Type
1	D	312	ARG
1	D	315	SER
1	D	322	THR
1	E	34	LYS
1	E	44	LEU
1	E	55	VAL
1	E	82	VAL
1	E	84	GLN
1	E	90	GLU
1	E	127	PHE
1	E	139	ILE
1	E	188	VAL
1	E	202	LEU
1	E	219	SER
1	E	230	LEU
1	E	232	VAL
1	E	246	LEU
1	E	255	VAL
1	E	256	VAL
1	E	259	SER
1	E	282	VAL
1	E	288	LEU
1	E	299	LEU
1	E	310	LEU
1	E	312	ARG
1	F	44	LEU
1	F	49	ASN
1	F	55	VAL
1	F	70	VAL
1	F	82	VAL
1	F	90	GLU
1	F	155	LYS
1	F	192	PHE
1	F	210	THR
1	F	256	VAL
1	F	282	VAL
1	F	288	LEU
1	F	297	GLN
1	F	299	LEU
1	F	312	ARG
1	F	322	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (60)

such sidechains are listed below:

Mol	Chain	Res	Type
1	A	84	GLN
1	A	96	GLN
1	A	138	ASN
1	A	140	ASN
1	A	161	ASN
1	A	223	GLN
1	A	244	ASN
1	A	250	GLN
1	A	297	GLN
1	A	311	ASN
1	A	320	GLN
1	B	46	ASN
1	B	138	ASN
1	B	223	GLN
1	B	244	ASN
1	B	250	GLN
1	B	297	GLN
1	B	308	GLN
1	B	311	ASN
1	B	320	GLN
1	C	45	GLN
1	C	84	GLN
1	C	88	HIS
1	C	138	ASN
1	C	140	ASN
1	C	223	GLN
1	C	250	GLN
1	C	311	ASN
1	D	45	GLN
1	D	84	GLN
1	D	88	HIS
1	D	138	ASN
1	D	140	ASN
1	D	161	ASN
1	D	244	ASN
1	D	250	GLN
1	D	291	HIS
1	D	311	ASN
1	D	320	GLN
1	E	46	ASN
1	E	84	GLN
1	E	96	GLN

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Mol	Chain	Res	Type
1	E	102	ASN
1	E	140	ASN
1	E	161	ASN
1	E	223	GLN
1	E	244	ASN
1	E	250	GLN
1	E	275	ASN
1	E	291	HIS
1	F	102	ASN
1	F	138	ASN
1	F	223	GLN
1	F	244	ASN
1	F	250	GLN
1	F	297	GLN
1	F	304	ASN
1	F	308	GLN
1	F	311	ASN
1	F	320	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 6 are monoatomic - leaving 2 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	$\# Z > 2$
2	3HG	A	400	3	9,9,9	1.47	0	8,11,11	1.89	2 (25%)
4	HGG	B	399	3	57,59,59	3.81	27 (47%)	79,87,87	2.96	29 (36%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	3HG	A	400	3	-	0/8/8/8	-
4	HGG	B	399	3	-	19/59/75/75	0/3/3/3

All (27) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	399	HGG	C2-N3	12.08	1.55	1.33
4	B	399	HGG	C4-N3	9.55	1.52	1.34
4	B	399	HGG	P31-O32	8.72	1.68	1.59
4	B	399	HGG	C6-N1	8.58	1.59	1.35
4	B	399	HGG	P35-O32	7.65	1.67	1.59
4	B	399	HGG	C2-N1	7.20	1.46	1.33
4	B	399	HGG	C5-C4	6.95	1.51	1.39
4	B	399	HGG	C69-N71	5.84	1.47	1.33
4	B	399	HGG	P39-O21	4.86	1.68	1.59
4	B	399	HGG	C5-C6	4.58	1.53	1.41
4	B	399	HGG	O72-C69	4.30	1.31	1.23
4	B	399	HGG	C8-N9	4.19	1.44	1.37
4	B	399	HGG	O4'-C5'	3.96	1.35	1.22
4	B	399	HGG	C16-C17	3.90	1.61	1.53
4	B	399	HGG	O51-C43	3.71	1.30	1.23
4	B	399	HGG	O1'-C1'	3.28	1.26	1.21
4	B	399	HGG	C4'-C5'	-3.15	1.42	1.51
4	B	399	HGG	C61-C69	3.10	1.57	1.51
4	B	399	HGG	P35-O36	3.04	1.61	1.50
4	B	399	HGG	C43-N68	2.99	1.40	1.33
4	B	399	HGG	P39-O41	2.99	1.59	1.50
4	B	399	HGG	C57-C45	-2.92	1.47	1.53
4	B	399	HGG	C2'-C1'	-2.57	1.46	1.51
4	B	399	HGG	C73-C74	-2.31	1.41	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	399	HGG	O20-C16	2.29	1.48	1.43
4	B	399	HGG	C62-N68	-2.22	1.41	1.46
4	B	399	HGG	O22-C15	2.01	1.46	1.42

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	399	HGG	P39-O21-C17	10.41	151.24	123.43
4	B	399	HGG	O38-C46-C45	10.19	126.94	110.55
4	B	399	HGG	C2'-C1'-S81	6.91	122.28	113.56
4	B	399	HGG	O1'-C1'-S81	-6.83	113.99	122.68
4	B	399	HGG	C74-N71-C69	6.80	135.49	122.82
4	B	399	HGG	N3-C2-N1	-6.56	118.65	128.58
4	B	399	HGG	C73-S81-C1'	5.78	118.92	101.84
4	B	399	HGG	C62-C61-C69	5.64	121.79	112.39
4	B	399	HGG	C73-C74-N71	4.49	121.78	112.41
4	B	399	HGG	O22-C18-C19	-4.18	95.95	109.33
4	B	399	HGG	C6-C5-C4	-3.57	112.30	117.18
4	B	399	HGG	C18-O22-C15	3.45	117.08	109.47
2	A	400	3HG	O5-C5-C4	3.18	123.88	114.00
4	B	399	HGG	O23-C19-C18	3.15	119.73	108.99
4	B	399	HGG	C5-C4-N9	-3.14	102.39	105.81
4	B	399	HGG	O72-C69-N71	3.13	129.17	123.03
4	B	399	HGG	C4-N9-C15	-2.96	119.72	126.63
4	B	399	HGG	C15-N9-C8	2.83	133.37	127.09
2	A	400	3HG	O1-C1-O3	-2.75	116.27	123.33
4	B	399	HGG	C53-C45-C46	-2.66	103.84	108.22
4	B	399	HGG	O37-P35-O32	2.57	114.22	107.27
4	B	399	HGG	O32-P35-O36	-2.41	103.47	110.70
4	B	399	HGG	C2-N3-C4	2.35	117.56	111.83
4	B	399	HGG	C61-C62-N68	-2.29	107.12	112.00
4	B	399	HGG	C61-C69-N71	-2.28	112.18	116.34
4	B	399	HGG	O20-C16-C17	2.27	117.53	111.19
4	B	399	HGG	C16-C15-N9	2.17	118.70	113.30
4	B	399	HGG	O5'-C5'-C4'	2.12	120.59	114.00
4	B	399	HGG	C44-C43-N68	-2.07	112.56	116.48
4	B	399	HGG	C19-C18-C17	2.06	121.24	114.38
4	B	399	HGG	O22-C15-C16	-2.03	102.27	106.62

There are no chirality outliers.

All (19) torsion outliers are listed below:

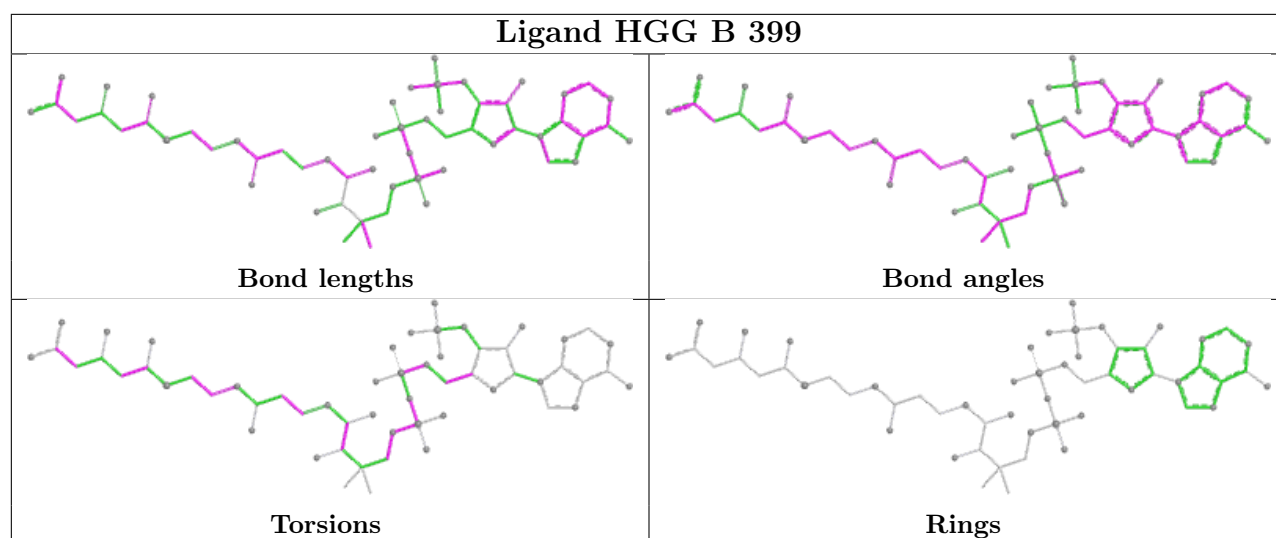
Mol	Chain	Res	Type	Atoms
4	B	399	HGG	S81-C1'-C2'-C3'
4	B	399	HGG	C19-O23-P31-O32
4	B	399	HGG	C46-O38-P35-O32
4	B	399	HGG	C46-O38-P35-O36
4	B	399	HGG	C46-O38-P35-O37
4	B	399	HGG	O51-C43-C44-O52
4	B	399	HGG	N68-C43-C44-O52
4	B	399	HGG	C69-C61-C62-N68
4	B	399	HGG	S81-C73-C74-N71
4	B	399	HGG	C17-C18-C19-O23
4	B	399	HGG	O22-C18-C19-O23
4	B	399	HGG	O51-C43-C44-C45
4	B	399	HGG	N68-C43-C44-C45
4	B	399	HGG	P31-O32-P35-O38
4	B	399	HGG	C73-C74-N71-C69
4	B	399	HGG	O1'-C1'-C2'-C3'
4	B	399	HGG	C3'-C4'-C5'-O5'
4	B	399	HGG	C3'-C4'-C5'-O4'
4	B	399	HGG	C45-C46-O38-P35

There are no ring outliers.

1 monomer is involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	399	HGG	10	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	296/298 (99%)	-1.73	0 100 100	8, 21, 44, 60	0
1	B	296/298 (99%)	-1.65	0 100 100	10, 27, 59, 74	0
1	C	296/298 (99%)	-1.63	0 100 100	10, 30, 52, 66	0
1	D	289/298 (96%)	-1.56	0 100 100	17, 37, 62, 79	0
1	E	296/298 (99%)	-1.67	0 100 100	11, 26, 56, 69	0
1	F	296/298 (99%)	-1.50	0 100 100	16, 42, 64, 73	0
All	All	1769/1788 (98%)	-1.62	0 100 100	8, 29, 59, 79	0

There are no RSRZ outliers to report.

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

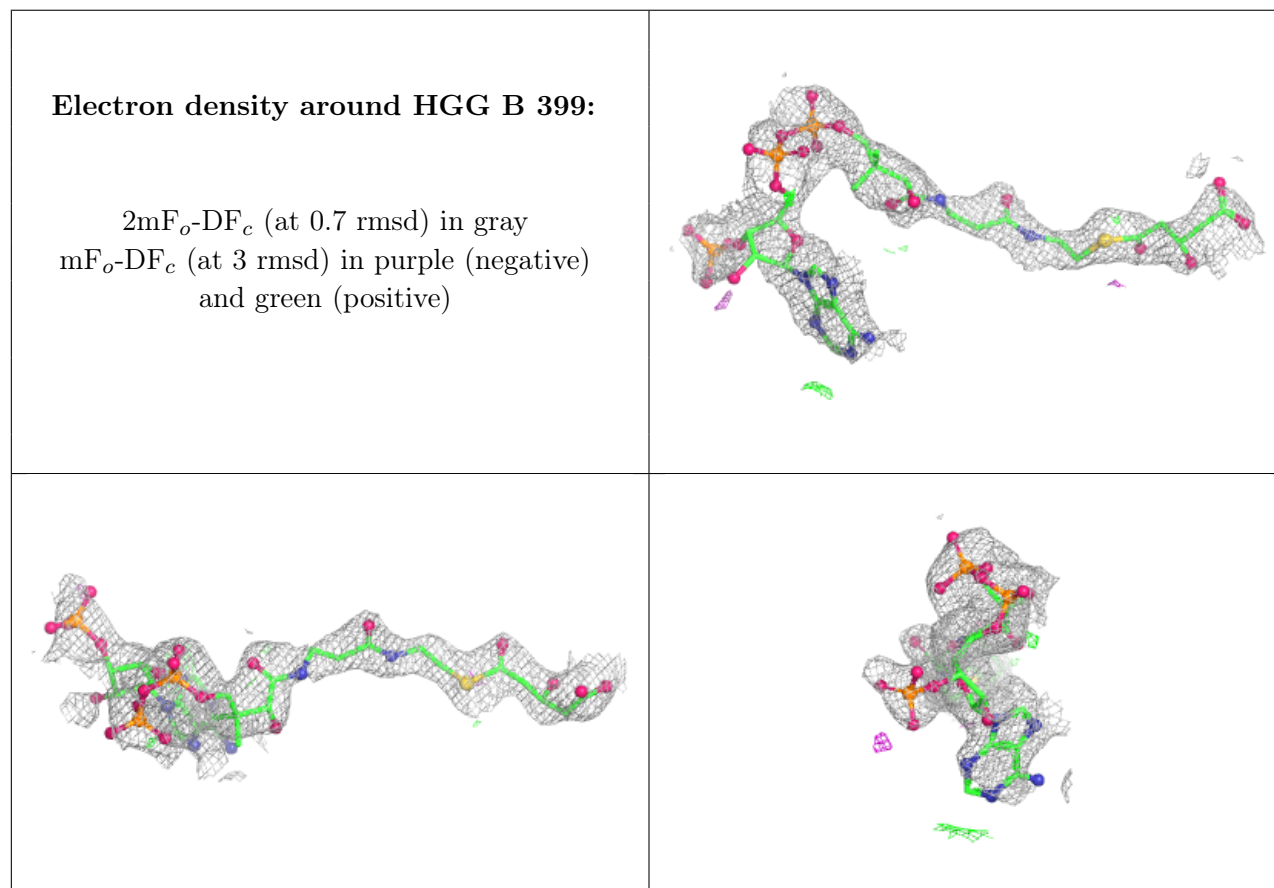
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	HGG	B	399	57/57	0.98	0.04	58,79,96,97	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	MG	A	403	1/1	0.99	0.01	16,16,16,16	0
3	MG	E	402	1/1	0.99	0.04	13,13,13,13	0
3	MG	F	406	1/1	0.99	0.04	30,30,30,30	0
2	3HG	A	400	10/10	0.99	0.05	61,63,66,67	0
3	MG	B	401	1/1	1.00	0.03	17,17,17,17	0
3	MG	C	404	1/1	1.00	0.03	15,15,15,15	0
3	MG	D	405	1/1	1.00	0.01	28,28,28,28	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.



6.5 Other polymers ⓘ

There are no such residues in this entry.