



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 2, 2026 – 01:17 AM UTC

PDB ID : 1XDY / pdb_00001xdy
Title : Structural and Biochemical Identification of a Novel Bacterial Oxidoreductase, W-containing cofactor
Authors : Loschi, L.; Brokx, S.J.; Hills, T.L.; Zhang, G.; Bertero, M.G.; Lovering, A.L.; Weiner, J.H.; Strynadka, N.C.
Deposited on : 2004-09-08
Resolution : 2.20 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
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

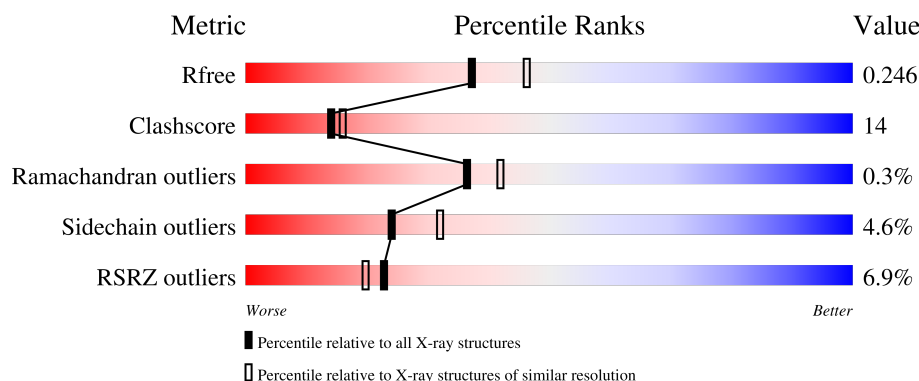
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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

Mol	Chain	Length	Quality of chain
1	A	298	3% 64% 19% 5% 12%
1	B	298	9% 59% 26% • 12%
1	C	298	3% 62% 24% •• 11%
1	D	298	% 60% 24% • 12%

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Mol	Chain	Length	Quality of chain
1	E	298	
1	F	298	
1	G	298	
1	H	298	
1	I	298	
1	J	298	

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
3	MTE	A	301	X	-	-	-
3	MTE	B	302	X	-	-	-
3	MTE	C	303	X	-	-	-
3	MTE	D	304	X	-	-	-
3	MTE	E	305	X	-	-	-
3	MTE	F	306	X	-	-	-
3	MTE	G	307	X	-	-	-
3	MTE	H	308	X	-	-	-
3	MTE	I	309	X	-	-	-
3	MTE	J	310	X	-	-	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 21369 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 Bacterial Sulfite Oxidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	263	Total	C	N	O	S	0	0	0
			2093	1358	354	375	6			
1	B	262	Total	C	N	O	S	0	0	0
			2088	1354	353	375	6			
1	C	264	Total	C	N	O	S	0	0	0
			2102	1363	356	377	6			
1	D	263	Total	C	N	O	S	0	0	0
			2094	1357	355	376	6			
1	E	262	Total	C	N	O	S	0	0	0
			2085	1351	353	375	6			
1	F	262	Total	C	N	O	S	1	0	0
			2084	1352	352	374	6			
1	G	261	Total	C	N	O	S	0	0	0
			2079	1349	351	373	6			
1	H	264	Total	C	N	O	S	11	0	0
			2102	1363	356	377	6			
1	I	262	Total	C	N	O	S	0	0	0
			2085	1351	353	375	6			
1	J	262	Total	C	N	O	S	0	0	0
			2085	1351	353	375	6			

There are 90 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	275	ASP	ALA	SEE REMARK 999	UNP P76342
A	291	LEU	-	expression tag	UNP P76342
A	292	GLU	-	expression tag	UNP P76342
A	293	HIS	-	expression tag	UNP P76342
A	294	HIS	-	expression tag	UNP P76342
A	295	HIS	-	expression tag	UNP P76342
A	296	HIS	-	expression tag	UNP P76342
A	297	HIS	-	expression tag	UNP P76342
A	298	HIS	-	expression tag	UNP P76342

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Chain	Residue	Modelled	Actual	Comment	Reference
B	275	ASP	ALA	SEE REMARK 999	UNP P76342
B	291	LEU	-	expression tag	UNP P76342
B	292	GLU	-	expression tag	UNP P76342
B	293	HIS	-	expression tag	UNP P76342
B	294	HIS	-	expression tag	UNP P76342
B	295	HIS	-	expression tag	UNP P76342
B	296	HIS	-	expression tag	UNP P76342
B	297	HIS	-	expression tag	UNP P76342
B	298	HIS	-	expression tag	UNP P76342
C	275	ASP	ALA	SEE REMARK 999	UNP P76342
C	291	LEU	-	expression tag	UNP P76342
C	292	GLU	-	expression tag	UNP P76342
C	293	HIS	-	expression tag	UNP P76342
C	294	HIS	-	expression tag	UNP P76342
C	295	HIS	-	expression tag	UNP P76342
C	296	HIS	-	expression tag	UNP P76342
C	297	HIS	-	expression tag	UNP P76342
C	298	HIS	-	expression tag	UNP P76342
D	275	ASP	ALA	SEE REMARK 999	UNP P76342
D	291	LEU	-	expression tag	UNP P76342
D	292	GLU	-	expression tag	UNP P76342
D	293	HIS	-	expression tag	UNP P76342
D	294	HIS	-	expression tag	UNP P76342
D	295	HIS	-	expression tag	UNP P76342
D	296	HIS	-	expression tag	UNP P76342
D	297	HIS	-	expression tag	UNP P76342
D	298	HIS	-	expression tag	UNP P76342
E	275	ASP	ALA	SEE REMARK 999	UNP P76342
E	291	LEU	-	expression tag	UNP P76342
E	292	GLU	-	expression tag	UNP P76342
E	293	HIS	-	expression tag	UNP P76342
E	294	HIS	-	expression tag	UNP P76342
E	295	HIS	-	expression tag	UNP P76342
E	296	HIS	-	expression tag	UNP P76342
E	297	HIS	-	expression tag	UNP P76342
E	298	HIS	-	expression tag	UNP P76342
F	275	ASP	ALA	SEE REMARK 999	UNP P76342
F	291	LEU	-	expression tag	UNP P76342
F	292	GLU	-	expression tag	UNP P76342
F	293	HIS	-	expression tag	UNP P76342
F	294	HIS	-	expression tag	UNP P76342
F	295	HIS	-	expression tag	UNP P76342

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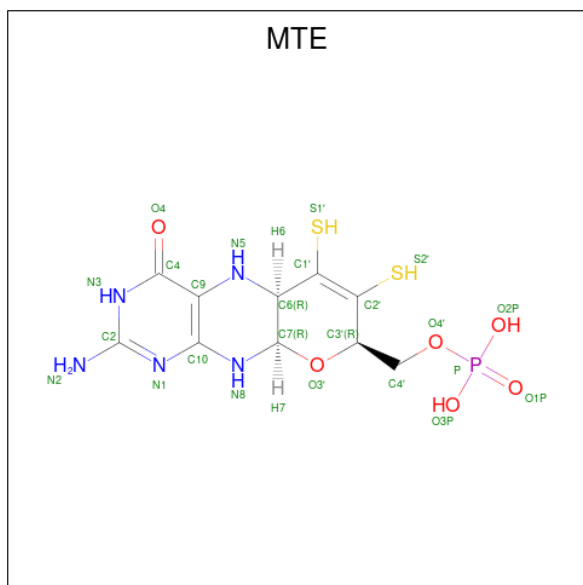
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Chain	Residue	Modelled	Actual	Comment	Reference
F	296	HIS	-	expression tag	UNP P76342
F	297	HIS	-	expression tag	UNP P76342
F	298	HIS	-	expression tag	UNP P76342
G	275	ASP	ALA	SEE REMARK 999	UNP P76342
G	291	LEU	-	expression tag	UNP P76342
G	292	GLU	-	expression tag	UNP P76342
G	293	HIS	-	expression tag	UNP P76342
G	294	HIS	-	expression tag	UNP P76342
G	295	HIS	-	expression tag	UNP P76342
G	296	HIS	-	expression tag	UNP P76342
G	297	HIS	-	expression tag	UNP P76342
G	298	HIS	-	expression tag	UNP P76342
H	275	ASP	ALA	SEE REMARK 999	UNP P76342
H	291	LEU	-	expression tag	UNP P76342
H	292	GLU	-	expression tag	UNP P76342
H	293	HIS	-	expression tag	UNP P76342
H	294	HIS	-	expression tag	UNP P76342
H	295	HIS	-	expression tag	UNP P76342
H	296	HIS	-	expression tag	UNP P76342
H	297	HIS	-	expression tag	UNP P76342
H	298	HIS	-	expression tag	UNP P76342
I	275	ASP	ALA	SEE REMARK 999	UNP P76342
I	291	LEU	-	expression tag	UNP P76342
I	292	GLU	-	expression tag	UNP P76342
I	293	HIS	-	expression tag	UNP P76342
I	294	HIS	-	expression tag	UNP P76342
I	295	HIS	-	expression tag	UNP P76342
I	296	HIS	-	expression tag	UNP P76342
I	297	HIS	-	expression tag	UNP P76342
I	298	HIS	-	expression tag	UNP P76342
J	275	ASP	ALA	SEE REMARK 999	UNP P76342
J	291	LEU	-	expression tag	UNP P76342
J	292	GLU	-	expression tag	UNP P76342
J	293	HIS	-	expression tag	UNP P76342
J	294	HIS	-	expression tag	UNP P76342
J	295	HIS	-	expression tag	UNP P76342
J	296	HIS	-	expression tag	UNP P76342
J	297	HIS	-	expression tag	UNP P76342
J	298	HIS	-	expression tag	UNP P76342

- Molecule 2 is TUNGSTEN ION (CCD ID: W) (formula: W).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total W 1 1	0	0
2	B	1	Total W 1 1	0	0
2	C	1	Total W 1 1	0	0
2	D	1	Total W 1 1	0	0
2	E	1	Total W 1 1	0	0
2	F	1	Total W 1 1	0	0
2	G	1	Total W 1 1	0	0
2	H	1	Total W 1 1	0	0
2	I	1	Total W 1 1	0	0
2	J	1	Total W 1 1	0	0

- Molecule 3 is PHOSPHONIC ACIDMONO-(2-AMINO-5,6-DIMERCAPTO-4-OXO-3,7,8A, 9,10,10A-HEXAHYDRO-4H-8-OXA-1,3,9,10-TETRAAZA-ANTHRACEN-7-YLMETHYL) ESTER (CCD ID: MTE) (formula: C₁₀H₁₄N₅O₆PS₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	S	
			24	10	5	6	1	2	0
									0

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Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	B	1	Total	C	N	O	P	S	0	0
			24	10	5	6	1	2		
3	C	1	Total	C	N	O	P	S	0	0
			24	10	5	6	1	2		
3	D	1	Total	C	N	O	P	S	0	0
			24	10	5	6	1	2		
3	E	1	Total	C	N	O	P	S	0	0
			24	10	5	6	1	2		
3	F	1	Total	C	N	O	P	S	0	0
			24	10	5	6	1	2		
3	G	1	Total	C	N	O	P	S	0	0
			24	10	5	6	1	2		
3	H	1	Total	C	N	O	P	S	0	0
			24	10	5	6	1	2		
3	I	1	Total	C	N	O	P	S	0	0
			24	10	5	6	1	2		
3	J	1	Total	C	N	O	P	S	0	0
			24	10	5	6	1	2		

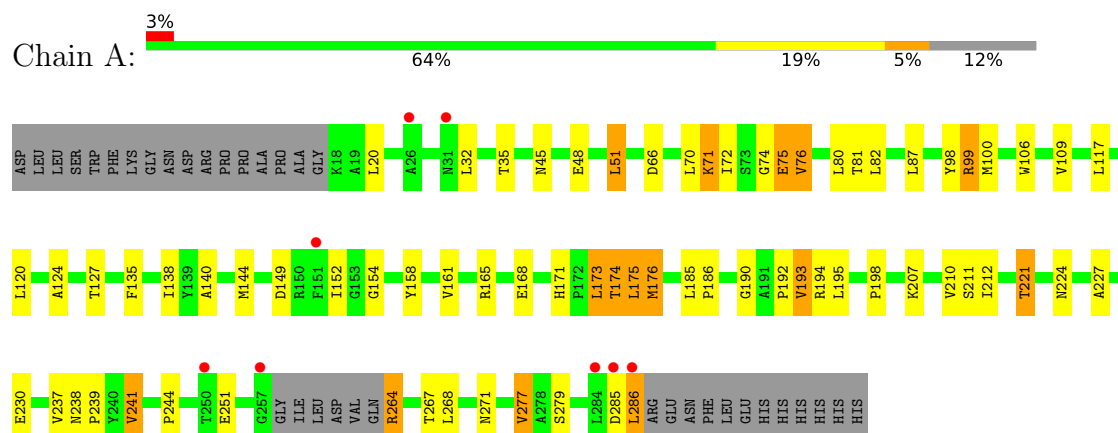
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	35	Total	O	0	0
			35	35		
4	B	13	Total	O	0	0
			13	13		
4	C	32	Total	O	0	0
			32	32		
4	D	30	Total	O	0	0
			30	30		
4	E	29	Total	O	0	0
			29	29		
4	F	10	Total	O	0	0
			10	10		
4	G	22	Total	O	0	0
			22	22		
4	H	4	Total	O	0	0
			4	4		
4	I	12	Total	O	0	0
			12	12		
4	J	35	Total	O	0	0
			35	35		

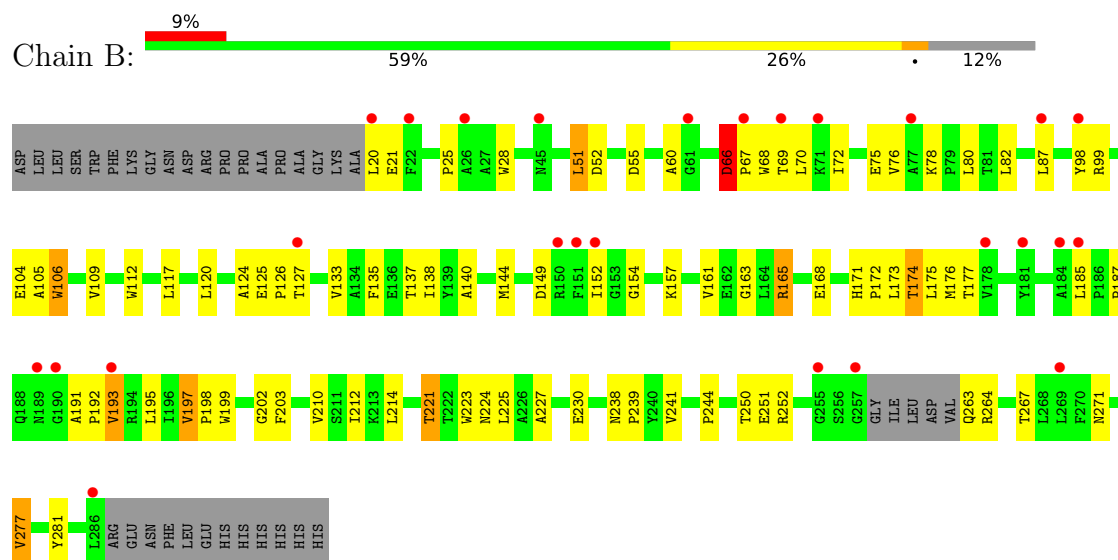
3 Residue-property plots [i](#)

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.

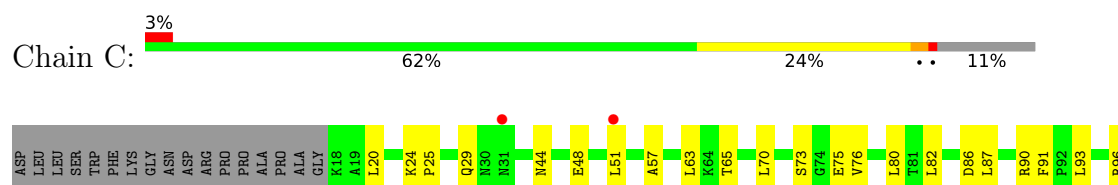
• Molecule 1: Bacterial Sulfite Oxidase

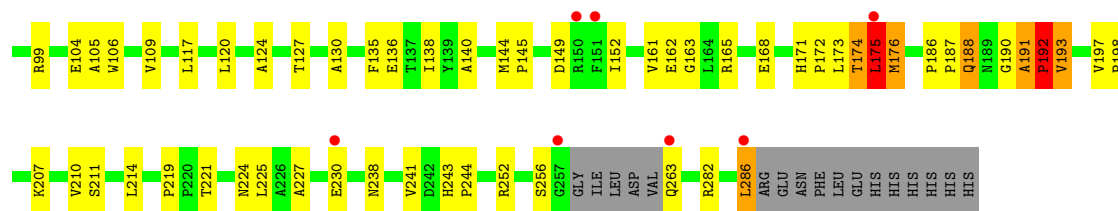


• Molecule 1: Bacterial Sulfite Oxidase

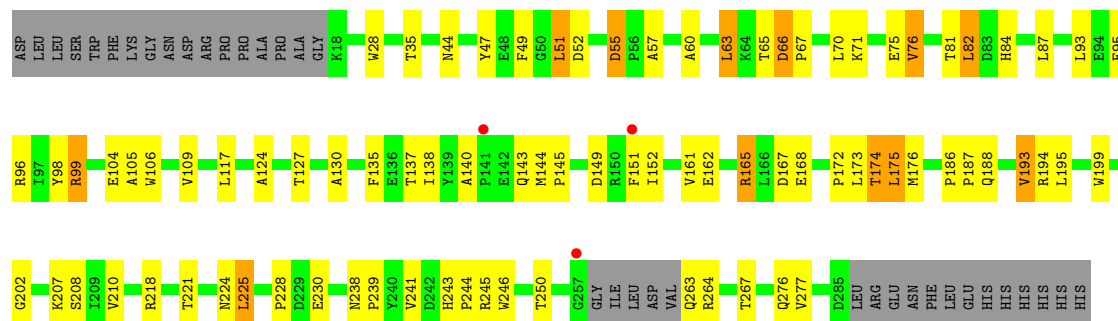


• Molecule 1: Bacterial Sulfite Oxidase

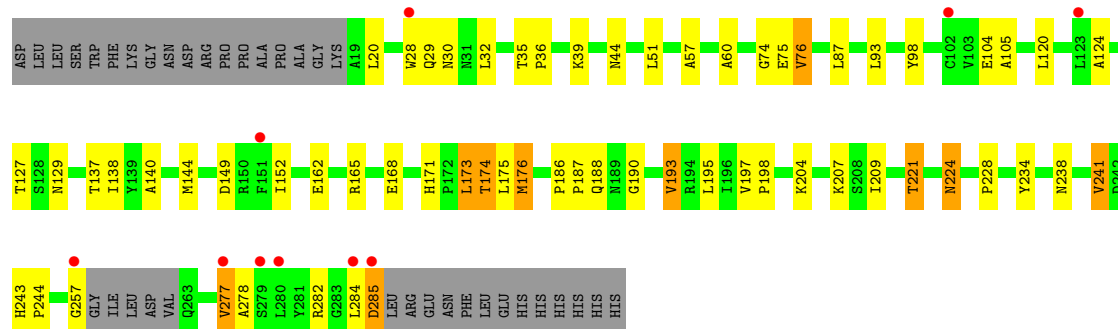




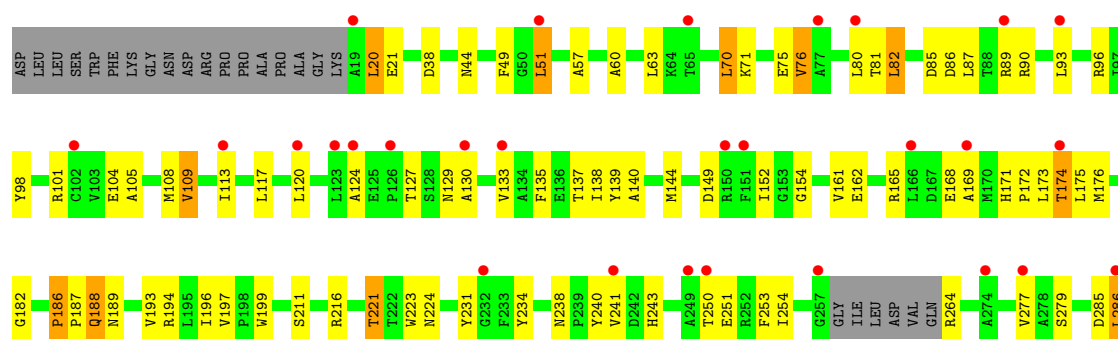
• Molecule 1: Bacterial Sulfite Oxidase



• Molecule 1: Bacterial Sulfite Oxidase



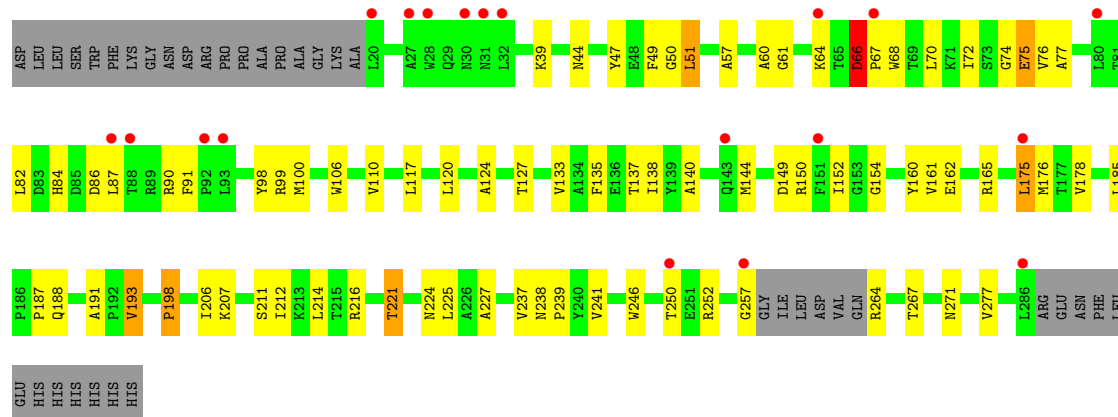
• Molecule 1: Bacterial Sulfite Oxidase



ARG
GLU
ASN
PHE
LEU
GLU
HIS
HIS
HIS
HIS
HIS

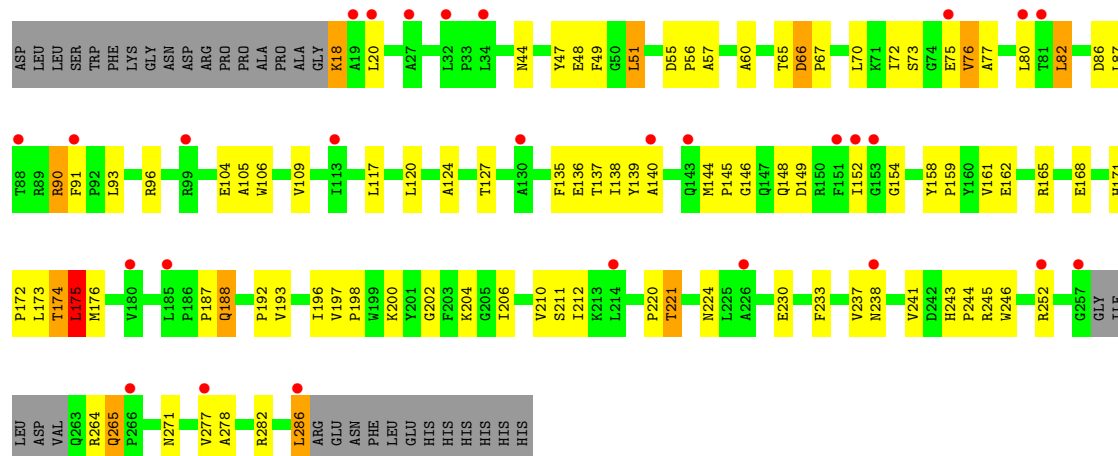
● Molecule 1: Bacterial Sulfite Oxidase

Chain G: 6% 61% 24% 12%



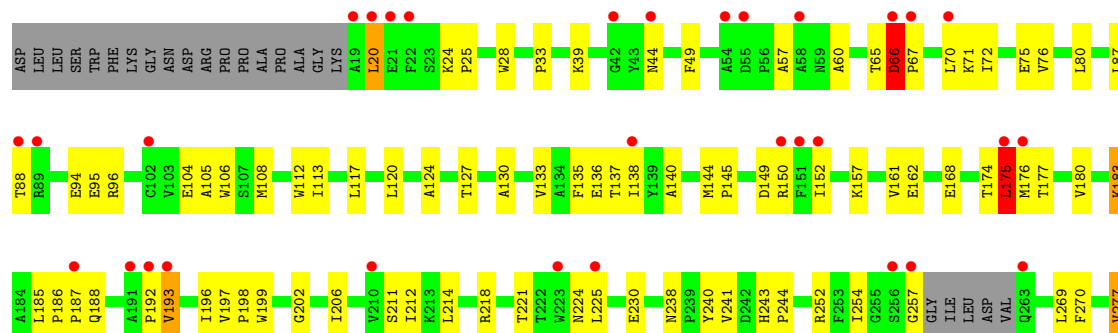
● Molecule 1: Bacterial Sulfite Oxidase

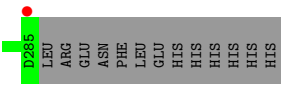
Chain H: 9% 57% 28% 11%



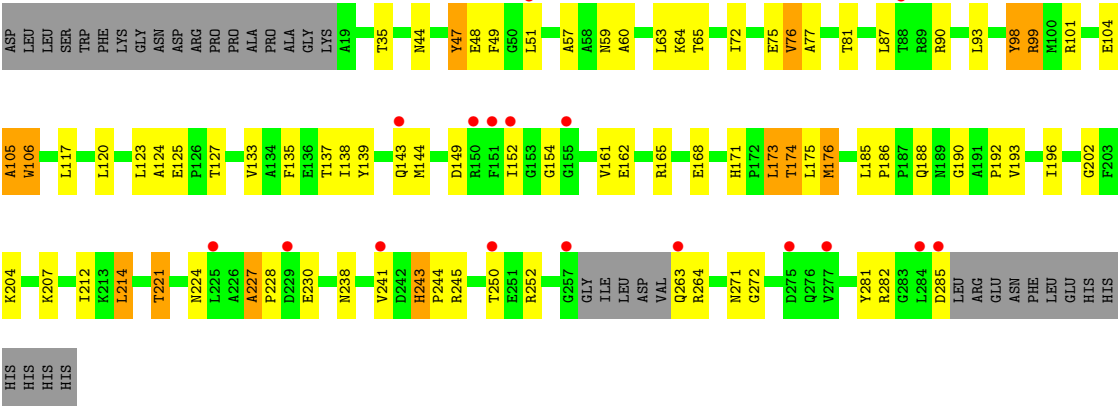
● Molecule 1: Bacterial Sulfite Oxidase

Chain I: 11% 59% 27% 12%





● Molecule 1: Bacterial Sulfite Oxidase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	84.84Å 177.51Å 104.21Å 90.00° 109.73° 90.00°	Depositor
Resolution (Å)	38.96 – 2.20 38.96 – 2.20	Depositor EDS
% Data completeness (in resolution range)	98.5 (38.96-2.20) 98.6 (38.96-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.69 (at 2.20Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.226 , 0.258 0.214 , 0.246	Depositor DCC
R_{free} test set	4389 reflections (3.03%)	wwPDB-VP
Wilson B-factor (Å ²)	43.8	Xtriage
Anisotropy	0.141	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 45.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	21369	wwPDB-VP
Average B, all atoms (Å ²)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 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: MTE, W

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	0.65	0/2158	1.06	17/2942 (0.6%)
1	B	0.53	1/2153 (0.0%)	1.05	12/2936 (0.4%)
1	C	0.61	0/2167	1.63	18/2954 (0.6%)
1	D	0.63	0/2159	1.14	17/2943 (0.6%)
1	E	0.65	0/2150	1.06	12/2932 (0.4%)
1	F	0.50	0/2149	0.99	10/2931 (0.3%)
1	G	0.62	0/2144	1.11	12/2924 (0.4%)
1	H	0.49	0/2167	1.00	12/2954 (0.4%)
1	I	0.55	0/2150	1.02	10/2932 (0.3%)
1	J	0.69	0/2150	1.09	16/2932 (0.5%)
All	All	0.59	1/21547 (0.0%)	1.13	136/29380 (0.5%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	E	0	1
1	J	0	1
All	All	0	3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	126	PRO	CA-C	5.36	1.59	1.52

All (136) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	191	ALA	CA-C-N	47.15	178.78	119.84
1	C	191	ALA	C-N-CA	47.15	178.78	119.84
1	D	66	ASP	CA-C-N	-14.41	104.77	119.90
1	D	66	ASP	C-N-CA	-14.41	104.77	119.90
1	G	66	ASP	CA-C-N	-12.10	104.71	119.84
1	G	66	ASP	C-N-CA	-12.10	104.71	119.84
1	C	191	ALA	C-N-CD	-11.87	76.34	125.00
1	B	66	ASP	CA-C-N	-11.65	106.47	119.98
1	B	66	ASP	C-N-CA	-11.65	106.47	119.98
1	H	66	ASP	CA-C-N	-10.75	106.41	119.84
1	H	66	ASP	C-N-CA	-10.75	106.41	119.84
1	I	66	ASP	CA-C-N	-10.42	106.81	119.84
1	I	66	ASP	C-N-CA	-10.42	106.81	119.84
1	E	277	VAL	N-CA-C	8.40	119.04	111.81
1	F	51	LEU	N-CA-C	8.11	120.11	111.28
1	D	35	THR	N-CA-C	-7.97	99.76	109.72
1	G	165	ARG	N-CA-C	-7.93	100.06	110.53
1	C	51	LEU	N-CA-C	7.72	119.70	111.28
1	C	44	ASN	N-CA-C	7.67	121.41	109.52
1	B	227	ALA	CA-C-N	7.53	128.20	119.47
1	B	227	ALA	C-N-CA	7.53	128.20	119.47
1	B	109	VAL	N-CA-C	-7.49	96.69	107.77
1	H	109	VAL	N-CA-C	-7.36	96.88	107.77
1	A	186	PRO	N-CA-C	-7.35	102.71	110.58
1	G	198	PRO	N-CA-C	7.33	123.36	113.84
1	B	165	ARG	N-CA-C	-7.24	99.57	110.28
1	J	202	GLY	N-CA-C	7.23	121.41	112.73
1	F	165	ARG	N-CA-C	-7.22	100.32	110.50
1	E	51	LEU	N-CA-C	7.21	120.00	111.71
1	I	175	LEU	N-CA-C	7.12	120.80	108.76
1	D	44	ASN	N-CA-C	7.12	120.55	109.52
1	D	186	PRO	N-CA-C	-7.09	102.04	110.70
1	I	186	PRO	N-CA-C	-7.00	103.09	110.58
1	J	44	ASN	N-CA-C	6.95	120.29	109.52
1	B	197	VAL	CA-C-N	6.90	127.47	119.47
1	B	197	VAL	C-N-CA	6.90	127.47	119.47
1	H	165	ARG	N-CA-C	-6.87	100.81	110.50
1	E	198	PRO	N-CA-C	6.84	122.74	113.84
1	E	165	ARG	N-CA-C	-6.84	100.85	110.50
1	C	175	LEU	N-CA-C	6.78	120.56	109.24
1	G	175	LEU	N-CA-C	6.76	120.52	109.24
1	A	176	MET	N-CA-C	-6.64	97.72	108.41
1	G	51	LEU	N-CA-C	6.61	120.61	112.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	165	ARG	N-CA-C	-6.61	101.19	110.50
1	D	109	VAL	N-CA-C	-6.61	98.86	108.11
1	F	20	LEU	N-CA-C	6.55	118.72	108.96
1	I	44	ASN	N-CA-C	6.48	119.72	109.81
1	G	44	ASN	N-CA-C	6.42	118.93	109.24
1	C	165	ARG	N-CA-C	-6.39	101.50	110.50
1	D	243	HIS	N-CA-C	-6.38	101.50	110.31
1	A	109	VAL	N-CA-C	-6.35	98.38	107.77
1	C	186	PRO	N-CA-C	-6.33	103.81	110.58
1	E	197	VAL	N-CA-C	-6.29	95.29	108.88
1	B	277	VAL	N-CA-C	6.29	119.25	112.96
1	C	192	PRO	CA-N-CD	-6.28	103.21	112.00
1	E	209	ILE	N-CA-C	6.19	117.04	109.30
1	B	51	LEU	N-CA-C	6.17	118.98	111.82
1	A	175	LEU	N-CA-C	6.14	118.96	109.07
1	G	110	VAL	N-CA-C	6.14	114.91	107.61
1	F	186	PRO	N-CA-C	-6.13	103.22	110.70
1	A	227	ALA	CA-C-N	6.13	126.58	119.47
1	A	227	ALA	C-N-CA	6.13	126.58	119.47
1	E	44	ASN	N-CA-C	6.05	119.06	109.81
1	A	194	ARG	CG-CD-NE	-6.03	98.72	112.00
1	J	165	ARG	N-CA-C	-6.01	101.58	110.48
1	C	227	ALA	CA-C-N	5.93	127.26	119.84
1	C	227	ALA	C-N-CA	5.93	127.26	119.84
1	I	202	GLY	N-CA-C	5.89	120.01	112.83
1	C	176	MET	N-CA-C	-5.85	100.99	109.59
1	E	186	PRO	N-CA-C	-5.81	103.61	110.70
1	J	106	TRP	N-CA-C	-5.79	100.36	109.50
1	F	44	ASN	N-CA-C	5.78	118.66	109.81
1	H	51	LEU	N-CA-C	5.78	117.66	111.36
1	D	143	GLN	N-CA-C	-5.78	107.22	114.56
1	H	73	SER	N-CA-C	5.75	115.61	108.19
1	H	202	GLY	N-CA-C	5.75	120.91	113.27
1	J	185	LEU	N-CA-C	5.72	116.18	109.60
1	J	245	ARG	N-CA-C	5.72	120.25	113.16
1	A	71	LYS	N-CA-C	5.70	117.82	108.52
1	C	99	ARG	N-CA-C	-5.70	101.25	110.32
1	I	198	PRO	N-CA-C	5.68	122.05	113.81
1	F	197	VAL	CA-C-N	5.68	126.06	119.47
1	F	197	VAL	C-N-CA	5.68	126.06	119.47
1	G	227	ALA	CA-C-N	5.68	126.06	119.47
1	G	227	ALA	C-N-CA	5.68	126.06	119.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	106	TRP	N-CA-C	-5.67	100.73	109.52
1	E	224	ASN	N-CA-C	-5.66	105.26	111.82
1	B	202	GLY	N-CA-C	5.65	119.73	112.83
1	C	109	VAL	N-CA-C	-5.64	99.43	107.77
1	D	210	VAL	N-CA-C	-5.63	108.36	113.71
1	J	176	MET	N-CA-C	-5.56	100.75	109.25
1	A	198	PRO	N-CA-C	5.53	121.03	113.84
1	J	186	PRO	N-CA-C	-5.53	103.96	110.70
1	A	99	ARG	N-CA-C	-5.52	101.29	110.17
1	J	227	ALA	CA-C-N	5.51	126.73	119.84
1	J	227	ALA	C-N-CA	5.51	126.73	119.84
1	A	277	VAL	N-CA-C	5.50	119.02	113.47
1	B	106	TRP	N-CA-C	-5.46	100.87	109.50
1	F	243	HIS	N-CA-C	-5.45	103.20	110.07
1	C	73	SER	N-CA-C	5.45	115.22	108.19
1	E	35	THR	N-CA-C	-5.43	102.93	109.72
1	G	207	LYS	N-CA-C	5.42	118.55	110.52
1	H	243	HIS	N-CA-C	-5.42	103.25	110.07
1	A	45	ASN	N-CA-C	-5.38	99.64	108.41
1	A	106	TRP	N-CA-C	-5.33	101.27	109.52
1	I	197	VAL	CA-C-N	5.32	125.64	119.47
1	I	197	VAL	C-N-CA	5.32	125.64	119.47
1	J	59	ASN	N-CA-C	5.31	120.77	113.97
1	A	185	LEU	N-CA-C	5.29	117.10	109.48
1	D	165	ARG	N-CA-C	-5.27	103.57	110.53
1	J	47	TYR	N-CA-C	5.27	117.02	111.28
1	D	175	LEU	N-CA-C	5.26	117.82	109.50
1	J	35	THR	N-CA-C	-5.26	103.02	109.65
1	C	243	HIS	N-CA-C	-5.24	103.06	110.29
1	D	202	GLY	N-CA-C	5.24	120.10	113.24
1	J	105	ALA	N-CA-C	5.21	118.90	111.92
1	C	198	PRO	N-CA-C	5.18	120.64	113.65
1	E	29	GLN	N-CA-C	-5.17	101.84	110.17
1	F	109	VAL	N-CA-C	-5.17	100.21	107.75
1	G	246	TRP	N-CA-C	5.17	116.37	108.52
1	A	35	THR	N-CA-C	-5.16	103.15	109.65
1	A	51	LEU	N-CA-C	5.16	117.80	111.82
1	D	194	ARG	N-CA-CB	-5.14	101.96	111.52
1	I	243	HIS	N-CA-C	-5.13	102.57	110.32
1	C	211	SER	N-CA-C	5.10	117.67	108.48
1	H	206	ILE	N-CA-C	5.09	115.97	109.30
1	D	99	ARG	N-CA-C	-5.08	102.24	110.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	175	LEU	N-CA-C	5.08	117.52	109.50
1	H	44	ASN	N-CA-C	5.04	117.32	109.52
1	J	243	HIS	N-CA-C	-5.03	103.73	110.07
1	D	55	ASP	CA-C-N	-5.03	113.47	119.05
1	D	55	ASP	C-N-CA	-5.03	113.47	119.05
1	H	210	VAL	N-CA-C	-5.03	108.93	113.71
1	E	176	MET	N-CA-C	-5.01	100.90	108.67
1	J	143	GLN	N-CA-C	-5.01	107.84	114.31
1	F	21	GLU	N-CA-C	5.00	117.40	109.24

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	98	TYR	Sidechain
1	E	234	TYR	Sidechain
1	J	98	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	2093	0	2067	54	0
1	B	2088	0	2057	68	0
1	C	2102	0	2075	52	0
1	D	2094	0	2064	68	0
1	E	2085	0	2051	48	0
1	F	2084	0	2054	68	0
1	G	2079	0	2049	57	0
1	H	2102	0	2075	63	0
1	I	2085	0	2051	70	0
1	J	2085	0	2051	63	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	F	1	0	0	0	0
2	G	1	0	0	0	0
2	H	1	0	0	0	0
2	I	1	0	0	0	0
2	J	1	0	0	0	0
3	A	24	0	9	0	0
3	B	24	0	9	0	0
3	C	24	0	9	0	0
3	D	24	0	9	0	0
3	E	24	0	9	0	0
3	F	24	0	9	0	0
3	G	24	0	9	0	0
3	H	24	0	9	0	0
3	I	24	0	9	0	0
3	J	24	0	9	0	0
4	A	35	0	0	1	0
4	B	13	0	0	1	0
4	C	32	0	0	2	0
4	D	30	0	0	0	0
4	E	29	0	0	1	0
4	F	10	0	0	0	0
4	G	22	0	0	1	0
4	H	4	0	0	0	0
4	I	12	0	0	0	0
4	J	35	0	0	2	0
All	All	21369	0	20684	599	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 14.

All (599) 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:B:117:LEU:HD13	1:B:176:MET:HE3	1.45	0.98
1:J:117:LEU:HD13	1:J:176:MET:HE3	1.47	0.94
1:D:250:THR:HG23	1:D:264:ARG:HH21	1.33	0.92
1:E:140:ALA:HB3	1:E:144:MET:HE3	1.50	0.92
1:H:57:ALA:HA	1:H:188:GLN:HG2	1.51	0.92
1:B:20:LEU:HG	1:B:21:GLU:H	1.35	0.90
1:G:117:LEU:HD13	1:G:176:MET:HE3	1.51	0.90
1:C:138:ILE:HG23	1:C:144:MET:HE1	1.56	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:117:LEU:HD13	1:C:176:MET:HE3	1.57	0.87
1:I:117:LEU:HD13	1:I:176:MET:HE3	1.55	0.86
1:A:140:ALA:HB3	1:A:144:MET:HE3	1.58	0.86
1:I:57:ALA:HA	1:I:188:GLN:HG2	1.55	0.85
1:E:57:ALA:HA	1:E:188:GLN:HG2	1.60	0.84
1:J:221:THR:HG21	4:J:440:HOH:O	1.76	0.84
1:F:117:LEU:HD13	1:F:176:MET:HE3	1.60	0.83
1:E:87:LEU:HD13	1:E:193:VAL:HG22	1.59	0.83
1:A:138:ILE:HG23	1:A:144:MET:HE1	1.60	0.83
1:J:250:THR:HG23	1:J:264:ARG:HE	1.44	0.82
1:G:250:THR:HG23	1:G:264:ARG:HH21	1.44	0.81
1:H:66:ASP:HB3	1:H:67:PRO:HD3	1.63	0.81
1:E:30:ASN:HD22	1:E:32:LEU:H	1.29	0.81
1:H:138:ILE:HG23	1:H:144:MET:HE1	1.62	0.81
1:C:57:ALA:HA	1:C:188:GLN:HG2	1.60	0.81
1:D:230:GLU:HG2	1:D:244:PRO:HG2	1.61	0.80
1:G:76:VAL:HG13	1:G:124:ALA:O	1.82	0.79
1:E:57:ALA:HA	1:E:188:GLN:CG	2.12	0.79
1:E:138:ILE:HG23	1:E:144:MET:HE1	1.64	0.79
1:A:117:LEU:HD13	1:A:176:MET:HE3	1.65	0.79
1:J:57:ALA:HA	1:J:188:GLN:CG	2.13	0.79
1:A:87:LEU:HD13	1:A:193:VAL:HG22	1.62	0.79
1:F:211:SER:OG	1:H:282:ARG:HD3	1.84	0.78
1:A:76:VAL:HG13	1:A:124:ALA:O	1.83	0.78
1:D:75:GLU:HG3	1:D:127:THR:OG1	1.84	0.78
1:H:286:LEU:HD22	1:H:286:LEU:H	1.47	0.78
1:H:76:VAL:HG13	1:H:124:ALA:O	1.84	0.78
1:I:75:GLU:HG3	1:I:127:THR:OG1	1.84	0.78
1:I:230:GLU:HG2	1:I:244:PRO:HG2	1.66	0.78
1:I:28:TRP:CG	1:I:96:ARG:HG2	2.19	0.77
1:H:168:GLU:O	1:H:174:THR:HG21	1.83	0.77
1:C:57:ALA:HA	1:C:188:GLN:CG	2.15	0.77
1:J:230:GLU:HG2	1:J:244:PRO:HG2	1.64	0.77
1:G:87:LEU:HD13	1:G:193:VAL:HG22	1.66	0.77
1:I:72:ILE:HD13	1:I:212:ILE:HB	1.67	0.76
1:J:149:ASP:HB3	1:J:152:ILE:HB	1.68	0.76
1:C:65:THR:HB	1:C:192:PRO:CD	2.16	0.76
1:D:57:ALA:HA	1:D:188:GLN:CG	2.16	0.76
1:F:140:ALA:H	1:F:144:MET:HE3	1.50	0.76
1:D:87:LEU:HD13	1:D:193:VAL:HG22	1.67	0.75
1:H:57:ALA:HA	1:H:188:GLN:CG	2.16	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:87:LEU:HD13	1:J:193:VAL:CG2	2.17	0.74
1:B:99:ARG:HG3	1:B:271:ASN:ND2	2.03	0.74
1:C:221:THR:HG21	4:C:426:HOH:O	1.85	0.74
1:D:117:LEU:HD13	1:D:176:MET:HE3	1.68	0.74
1:I:57:ALA:HA	1:I:188:GLN:CG	2.18	0.73
1:J:221:THR:HG22	1:J:224:ASN:H	1.53	0.73
1:H:230:GLU:HG2	1:H:244:PRO:HG2	1.68	0.73
1:H:87:LEU:HD13	1:H:193:VAL:HG22	1.69	0.72
1:C:238:ASN:HB3	1:C:241:VAL:HG12	1.70	0.72
1:J:238:ASN:HB3	1:J:241:VAL:HG12	1.72	0.72
1:A:221:THR:HG22	1:A:224:ASN:H	1.55	0.71
1:A:168:GLU:O	1:A:174:THR:HG21	1.91	0.71
1:I:87:LEU:HD13	1:I:193:VAL:HG22	1.73	0.71
1:H:238:ASN:HB3	1:H:241:VAL:HG12	1.72	0.71
1:F:57:ALA:HA	1:F:188:GLN:CG	2.21	0.70
1:I:137:THR:HG22	1:I:138:ILE:N	2.06	0.70
1:I:39:LYS:NZ	1:I:257:GLY:H	1.89	0.70
1:G:57:ALA:HA	1:G:188:GLN:HG3	1.73	0.70
1:C:168:GLU:O	1:C:174:THR:HG21	1.92	0.70
1:I:140:ALA:H	1:I:144:MET:HE3	1.56	0.70
1:D:162:GLU:OE2	1:D:221:THR:HG23	1.91	0.70
1:E:238:ASN:HB3	1:E:241:VAL:CG1	2.22	0.70
1:A:99:ARG:HD2	1:A:271:ASN:O	1.91	0.69
1:H:60:ALA:HB1	1:H:187:PRO:HB2	1.75	0.69
1:F:60:ALA:HB1	1:F:187:PRO:HB2	1.72	0.69
1:C:187:PRO:HA	1:C:191:ALA:HB2	1.75	0.69
1:B:87:LEU:HD13	1:B:193:VAL:HG22	1.75	0.68
1:C:65:THR:HB	1:C:192:PRO:HD3	1.74	0.68
1:H:96:ARG:HD2	1:H:175:LEU:HD13	1.75	0.68
1:I:133:VAL:HG22	1:I:214:LEU:HD13	1.75	0.68
1:I:162:GLU:OE2	1:I:221:THR:HG23	1.94	0.68
1:J:75:GLU:HG3	1:J:127:THR:OG1	1.93	0.68
1:H:162:GLU:OE2	1:H:221:THR:HG23	1.94	0.68
1:J:168:GLU:O	1:J:174:THR:HG21	1.93	0.68
1:B:230:GLU:HG2	1:B:244:PRO:HG2	1.74	0.68
1:E:87:LEU:HD13	1:E:193:VAL:CG2	2.24	0.68
1:G:51:LEU:HD21	1:G:154:GLY:CA	2.23	0.68
1:F:120:LEU:O	1:F:120:LEU:HD23	1.93	0.68
1:H:149:ASP:HB3	1:H:152:ILE:HB	1.75	0.67
1:I:66:ASP:HB3	1:I:67:PRO:HD3	1.75	0.67
1:A:32:LEU:HD11	1:A:99:ARG:HH11	1.59	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:238:ASN:HB3	1:D:241:VAL:HG12	1.77	0.67
1:G:51:LEU:HD21	1:G:154:GLY:N	2.09	0.67
1:A:51:LEU:HD21	1:A:154:GLY:CA	2.25	0.66
1:J:87:LEU:HD22	1:J:193:VAL:HG21	1.77	0.66
1:E:221:THR:HG22	1:E:224:ASN:H	1.60	0.66
1:B:241:VAL:HG13	4:B:413:HOH:O	1.96	0.66
1:D:250:THR:HG23	1:D:264:ARG:NH2	2.07	0.66
1:F:87:LEU:HD13	1:F:193:VAL:CG2	2.25	0.66
1:J:250:THR:CG2	1:J:264:ARG:HE	2.08	0.66
1:B:238:ASN:HB3	1:B:241:VAL:CG1	2.26	0.66
1:F:238:ASN:HB3	1:F:241:VAL:HG12	1.78	0.66
1:J:57:ALA:HA	1:J:188:GLN:HG3	1.77	0.66
1:A:99:ARG:HG3	1:A:271:ASN:ND2	2.11	0.65
1:G:221:THR:HG22	1:G:224:ASN:H	1.61	0.65
1:G:138:ILE:HG23	1:G:144:MET:HE1	1.78	0.65
1:D:140:ALA:H	1:D:144:MET:HE3	1.60	0.65
1:E:241:VAL:HG13	4:E:427:HOH:O	1.95	0.65
1:D:57:ALA:HA	1:D:188:GLN:HG2	1.80	0.64
1:I:76:VAL:HG13	1:I:124:ALA:O	1.97	0.64
1:C:238:ASN:HB3	1:C:241:VAL:CG1	2.27	0.64
1:F:51:LEU:HD21	1:F:154:GLY:CA	2.27	0.64
1:D:149:ASP:OD2	1:D:152:ILE:HG12	1.97	0.64
1:J:63:LEU:HD12	1:J:65:THR:HG23	1.80	0.64
1:I:180:VAL:HG12	1:I:185:LEU:HD23	1.78	0.64
1:D:135:PHE:O	1:D:161:VAL:HG23	1.98	0.64
1:H:57:ALA:CA	1:H:188:GLN:HG2	2.25	0.63
1:F:51:LEU:HD21	1:F:154:GLY:HA3	1.81	0.63
1:F:162:GLU:OE2	1:F:221:THR:HG23	1.97	0.63
1:B:168:GLU:O	1:B:174:THR:HG21	1.99	0.63
1:E:57:ALA:CA	1:E:188:GLN:HG2	2.26	0.63
1:C:140:ALA:HB3	1:C:144:MET:HE3	1.81	0.62
1:H:286:LEU:HD22	1:H:286:LEU:N	2.14	0.62
1:I:57:ALA:CA	1:I:188:GLN:HG2	2.29	0.62
1:B:76:VAL:HG11	1:B:124:ALA:HB1	1.82	0.62
1:F:75:GLU:HG3	1:F:127:THR:OG1	1.98	0.62
1:E:162:GLU:OE2	1:E:221:THR:HG23	1.99	0.62
1:B:238:ASN:HB3	1:B:241:VAL:HG12	1.80	0.62
1:D:28:TRP:CG	1:D:96:ARG:HG2	2.34	0.62
1:B:75:GLU:HG3	1:B:127:THR:OG1	2.00	0.62
1:D:76:VAL:HG13	1:D:124:ALA:O	2.00	0.61
1:C:57:ALA:CA	1:C:188:GLN:HG2	2.30	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:221:THR:HG22	1:H:224:ASN:H	1.65	0.61
1:A:149:ASP:HB3	1:A:152:ILE:HB	1.82	0.61
1:H:197:VAL:HG11	1:H:200:LYS:HD2	1.82	0.61
1:J:57:ALA:HA	1:J:188:GLN:HG2	1.82	0.61
1:C:187:PRO:HA	1:C:191:ALA:CB	2.30	0.61
1:G:72:ILE:HD13	1:G:212:ILE:HB	1.81	0.60
1:I:221:THR:HG22	1:I:224:ASN:H	1.67	0.60
1:I:135:PHE:HE2	1:I:176:MET:HE1	1.67	0.60
1:J:99:ARG:HD3	1:J:271:ASN:O	2.01	0.60
1:E:60:ALA:HB1	1:E:187:PRO:HB2	1.84	0.60
1:J:149:ASP:OD2	1:J:152:ILE:HG12	2.01	0.60
1:B:99:ARG:HD2	1:B:271:ASN:O	2.02	0.60
1:A:75:GLU:HG3	1:A:127:THR:OG1	2.02	0.60
1:G:135:PHE:O	1:G:161:VAL:HG23	2.01	0.60
1:G:72:ILE:HG13	1:G:120:LEU:HD21	1.84	0.60
1:I:20:LEU:HD23	1:I:88:THR:HG21	1.84	0.59
1:B:250:THR:HG23	1:B:264:ARG:NH2	2.16	0.59
1:H:82:LEU:N	1:H:82:LEU:HD12	2.17	0.59
1:C:241:VAL:HG13	4:C:435:HOH:O	2.01	0.59
1:D:51:LEU:HD23	1:D:152:ILE:O	2.03	0.59
1:H:140:ALA:HB3	1:H:144:MET:HE3	1.84	0.59
1:D:250:THR:CG2	1:D:264:ARG:HH21	2.12	0.59
1:F:221:THR:HG22	1:F:224:ASN:H	1.67	0.59
1:F:76:VAL:HG11	1:F:124:ALA:HB1	1.84	0.58
1:C:87:LEU:HD13	1:C:193:VAL:HG22	1.85	0.58
1:F:82:LEU:N	1:F:82:LEU:HD12	2.18	0.58
1:G:140:ALA:HB3	1:G:144:MET:HE3	1.84	0.58
1:I:199:TRP:CD2	1:I:277:VAL:HG21	2.37	0.58
1:H:18:LYS:NZ	1:H:18:LYS:HA	2.18	0.58
1:A:176:MET:CE	1:A:195:LEU:HD13	2.32	0.58
1:E:176:MET:HE2	1:E:176:MET:HA	1.85	0.58
1:D:168:GLU:O	1:D:174:THR:HG21	2.04	0.58
1:H:75:GLU:HG3	1:H:127:THR:OG1	2.04	0.58
1:B:98:TYR:HE1	1:B:175:LEU:HD22	1.68	0.57
1:D:57:ALA:HA	1:D:188:GLN:HG3	1.85	0.57
1:I:137:THR:CG2	1:I:138:ILE:N	2.67	0.57
1:D:66:ASP:HB3	1:D:67:PRO:HD3	1.86	0.57
1:G:66:ASP:O	1:G:84:HIS:HB2	2.04	0.57
1:G:99:ARG:HG3	1:G:271:ASN:ND2	2.18	0.57
1:G:250:THR:CG2	1:G:264:ARG:HH21	2.16	0.57
1:A:99:ARG:NH2	1:A:268:LEU:HD12	2.20	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:129:ASN:HB3	1:F:216:ARG:NH1	2.20	0.57
1:A:285:ASP:O	1:A:286:LEU:HB3	2.05	0.57
1:B:135:PHE:HE2	1:B:176:MET:HE1	1.70	0.57
1:F:135:PHE:O	1:F:161:VAL:HG23	2.05	0.57
1:H:238:ASN:HB3	1:H:241:VAL:CG1	2.33	0.57
1:E:120:LEU:HD23	1:E:120:LEU:O	2.05	0.57
1:F:250:THR:HG23	1:F:264:ARG:HH21	1.70	0.57
1:I:168:GLU:O	1:I:174:THR:HG21	2.05	0.57
1:A:238:ASN:HB3	1:A:241:VAL:CG1	2.35	0.56
1:E:149:ASP:HB3	1:E:152:ILE:HB	1.85	0.56
1:C:120:LEU:C	1:C:120:LEU:HD23	2.30	0.56
1:H:120:LEU:C	1:H:120:LEU:HD23	2.30	0.56
1:I:28:TRP:CB	1:I:96:ARG:HG2	2.35	0.56
1:I:238:ASN:HB3	1:I:241:VAL:HG12	1.86	0.56
1:A:80:LEU:HD23	1:A:80:LEU:H	1.71	0.56
1:B:140:ALA:H	1:B:144:MET:HE3	1.69	0.56
1:F:57:ALA:HA	1:F:188:GLN:HG3	1.86	0.56
1:I:39:LYS:HZ1	1:I:257:GLY:H	1.53	0.56
1:G:133:VAL:HG22	1:G:214:LEU:HD13	1.88	0.56
1:B:60:ALA:HB1	1:B:187:PRO:HB2	1.87	0.56
1:B:120:LEU:HD23	1:B:120:LEU:O	2.06	0.56
1:C:65:THR:HB	1:C:192:PRO:CG	2.36	0.56
1:A:176:MET:HE2	1:A:195:LEU:CD1	2.36	0.56
1:A:239:PRO:HB3	1:A:267:THR:HB	1.88	0.56
1:H:120:LEU:HD23	1:H:120:LEU:O	2.05	0.56
1:B:69:THR:O	1:B:210:VAL:HG23	2.06	0.55
1:E:285:ASP:OD2	1:E:285:ASP:N	2.39	0.55
1:G:57:ALA:HA	1:G:188:GLN:CG	2.36	0.55
1:A:285:ASP:OD1	1:A:286:LEU:N	2.38	0.55
1:B:20:LEU:HG	1:B:21:GLU:N	2.14	0.55
1:F:211:SER:HB2	1:H:282:ARG:NH2	2.21	0.55
1:I:49:PHE:CZ	1:I:144:MET:HG2	2.41	0.55
1:J:63:LEU:C	1:J:63:LEU:HD13	2.32	0.55
1:E:39:LYS:NZ	1:E:257:GLY:H	2.05	0.55
1:E:140:ALA:HB3	1:E:144:MET:CE	2.32	0.55
1:F:234:TYR:CD2	1:F:286:LEU:HB3	2.41	0.55
1:C:93:LEU:N	1:C:93:LEU:HD12	2.21	0.55
1:C:149:ASP:OD2	1:C:152:ILE:HG12	2.06	0.55
1:C:65:THR:HB	1:C:192:PRO:HG3	1.88	0.54
1:J:133:VAL:HG22	1:J:214:LEU:HD13	1.89	0.54
1:D:96:ARG:NH1	1:D:172:PRO:O	2.40	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:225:LEU:O	1:D:228:PRO:HD3	2.07	0.54
1:F:171:HIS:ND1	1:F:172:PRO:HD2	2.22	0.54
1:J:138:ILE:HG12	1:J:139:TYR:N	2.22	0.54
1:D:75:GLU:HG3	1:D:127:THR:HG1	1.70	0.54
1:J:117:LEU:CD1	1:J:176:MET:HE3	2.31	0.54
1:F:82:LEU:N	1:F:82:LEU:CD1	2.71	0.54
1:G:60:ALA:HB1	1:G:187:PRO:HB2	1.90	0.54
1:H:87:LEU:HD13	1:H:193:VAL:CG2	2.36	0.54
1:D:263:GLN:HE21	1:D:264:ARG:H	1.54	0.54
1:G:238:ASN:HB3	1:G:241:VAL:HG12	1.90	0.54
1:D:238:ASN:HB3	1:D:241:VAL:CG1	2.37	0.54
1:E:30:ASN:ND2	1:E:32:LEU:H	2.03	0.54
1:D:28:TRP:CB	1:D:96:ARG:HG2	2.38	0.54
1:J:51:LEU:HD21	1:J:154:GLY:CA	2.38	0.53
1:A:87:LEU:HD13	1:A:193:VAL:CG2	2.37	0.53
1:C:76:VAL:HG13	1:C:124:ALA:O	2.08	0.53
1:F:87:LEU:HD22	1:F:193:VAL:HG21	1.89	0.53
1:F:140:ALA:N	1:F:144:MET:HE3	2.20	0.53
1:H:66:ASP:HB3	1:H:67:PRO:CD	2.37	0.53
1:J:238:ASN:HB3	1:J:241:VAL:CG1	2.37	0.53
1:F:20:LEU:HD11	1:F:113:ILE:HG21	1.91	0.53
1:F:57:ALA:HA	1:F:188:GLN:HG2	1.90	0.53
1:I:140:ALA:H	1:I:144:MET:CE	2.21	0.53
1:A:176:MET:HE2	1:A:195:LEU:HD13	1.90	0.53
1:C:75:GLU:HG3	1:C:127:THR:OG1	2.09	0.53
1:E:76:VAL:HG13	1:E:124:ALA:O	2.08	0.53
1:I:20:LEU:HD12	1:I:20:LEU:H	1.73	0.53
1:B:192:PRO:HG2	1:B:193:VAL:HG23	1.90	0.53
1:C:190:GLY:HA3	1:C:207:LYS:HB3	1.91	0.53
1:E:238:ASN:HB3	1:E:241:VAL:HG13	1.92	0.52
1:J:90:ARG:NH1	1:J:123:LEU:HD11	2.24	0.52
1:H:220:PRO:HB3	1:J:285:ASP:OD1	2.10	0.52
1:C:230:GLU:HG2	1:C:244:PRO:HG2	1.91	0.52
1:I:138:ILE:O	1:I:144:MET:HE1	2.10	0.52
1:F:98:TYR:HE1	1:F:175:LEU:HD22	1.74	0.52
1:I:20:LEU:HD23	1:I:88:THR:CG2	2.39	0.52
1:D:104:GLU:O	1:D:105:ALA:HB3	2.10	0.52
1:D:221:THR:HG22	1:D:224:ASN:H	1.74	0.52
1:G:162:GLU:OE2	1:G:221:THR:HG23	2.10	0.52
1:B:149:ASP:HB3	1:B:152:ILE:HB	1.92	0.52
1:E:98:TYR:HE1	1:E:175:LEU:HD22	1.75	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:99:ARG:HG3	1:B:271:ASN:HD21	1.75	0.51
1:I:149:ASP:HB3	1:I:152:ILE:HB	1.92	0.51
1:D:241:VAL:O	1:D:241:VAL:HG13	2.11	0.51
1:F:238:ASN:HB3	1:F:241:VAL:CG1	2.40	0.51
1:G:106:TRP:HA	1:G:252:ARG:O	2.10	0.51
1:J:57:ALA:CA	1:J:188:GLN:HG2	2.40	0.51
1:A:140:ALA:CB	1:A:144:MET:HE3	2.37	0.51
1:F:120:LEU:HD23	1:F:120:LEU:C	2.36	0.51
1:G:82:LEU:HD12	1:G:82:LEU:N	2.25	0.51
1:I:106:TRP:HA	1:I:252:ARG:O	2.11	0.51
1:J:76:VAL:HG13	1:J:124:ALA:O	2.10	0.51
1:B:51:LEU:HD21	1:B:154:GLY:CA	2.40	0.51
1:B:104:GLU:O	1:B:105:ALA:HB3	2.10	0.51
1:A:99:ARG:HH21	1:A:268:LEU:HD12	1.75	0.51
1:B:277:VAL:CG1	1:B:281:TYR:HE2	2.24	0.51
1:G:150:ARG:NH1	1:I:240:TYR:HA	2.26	0.51
1:B:137:THR:HG22	1:B:138:ILE:N	2.25	0.51
1:H:136:GLU:HB3	1:H:211:SER:HB3	1.92	0.51
1:B:171:HIS:CE1	1:B:172:PRO:HG2	2.46	0.51
1:C:190:GLY:CA	1:C:207:LYS:HB3	2.40	0.51
1:C:75:GLU:HB3	1:C:130:ALA:HB2	1.92	0.50
1:F:76:VAL:CG1	1:F:124:ALA:HB1	2.40	0.50
1:F:137:THR:HG22	1:F:138:ILE:N	2.26	0.50
1:F:138:ILE:HG12	1:F:139:TYR:N	2.26	0.50
1:J:241:VAL:HG13	1:J:241:VAL:O	2.10	0.50
1:A:48:GLU:HG3	1:A:158:TYR:CE2	2.46	0.50
1:G:239:PRO:HG3	1:G:267:THR:O	2.12	0.50
1:D:149:ASP:HB3	1:D:152:ILE:HB	1.93	0.50
1:A:192:PRO:HG2	1:A:193:VAL:HG23	1.93	0.50
1:G:39:LYS:NZ	1:G:257:GLY:O	2.34	0.50
1:J:65:THR:HB	1:J:192:PRO:HD3	1.93	0.50
1:A:190:GLY:HA3	1:A:207:LYS:HG2	1.92	0.50
1:B:176:MET:CE	1:B:195:LEU:HD13	2.41	0.50
1:D:140:ALA:N	1:D:144:MET:HE3	2.27	0.50
1:F:96:ARG:HD2	1:F:175:LEU:HD13	1.92	0.50
1:A:286:LEU:C	1:A:286:LEU:HD22	2.36	0.50
1:H:241:VAL:HG13	1:H:241:VAL:O	2.12	0.50
1:D:49:PHE:CE2	1:D:60:ALA:HA	2.47	0.50
1:J:98:TYR:HE1	1:J:175:LEU:HD22	1.77	0.50
1:A:72:ILE:HD13	1:A:212:ILE:HB	1.94	0.50
1:D:221:THR:HB	1:D:224:ASN:HB3	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:86:ASP:O	1:F:90:ARG:HB2	2.11	0.50
1:A:140:ALA:HB3	1:A:144:MET:CE	2.35	0.49
1:D:98:TYR:HE1	1:D:175:LEU:HD22	1.76	0.49
1:B:251:GLU:O	1:B:264:ARG:HA	2.12	0.49
1:F:104:GLU:O	1:F:105:ALA:HB3	2.12	0.49
1:F:223:TRP:HE3	1:F:231:TYR:CE2	2.31	0.49
1:C:93:LEU:N	1:C:93:LEU:CD1	2.74	0.49
1:D:173:LEU:HD11	1:D:276:GLN:NE2	2.28	0.49
1:G:239:PRO:HB3	1:G:267:THR:HB	1.93	0.49
1:E:137:THR:HG22	1:E:138:ILE:N	2.26	0.49
1:J:63:LEU:HD12	1:J:65:THR:CG2	2.41	0.49
1:D:28:TRP:HB3	1:D:96:ARG:HG2	1.93	0.49
1:F:171:HIS:CG	1:F:172:PRO:HD2	2.47	0.49
1:D:75:GLU:HB3	1:D:130:ALA:HB2	1.94	0.49
1:D:82:LEU:N	1:D:82:LEU:HD12	2.28	0.49
1:H:139:TYR:HB2	1:H:159:PRO:HB3	1.95	0.49
1:J:224:ASN:O	1:J:228:PRO:HG3	2.13	0.49
1:H:245:ARG:O	1:H:246:TRP:HB3	2.13	0.48
1:I:96:ARG:HD3	1:I:175:LEU:HD13	1.95	0.48
1:J:77:ALA:HB3	1:J:125:GLU:HB2	1.95	0.48
1:E:168:GLU:O	1:E:174:THR:HG21	2.13	0.48
1:F:250:THR:HG23	1:F:264:ARG:NH2	2.28	0.48
1:J:243:HIS:O	1:J:244:PRO:C	2.57	0.48
1:G:185:LEU:O	1:G:191:ALA:HB2	2.12	0.48
1:H:90:ARG:HB3	1:H:91:PHE:CD1	2.48	0.48
1:A:251:GLU:HA	1:A:264:ARG:HH22	1.78	0.48
1:B:210:VAL:O	1:B:210:VAL:HG22	2.12	0.48
1:C:80:LEU:HD23	1:C:80:LEU:N	2.28	0.48
1:C:286:LEU:HD13	1:C:286:LEU:N	2.28	0.48
1:D:176:MET:CE	1:D:195:LEU:HD13	2.43	0.48
1:A:80:LEU:HD23	1:A:80:LEU:N	2.29	0.48
1:A:211:SER:OG	1:C:282:ARG:HD3	2.12	0.48
1:B:106:TRP:HA	1:B:252:ARG:O	2.13	0.48
1:B:66:ASP:O	1:B:67:PRO:C	2.51	0.48
1:H:48:GLU:HG3	1:H:158:TYR:CE2	2.48	0.48
1:D:176:MET:HE2	1:D:195:LEU:HD13	1.95	0.48
1:D:263:GLN:HA	1:D:263:GLN:NE2	2.27	0.48
1:H:135:PHE:O	1:H:161:VAL:HG23	2.14	0.48
1:G:120:LEU:HD23	1:G:120:LEU:O	2.14	0.47
1:C:136:GLU:OE1	1:E:282:ARG:HD3	2.14	0.47
1:H:171:HIS:O	1:H:174:THR:HG23	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:87:LEU:HD13	1:J:193:VAL:HG22	1.93	0.47
1:B:176:MET:HE2	1:B:195:LEU:HD13	1.95	0.47
1:E:140:ALA:CB	1:E:144:MET:HE3	2.34	0.47
1:G:74:GLY:O	1:G:76:VAL:N	2.43	0.47
1:G:99:ARG:HD2	1:G:271:ASN:O	2.14	0.47
1:H:80:LEU:HD23	1:H:80:LEU:H	1.78	0.47
1:J:171:HIS:ND1	1:J:173:LEU:HB2	2.28	0.47
1:J:175:LEU:HD23	1:J:175:LEU:H	1.79	0.47
1:J:272:GLY:HA2	4:J:434:HOH:O	2.14	0.47
1:B:87:LEU:CD1	1:B:193:VAL:HG22	2.43	0.47
1:C:29:GLN:HG2	1:G:61:GLY:HA3	1.95	0.47
1:D:245:ARG:O	1:D:246:TRP:HB3	2.14	0.47
1:B:51:LEU:HD21	1:B:154:GLY:HA3	1.97	0.47
1:B:199:TRP:CD1	1:B:277:VAL:HG11	2.49	0.47
1:G:75:GLU:HG3	1:G:127:THR:OG1	2.14	0.47
1:J:49:PHE:CE2	1:J:60:ALA:HA	2.49	0.47
1:B:149:ASP:OD2	1:B:152:ILE:HG12	2.14	0.47
1:D:51:LEU:HD21	1:D:151:PHE:O	2.14	0.47
1:D:140:ALA:HB3	1:D:144:MET:HE2	1.97	0.47
1:F:168:GLU:O	1:F:174:THR:HG21	2.15	0.47
1:F:240:TYR:HA	1:I:150:ARG:NH1	2.30	0.47
1:G:66:ASP:O	1:G:68:TRP:N	2.47	0.47
1:E:104:GLU:O	1:E:105:ALA:HB3	2.13	0.47
1:I:60:ALA:HB1	1:I:187:PRO:HB2	1.95	0.47
1:G:66:ASP:O	1:G:67:PRO:C	2.52	0.47
1:I:112:TRP:CE3	1:I:177:THR:HG21	2.50	0.47
1:C:96:ARG:HD2	1:C:175:LEU:HD13	1.97	0.47
1:E:243:HIS:O	1:E:244:PRO:C	2.58	0.47
1:E:277:VAL:O	1:E:278:ALA:C	2.58	0.47
1:G:178:VAL:C	1:G:185:LEU:HG	2.40	0.47
1:J:168:GLU:OE2	1:J:281:TYR:OH	2.30	0.47
1:A:135:PHE:O	1:A:161:VAL:HG23	2.14	0.46
1:D:57:ALA:CA	1:D:188:GLN:HG2	2.42	0.46
1:E:75:GLU:HG3	1:E:127:THR:OG1	2.15	0.46
1:G:237:VAL:HG21	1:G:271:ASN:HB2	1.97	0.46
1:A:138:ILE:HG23	1:A:144:MET:CE	2.40	0.46
1:A:171:HIS:ND1	1:A:173:LEU:HB2	2.31	0.46
1:B:185:LEU:O	1:B:191:ALA:HB2	2.16	0.46
1:D:60:ALA:HB1	1:D:187:PRO:HB2	1.97	0.46
1:F:223:TRP:CE3	1:F:231:TYR:CE2	3.03	0.46
1:I:136:GLU:HB3	1:I:211:SER:HB3	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:175:LEU:HD23	1:I:196:ILE:HB	1.98	0.46
1:C:162:GLU:OE2	1:C:221:THR:HG23	2.16	0.46
1:F:138:ILE:HG23	1:F:144:MET:HE1	1.97	0.46
1:J:104:GLU:O	1:J:105:ALA:HB3	2.16	0.46
1:D:63:LEU:HD13	1:D:65:THR:CG2	2.45	0.46
1:B:133:VAL:HG22	1:B:214:LEU:HD13	1.97	0.46
1:H:117:LEU:HD13	1:H:176:MET:HE3	1.97	0.46
1:B:221:THR:HG22	1:B:224:ASN:H	1.81	0.46
1:F:109:VAL:HG23	1:F:253:PHE:HE1	1.80	0.46
1:H:211:SER:OG	1:J:282:ARG:HD3	2.16	0.46
1:H:264:ARG:O	1:H:265:GLN:HB3	2.15	0.46
1:I:133:VAL:HG22	1:I:214:LEU:CD1	2.44	0.46
1:I:221:THR:HB	1:I:224:ASN:HB3	1.98	0.46
1:B:117:LEU:CD1	1:B:176:MET:HE3	2.31	0.45
1:D:95:GLU:O	1:D:96:ARG:HG3	2.16	0.45
1:A:82:LEU:N	1:A:82:LEU:HD12	2.31	0.45
1:B:76:VAL:CG1	1:B:124:ALA:HB1	2.45	0.45
1:I:66:ASP:HB3	1:I:67:PRO:CD	2.44	0.45
1:A:76:VAL:HG11	1:A:124:ALA:HB1	1.97	0.45
1:A:277:VAL:O	1:A:277:VAL:HG12	2.15	0.45
1:B:25:PRO:HB2	1:B:28:TRP:CD1	2.51	0.45
1:B:171:HIS:CG	1:B:172:PRO:HD2	2.51	0.45
1:E:120:LEU:HD23	1:E:120:LEU:C	2.42	0.45
1:F:80:LEU:N	1:F:80:LEU:HD23	2.32	0.45
1:I:238:ASN:HB3	1:I:241:VAL:CG1	2.47	0.45
1:C:135:PHE:O	1:C:161:VAL:HG23	2.17	0.45
1:C:144:MET:HA	1:C:145:PRO:HD2	1.85	0.45
1:F:76:VAL:HG13	1:F:124:ALA:O	2.16	0.45
1:F:108:MET:HG2	1:F:254:ILE:HB	1.99	0.45
1:F:171:HIS:ND1	1:F:172:PRO:CD	2.80	0.45
1:J:47:TYR:C	1:J:49:PHE:N	2.73	0.45
1:A:81:THR:C	1:A:82:LEU:HD12	2.42	0.45
1:I:221:THR:O	1:I:225:LEU:HB2	2.17	0.45
1:B:138:ILE:HG23	1:B:144:MET:HE1	1.98	0.45
1:E:39:LYS:NZ	1:E:257:GLY:N	2.65	0.45
1:F:113:ILE:HD11	1:F:182:GLY:HA2	1.99	0.45
1:I:72:ILE:CD1	1:I:212:ILE:HB	2.41	0.45
1:J:72:ILE:HD13	1:J:212:ILE:HB	1.99	0.45
1:F:285:ASP:OD1	1:F:286:LEU:N	2.50	0.45
1:G:50:GLY:HA2	1:G:152:ILE:O	2.17	0.45
1:J:106:TRP:HA	1:J:252:ARG:O	2.18	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:112:TRP:CE3	1:B:177:THR:HG21	2.53	0.44
1:C:221:THR:HG22	1:C:224:ASN:H	1.82	0.44
1:D:207:LYS:O	1:D:208:SER:C	2.59	0.44
1:G:238:ASN:HB3	1:G:241:VAL:CG1	2.46	0.44
1:G:140:ALA:N	1:G:144:MET:HE3	2.32	0.44
1:I:135:PHE:O	1:I:161:VAL:HG23	2.16	0.44
1:D:81:THR:C	1:D:82:LEU:HD12	2.43	0.44
1:D:82:LEU:N	1:D:82:LEU:CD1	2.81	0.44
1:F:20:LEU:HD12	1:F:20:LEU:C	2.43	0.44
1:F:57:ALA:CA	1:F:188:GLN:HG2	2.47	0.44
1:G:90:ARG:HB3	1:G:91:PHE:CD1	2.52	0.44
1:B:76:VAL:HG12	1:B:78:LYS:N	2.33	0.44
1:C:82:LEU:N	1:C:82:LEU:HD12	2.32	0.44
1:F:70:LEU:C	1:F:70:LEU:CD2	2.91	0.44
1:B:82:LEU:HD12	1:B:82:LEU:N	2.32	0.44
1:C:63:LEU:CD1	1:C:65:THR:HG22	2.48	0.44
1:C:171:HIS:CG	1:C:172:PRO:HD2	2.53	0.44
1:D:199:TRP:CD1	1:D:277:VAL:HG11	2.53	0.44
1:F:101:ARG:NE	1:F:251:GLU:OE2	2.51	0.44
1:I:80:LEU:N	1:I:80:LEU:HD23	2.33	0.44
1:I:120:LEU:O	1:I:120:LEU:HD23	2.18	0.44
1:I:183:LYS:H	1:I:183:LYS:HG2	1.65	0.44
1:J:190:GLY:CA	1:J:207:LYS:HB3	2.48	0.44
1:B:133:VAL:O	1:B:163:GLY:HA2	2.18	0.44
1:B:176:MET:HE2	1:B:195:LEU:CD1	2.47	0.44
1:D:76:VAL:HG11	1:D:124:ALA:HB1	1.99	0.44
1:J:135:PHE:O	1:J:161:VAL:HG23	2.17	0.44
1:C:106:TRP:HA	1:C:252:ARG:O	2.18	0.44
1:G:76:VAL:HG11	1:G:124:ALA:HB1	1.99	0.44
1:H:82:LEU:N	1:H:82:LEU:CD1	2.81	0.44
1:C:57:ALA:HA	1:C:188:GLN:HG3	1.96	0.43
1:D:87:LEU:CD1	1:D:193:VAL:HG22	2.41	0.43
1:D:137:THR:HG22	1:D:138:ILE:N	2.32	0.43
1:D:165:ARG:NH2	1:D:167:ASP:OD2	2.47	0.43
1:E:57:ALA:HA	1:E:188:GLN:HG3	1.96	0.43
1:I:75:GLU:HB3	1:I:130:ALA:HB2	1.98	0.43
1:I:87:LEU:CD1	1:I:193:VAL:HG22	2.46	0.43
1:J:51:LEU:HD21	1:J:154:GLY:HA3	1.98	0.43
1:D:52:ASP:HB2	1:D:55:ASP:CG	2.43	0.43
1:H:137:THR:HG22	1:H:138:ILE:N	2.33	0.43
1:H:196:ILE:HG22	1:H:198:PRO:HD3	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:190:GLY:HA2	1:E:207:LYS:HB3	2.00	0.43
1:B:72:ILE:HD13	1:B:212:ILE:HB	2.00	0.43
1:D:76:VAL:CG1	1:D:124:ALA:HB1	2.49	0.43
1:A:230:GLU:HG2	1:A:244:PRO:HG2	1.99	0.43
1:D:188:GLN:H	1:D:188:GLN:CD	2.26	0.43
1:E:75:GLU:CG	1:E:129:ASN:HB2	2.49	0.43
1:G:98:TYR:CE1	1:G:198:PRO:HG3	2.53	0.43
1:G:149:ASP:HB3	1:G:152:ILE:HB	2.01	0.43
1:B:66:ASP:O	1:B:68:TRP:N	2.52	0.43
1:C:86:ASP:O	1:C:90:ARG:HB2	2.18	0.43
1:E:224:ASN:O	1:E:228:PRO:HG3	2.19	0.43
1:F:140:ALA:H	1:F:144:MET:CE	2.24	0.43
1:G:120:LEU:HD23	1:G:120:LEU:C	2.43	0.43
1:H:104:GLU:O	1:H:105:ALA:HB3	2.19	0.43
1:A:51:LEU:HD21	1:A:154:GLY:N	2.34	0.43
1:F:186:PRO:HG2	1:F:189:ASN:OD1	2.19	0.43
1:H:80:LEU:HD23	1:H:80:LEU:N	2.34	0.43
1:I:66:ASP:O	1:I:67:PRO:C	2.56	0.43
1:A:238:ASN:CG	1:A:241:VAL:HG12	2.44	0.43
1:B:80:LEU:N	1:B:80:LEU:HD23	2.34	0.43
1:C:163:GLY:HA3	1:C:219:PRO:HG2	2.00	0.43
1:D:140:ALA:HB3	1:D:144:MET:CE	2.49	0.43
1:I:104:GLU:O	1:I:105:ALA:HB3	2.18	0.43
1:B:238:ASN:CB	1:B:241:VAL:HG12	2.47	0.42
1:D:63:LEU:HD13	1:D:65:THR:HG22	2.00	0.42
1:G:82:LEU:HB3	1:G:86:ASP:HB2	2.00	0.42
1:G:90:ARG:HB3	1:G:91:PHE:CE1	2.53	0.42
1:J:63:LEU:HD13	1:J:64:LYS:N	2.34	0.42
1:J:214:LEU:N	1:J:214:LEU:CD2	2.82	0.42
1:C:149:ASP:HB3	1:C:152:ILE:HB	2.00	0.42
1:E:204:LYS:HE3	1:E:204:LYS:HB3	1.89	0.42
1:F:49:PHE:CE2	1:F:60:ALA:HA	2.53	0.42
1:G:160:TYR:CD1	1:G:206:ILE:HG13	2.53	0.42
1:H:55:ASP:HB2	1:H:56:PRO:HD3	2.01	0.42
1:H:144:MET:HA	1:H:145:PRO:HD2	1.83	0.42
1:I:20:LEU:HD12	1:I:20:LEU:N	2.33	0.42
1:J:263:GLN:HG3	1:J:264:ARG:H	1.83	0.42
1:A:71:LYS:HE3	1:A:71:LYS:HB2	1.84	0.42
1:F:133:VAL:HG21	1:F:169:ALA:CB	2.49	0.42
1:F:199:TRP:CD1	1:F:277:VAL:HG11	2.54	0.42
1:H:86:ASP:O	1:H:90:ARG:HB2	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:227:ALA:N	1:J:228:PRO:HD3	2.34	0.42
1:C:171:HIS:O	1:C:174:THR:HG23	2.19	0.42
1:F:238:ASN:ND2	1:I:157:LYS:HE3	2.34	0.42
1:I:28:TRP:CD1	1:I:96:ARG:HG2	2.52	0.42
1:J:175:LEU:CD2	1:J:196:ILE:HB	2.50	0.42
1:C:286:LEU:N	1:C:286:LEU:CD1	2.82	0.42
1:B:171:HIS:HE1	1:B:173:LEU:HD13	1.84	0.42
1:C:90:ARG:HB3	1:C:91:PHE:CD1	2.54	0.42
1:C:104:GLU:O	1:C:105:ALA:HB3	2.19	0.42
1:D:239:PRO:HB3	1:D:267:THR:HB	2.01	0.42
1:G:49:PHE:CE2	1:G:60:ALA:HA	2.54	0.42
1:G:140:ALA:HB3	1:G:144:MET:CE	2.48	0.42
1:H:72:ILE:HD13	1:H:212:ILE:HB	2.00	0.42
1:H:286:LEU:H	1:H:286:LEU:CD2	2.26	0.42
1:F:101:ARG:HD3	1:F:251:GLU:OE2	2.20	0.42
1:G:47:TYR:CE2	1:G:51:LEU:HD22	2.55	0.42
1:I:94:GLU:O	1:I:113:ILE:HA	2.20	0.42
1:I:135:PHE:CB	1:I:206:ILE:HG21	2.49	0.42
1:C:24:LYS:O	1:C:25:PRO:C	2.63	0.42
1:D:71:LYS:HE3	1:D:71:LYS:HB2	1.76	0.42
1:J:47:TYR:O	1:J:48:GLU:C	2.62	0.42
1:A:99:ARG:NH2	1:A:268:LEU:CD1	2.83	0.42
1:A:237:VAL:HG21	1:A:271:ASN:HB2	2.01	0.42
1:B:157:LYS:HE3	1:D:238:ASN:ND2	2.35	0.42
1:B:221:THR:O	1:B:225:LEU:HD13	2.19	0.42
1:E:98:TYR:CE1	1:E:175:LEU:HD22	2.55	0.42
1:I:65:THR:HB	1:I:192:PRO:HG3	2.01	0.42
1:A:100:MET:HG2	4:A:428:HOH:O	2.20	0.41
1:B:52:ASP:HB2	1:B:55:ASP:CG	2.45	0.41
1:B:263:GLN:OE1	1:B:263:GLN:HA	2.20	0.41
1:E:284:LEU:C	1:E:285:ASP:OD2	2.63	0.41
1:G:137:THR:HG22	1:G:138:ILE:N	2.35	0.41
1:H:51:LEU:HD21	1:H:154:GLY:CA	2.49	0.41
1:I:108:MET:HG2	1:I:254:ILE:HB	2.02	0.41
1:C:48:GLU:OE2	1:C:48:GLU:N	2.46	0.41
1:D:66:ASP:O	1:D:84:HIS:HB2	2.20	0.41
1:G:76:VAL:CG1	1:G:77:ALA:N	2.83	0.41
1:G:277:VAL:O	1:G:277:VAL:CG1	2.68	0.41
1:B:135:PHE:O	1:B:161:VAL:HG23	2.20	0.41
1:B:197:VAL:HA	1:B:198:PRO:HD2	1.86	0.41
1:E:171:HIS:ND1	1:E:173:LEU:HB2	2.36	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:149:ASP:HB3	1:F:152:ILE:HB	2.03	0.41
1:F:175:LEU:HD23	1:F:196:ILE:HB	2.03	0.41
1:G:140:ALA:H	1:G:144:MET:HE3	1.85	0.41
1:B:20:LEU:CG	1:B:21:GLU:H	2.12	0.41
1:F:71:LYS:HB2	1:F:71:LYS:HE3	1.81	0.41
1:H:106:TRP:HA	1:H:252:ARG:O	2.20	0.41
1:H:146:GLY:C	1:H:148:GLN:H	2.29	0.41
1:I:71:LYS:HB2	1:I:71:LYS:HE3	1.84	0.41
1:A:74:GLY:O	1:A:76:VAL:N	2.47	0.41
1:A:277:VAL:O	1:A:277:VAL:CG1	2.68	0.41
1:E:39:LYS:HZ1	1:E:257:GLY:H	1.68	0.41
1:F:194:ARG:HD3	1:F:194:ARG:C	2.46	0.41
1:I:137:THR:CG2	1:I:138:ILE:H	2.32	0.41
1:J:101:ARG:NH1	1:J:105:ALA:HA	2.35	0.41
1:D:47:TYR:N	1:D:47:TYR:CD1	2.89	0.41
1:J:137:THR:HG22	1:J:138:ILE:N	2.35	0.41
1:E:176:MET:CE	1:E:195:LEU:HD13	2.51	0.41
1:D:224:ASN:O	1:D:228:PRO:HG3	2.20	0.41
1:F:285:ASP:C	1:F:286:LEU:HG	2.46	0.41
1:G:161:VAL:HG22	1:G:162:GLU:N	2.36	0.41
1:H:65:THR:HB	1:H:192:PRO:HD3	2.03	0.41
1:H:93:LEU:N	1:H:93:LEU:HD12	2.36	0.41
1:J:214:LEU:N	1:J:214:LEU:HD22	2.35	0.41
1:A:251:GLU:C	1:A:264:ARG:HH21	2.28	0.41
1:A:285:ASP:OD1	1:A:286:LEU:HD13	2.20	0.41
1:B:239:PRO:HB3	1:B:267:THR:HB	2.03	0.41
1:H:76:VAL:CG1	1:H:77:ALA:N	2.83	0.41
1:B:120:LEU:HD23	1:B:120:LEU:C	2.45	0.41
1:F:85:ASP:O	1:F:89:ARG:HB2	2.21	0.41
1:I:70:LEU:HD12	1:I:193:VAL:HG21	2.03	0.41
1:I:144:MET:HA	1:I:145:PRO:HD2	1.87	0.41
1:I:269:LEU:O	1:I:270:PHE:HB2	2.20	0.41
1:J:162:GLU:OE1	1:J:204:LYS:HD2	2.21	0.41
1:B:66:ASP:HB3	1:B:67:PRO:CD	2.51	0.40
1:B:203:PHE:HB2	1:B:223:TRP:CD1	2.55	0.40
1:E:36:PRO:HG2	1:E:39:LYS:HD2	2.04	0.40
1:E:39:LYS:HZ1	1:E:257:GLY:N	2.19	0.40
1:J:138:ILE:HG23	1:J:144:MET:HE1	2.02	0.40
1:E:28:TRP:CH2	1:E:173:LEU:CD1	3.04	0.40
1:H:237:VAL:HG21	1:H:271:ASN:HB2	2.04	0.40
1:I:24:LYS:O	1:I:25:PRO:C	2.64	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:99:ARG:HG3	1:J:271:ASN:ND2	2.36	0.40
1:E:74:GLY:O	1:E:76:VAL:N	2.52	0.40
1:E:173:LEU:HA	1:E:173:LEU:HD12	1.78	0.40
1:G:100:MET:HG2	4:G:421:HOH:O	2.21	0.40
1:H:47:TYR:C	1:H:49:PHE:N	2.78	0.40
1:H:171:HIS:CG	1:H:172:PRO:HD2	2.57	0.40
1:H:277:VAL:O	1:H:278:ALA:C	2.63	0.40
1:J:190:GLY:HA2	1:J:207:LYS:HB3	2.03	0.40
1:A:51:LEU:HD21	1:A:154:GLY:HA3	2.00	0.40
1:B:70:LEU:HD23	1:B:70:LEU:C	2.47	0.40
1:F:75:GLU:HB3	1:F:130:ALA:HB2	2.02	0.40
1:A:66:ASP:CG	1:I:33:PRO:HB3	2.46	0.40
1:D:144:MET:HA	1:D:145:PRO:HD2	1.88	0.40
1:H:204:LYS:NZ	1:H:233:PHE:O	2.52	0.40
1:I:95:GLU:C	1:I:96:ARG:HG3	2.47	0.40
1:I:161:VAL:HG22	1:I:162:GLU:N	2.36	0.40
1:J:87:LEU:HD13	1:J:193:VAL:HG23	2.00	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	259/298 (87%)	249 (96%)	9 (4%)	1 (0%)	30	34
1	B	258/298 (87%)	242 (94%)	15 (6%)	1 (0%)	30	34
1	C	260/298 (87%)	246 (95%)	13 (5%)	1 (0%)	30	34
1	D	259/298 (87%)	248 (96%)	11 (4%)	0	100	100
1	E	258/298 (87%)	241 (93%)	17 (7%)	0	100	100
1	F	258/298 (87%)	243 (94%)	15 (6%)	0	100	100
1	G	257/298 (86%)	242 (94%)	13 (5%)	2 (1%)	16	16

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	H	260/298 (87%)	236 (91%)	23 (9%)	1 (0%)	30	34
1	I	258/298 (87%)	239 (93%)	18 (7%)	1 (0%)	30	34
1	J	258/298 (87%)	244 (95%)	14 (5%)	0	100	100
All	All	2585/2980 (87%)	2430 (94%)	148 (6%)	7 (0%)	36	42

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	192	PRO
1	G	75	GLU
1	A	75	GLU
1	B	66	ASP
1	G	66	ASP
1	H	265	GLN
1	I	66	ASP

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	219/249 (88%)	205 (94%)	14 (6%)	16	19
1	B	219/249 (88%)	214 (98%)	5 (2%)	44	59
1	C	220/249 (88%)	206 (94%)	14 (6%)	16	19
1	D	219/249 (88%)	208 (95%)	11 (5%)	22	28
1	E	218/249 (88%)	209 (96%)	9 (4%)	27	37
1	F	218/249 (88%)	205 (94%)	13 (6%)	17	21
1	G	218/249 (88%)	210 (96%)	8 (4%)	30	41
1	H	220/249 (88%)	208 (94%)	12 (6%)	19	24
1	I	218/249 (88%)	212 (97%)	6 (3%)	38	52
1	J	218/249 (88%)	209 (96%)	9 (4%)	27	37
All	All	2187/2490 (88%)	2086 (95%)	101 (5%)	24	32

All (101) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	20	LEU
1	A	70	LEU
1	A	76	VAL
1	A	120	LEU
1	A	173	LEU
1	A	174	THR
1	A	175	LEU
1	A	193	VAL
1	A	210	VAL
1	A	221	THR
1	A	241	VAL
1	A	264	ARG
1	A	279	SER
1	A	286	LEU
1	B	125	GLU
1	B	165	ARG
1	B	174	THR
1	B	193	VAL
1	B	221	THR
1	C	20	LEU
1	C	70	LEU
1	C	173	LEU
1	C	174	THR
1	C	175	LEU
1	C	188	GLN
1	C	193	VAL
1	C	197	VAL
1	C	210	VAL
1	C	214	LEU
1	C	225	LEU
1	C	256	SER
1	C	263	GLN
1	C	286	LEU
1	D	51	LEU
1	D	63	LEU
1	D	70	LEU
1	D	76	VAL
1	D	82	LEU
1	D	93	LEU
1	D	99	ARG
1	D	174	THR
1	D	193	VAL

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Mol	Chain	Res	Type
1	D	218	ARG
1	D	225	LEU
1	E	20	LEU
1	E	76	VAL
1	E	93	LEU
1	E	173	LEU
1	E	174	THR
1	E	193	VAL
1	E	221	THR
1	E	241	VAL
1	E	285	ASP
1	F	38	ASP
1	F	63	LEU
1	F	70	LEU
1	F	76	VAL
1	F	81	THR
1	F	82	LEU
1	F	93	LEU
1	F	173	LEU
1	F	174	THR
1	F	188	GLN
1	F	221	THR
1	F	279	SER
1	F	286	LEU
1	G	64	LYS
1	G	70	LEU
1	G	175	LEU
1	G	193	VAL
1	G	211	SER
1	G	216	ARG
1	G	221	THR
1	G	225	LEU
1	H	18	LYS
1	H	20	LEU
1	H	70	LEU
1	H	76	VAL
1	H	82	LEU
1	H	90	ARG
1	H	173	LEU
1	H	174	THR
1	H	175	LEU
1	H	188	GLN

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Mol	Chain	Res	Type
1	H	221	THR
1	H	286	LEU
1	I	20	LEU
1	I	175	LEU
1	I	183	LYS
1	I	193	VAL
1	I	218	ARG
1	I	277	VAL
1	J	76	VAL
1	J	81	THR
1	J	93	LEU
1	J	99	ARG
1	J	120	LEU
1	J	173	LEU
1	J	174	THR
1	J	214	LEU
1	J	221	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	148	GLN
1	B	29	GLN
1	C	29	GLN
1	C	44	ASN
1	C	143	GLN
1	C	148	GLN
1	D	44	ASN
1	D	84	HIS
1	D	143	GLN
1	D	148	GLN
1	D	263	GLN
1	E	29	GLN
1	E	30	ASN
1	E	44	ASN
1	E	189	ASN
1	F	143	GLN
1	G	84	HIS
1	G	276	GLN
1	H	143	GLN
1	I	31	ASN
1	I	143	GLN

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Mol	Chain	Res	Type
1	I	148	GLN
1	J	143	GLN
1	J	263	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 20 ligands modelled in this entry, 10 are monoatomic - leaving 10 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	$\# Z > 2$
3	MTE	B	302	2	21,26,26	3.40	12 (57%)	19,40,40	6.23	15 (78%)
3	MTE	J	310	2	21,26,26	3.65	9 (42%)	19,40,40	6.80	12 (63%)
3	MTE	D	304	2	21,26,26	3.31	9 (42%)	19,40,40	6.28	14 (73%)
3	MTE	H	308	2	21,26,26	3.67	8 (38%)	19,40,40	6.18	16 (84%)
3	MTE	I	309	2	21,26,26	3.80	10 (47%)	19,40,40	6.30	12 (63%)
3	MTE	C	303	2	21,26,26	3.26	10 (47%)	19,40,40	6.87	13 (68%)
3	MTE	F	306	2	21,26,26	3.55	9 (42%)	19,40,40	6.51	15 (78%)
3	MTE	G	307	2	21,26,26	3.45	10 (47%)	19,40,40	6.55	11 (57%)
3	MTE	A	301	2	21,26,26	3.46	10 (47%)	19,40,40	6.50	12 (63%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	MTE	E	305	2	21,26,26	3.17	10 (47%)	19,40,40	6.74	14 (73%)

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
3	MTE	B	302	2	1/1/6/8	0/6/34/34	0/3/3/3
3	MTE	J	310	2	1/1/6/8	0/6/34/34	0/3/3/3
3	MTE	D	304	2	1/1/6/8	0/6/34/34	0/3/3/3
3	MTE	H	308	2	1/1/6/8	0/6/34/34	0/3/3/3
3	MTE	I	309	2	1/1/6/8	0/6/34/34	0/3/3/3
3	MTE	C	303	2	1/1/6/8	0/6/34/34	0/3/3/3
3	MTE	F	306	2	1/1/6/8	0/6/34/34	0/3/3/3
3	MTE	G	307	2	1/1/6/8	0/6/34/34	0/3/3/3
3	MTE	A	301	2	1/1/6/8	0/6/34/34	0/3/3/3
3	MTE	E	305	2	1/1/6/8	0/6/34/34	0/3/3/3

All (97) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	I	309	MTE	C4'-C3'	-9.79	1.38	1.51
3	H	308	MTE	C9-C10	9.74	1.55	1.38
3	F	306	MTE	C9-C10	9.28	1.54	1.38
3	H	308	MTE	C4'-C3'	-9.15	1.39	1.51
3	G	307	MTE	C9-C10	9.03	1.54	1.38
3	J	310	MTE	C9-C10	9.00	1.54	1.38
3	J	310	MTE	C4'-C3'	-8.93	1.39	1.51
3	C	303	MTE	C9-C10	8.70	1.53	1.38
3	F	306	MTE	C4'-C3'	-8.69	1.39	1.51
3	I	309	MTE	C9-C10	8.66	1.53	1.38
3	G	307	MTE	C4'-C3'	-8.60	1.39	1.51
3	E	305	MTE	C9-C10	8.59	1.53	1.38
3	B	302	MTE	C9-C10	8.55	1.53	1.38
3	D	304	MTE	C9-C10	8.45	1.53	1.38
3	A	301	MTE	C4'-C3'	-8.12	1.40	1.51
3	B	302	MTE	C4'-C3'	-7.60	1.41	1.51
3	C	303	MTE	C4'-C3'	-7.58	1.41	1.51
3	A	301	MTE	C9-C10	7.52	1.51	1.38
3	D	304	MTE	C4'-C3'	-7.51	1.41	1.51
3	E	305	MTE	C4'-C3'	-7.34	1.41	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	I	309	MTE	C6-N5	6.29	1.53	1.46
3	I	309	MTE	P-O4'	-5.81	1.42	1.60
3	H	308	MTE	P-O4'	-5.53	1.42	1.60
3	J	310	MTE	C6-N5	5.41	1.52	1.46
3	J	310	MTE	P-O4'	-5.32	1.43	1.60
3	C	303	MTE	P-O4'	-5.20	1.43	1.60
3	F	306	MTE	C6-N5	5.08	1.52	1.46
3	F	306	MTE	P-O4'	-5.05	1.44	1.60
3	A	301	MTE	P-O1P	-5.03	1.34	1.50
3	B	302	MTE	P-O4'	-4.76	1.45	1.60
3	B	302	MTE	C6-N5	4.66	1.51	1.46
3	J	310	MTE	O3'-C7	4.65	1.50	1.43
3	A	301	MTE	C6-N5	4.59	1.51	1.46
3	D	304	MTE	P-O4'	-4.58	1.45	1.60
3	A	301	MTE	P-O4'	-4.57	1.45	1.60
3	E	305	MTE	P-O4'	-4.41	1.46	1.60
3	G	307	MTE	C6-N5	4.06	1.51	1.46
3	D	304	MTE	O3'-C7	4.04	1.49	1.43
3	G	307	MTE	C10-N8	3.86	1.39	1.35
3	A	301	MTE	C2-N3	-3.85	1.28	1.37
3	G	307	MTE	P-O4'	-3.82	1.48	1.60
3	H	308	MTE	C6-N5	3.78	1.50	1.46
3	H	308	MTE	P-O1P	-3.67	1.39	1.50
3	B	302	MTE	P-O1P	-3.59	1.39	1.50
3	E	305	MTE	C7-C6	3.49	1.56	1.53
3	H	308	MTE	O3'-C7	3.48	1.48	1.43
3	G	307	MTE	C9-N5	3.47	1.46	1.37
3	E	305	MTE	C6-N5	3.42	1.50	1.46
3	D	304	MTE	C6-N5	3.32	1.50	1.46
3	G	307	MTE	C2-N3	-3.31	1.29	1.37
3	D	304	MTE	O3'-C3'	3.25	1.49	1.43
3	F	306	MTE	C9-N5	3.24	1.45	1.37
3	J	310	MTE	C9-N5	3.19	1.45	1.37
3	I	309	MTE	C9-N5	3.18	1.45	1.37
3	B	302	MTE	C9-N5	3.15	1.45	1.37
3	F	306	MTE	P-O1P	-3.13	1.40	1.50
3	C	303	MTE	C9-N5	3.10	1.45	1.37
3	C	303	MTE	C6-N5	3.10	1.50	1.46
3	A	301	MTE	C9-N5	3.05	1.45	1.37
3	J	310	MTE	C2-N3	-3.01	1.30	1.37
3	I	309	MTE	C2-N3	-3.01	1.30	1.37
3	D	304	MTE	C2-N3	-2.93	1.30	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	H	308	MTE	C9-N5	2.76	1.44	1.37
3	A	301	MTE	O3'-C3'	2.76	1.48	1.43
3	C	303	MTE	O3'-C7	2.74	1.47	1.43
3	B	302	MTE	C2-N3	-2.73	1.31	1.37
3	D	304	MTE	P-O1P	-2.72	1.42	1.50
3	E	305	MTE	P-O1P	-2.72	1.42	1.50
3	I	309	MTE	P-O1P	-2.72	1.42	1.50
3	B	302	MTE	O3'-C7	2.69	1.47	1.43
3	C	303	MTE	C2-N3	-2.65	1.31	1.37
3	F	306	MTE	O3'-C7	2.59	1.47	1.43
3	J	310	MTE	P-O1P	-2.58	1.42	1.50
3	I	309	MTE	C10-N8	2.54	1.38	1.35
3	F	306	MTE	O3'-C3'	2.50	1.48	1.43
3	C	303	MTE	P-O1P	-2.47	1.42	1.50
3	I	309	MTE	O3'-C7	2.46	1.47	1.43
3	A	301	MTE	O3'-C7	2.44	1.47	1.43
3	E	305	MTE	C2-N3	-2.43	1.31	1.37
3	H	308	MTE	C2-N3	-2.41	1.31	1.37
3	B	302	MTE	C2-N1	2.39	1.39	1.33
3	D	304	MTE	C10-N1	-2.37	1.32	1.36
3	B	302	MTE	O4'-C4'	2.32	1.53	1.44
3	G	307	MTE	P-O1P	-2.26	1.43	1.50
3	I	309	MTE	C2-N1	2.24	1.38	1.33
3	J	310	MTE	C2-N2	-2.22	1.29	1.34
3	E	305	MTE	C10-N1	-2.22	1.32	1.36
3	G	307	MTE	C2-N1	2.20	1.38	1.33
3	E	305	MTE	C2-N2	-2.19	1.29	1.34
3	B	302	MTE	C10-N8	2.19	1.37	1.35
3	E	305	MTE	O4'-C4'	2.18	1.52	1.44
3	A	301	MTE	C2-N2	-2.17	1.29	1.34
3	C	303	MTE	O3'-C3'	2.11	1.47	1.43
3	F	306	MTE	C2-N3	-2.09	1.32	1.37
3	C	303	MTE	O4'-C4'	2.08	1.52	1.44
3	G	307	MTE	O3'-C3'	2.08	1.47	1.43
3	B	302	MTE	P-O2P	-2.02	1.47	1.54

All (134) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	305	MTE	O3'-C7-C6	-24.51	92.61	108.96
3	J	310	MTE	O3'-C7-C6	-24.48	92.63	108.96
3	C	303	MTE	O3'-C7-C6	-24.35	92.72	108.96

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	G	307	MTE	O3'-C7-C6	-23.05	93.59	108.96
3	A	301	MTE	O3'-C7-C6	-22.85	93.73	108.96
3	F	306	MTE	O3'-C7-C6	-22.67	93.84	108.96
3	D	304	MTE	O3'-C7-C6	-21.85	94.39	108.96
3	I	309	MTE	O3'-C7-C6	-21.57	94.58	108.96
3	B	302	MTE	O3'-C7-C6	-21.24	94.80	108.96
3	H	308	MTE	O3'-C7-C6	-20.93	95.00	108.96
3	C	303	MTE	C2-N1-C10	10.36	131.64	113.36
3	D	304	MTE	C2-N1-C10	10.31	131.56	113.36
3	A	301	MTE	C2-N1-C10	10.30	131.54	113.36
3	I	309	MTE	C2-N1-C10	10.13	131.25	113.36
3	G	307	MTE	C2-N1-C10	10.12	131.23	113.36
3	F	306	MTE	C2-N1-C10	10.00	131.01	113.36
3	H	308	MTE	C2-N1-C10	9.97	130.96	113.36
3	E	305	MTE	C2-N1-C10	9.93	130.89	113.36
3	J	310	MTE	C2-N1-C10	9.92	130.87	113.36
3	B	302	MTE	C2-N1-C10	9.76	130.59	113.36
3	B	302	MTE	O4-C4-C9	-7.12	110.10	127.26
3	F	306	MTE	O4-C4-C9	-7.03	110.31	127.26
3	J	310	MTE	O4-C4-C9	-6.82	110.81	127.26
3	C	303	MTE	O4-C4-C9	-6.77	110.93	127.26
3	I	309	MTE	O4-C4-C9	-6.71	111.08	127.26
3	H	308	MTE	O4-C4-C9	-6.71	111.08	127.26
3	D	304	MTE	O4-C4-C9	-6.66	111.20	127.26
3	G	307	MTE	O4-C4-C9	-6.53	111.52	127.26
3	E	305	MTE	O4-C4-C9	-6.49	111.61	127.26
3	A	301	MTE	O4-C4-C9	-6.43	111.76	127.26
3	C	303	MTE	C7-C6-N5	5.62	113.39	107.87
3	H	308	MTE	C7-C6-N5	5.37	113.14	107.87
3	D	304	MTE	C7-C6-N5	5.33	113.10	107.87
3	G	307	MTE	C7-C6-N5	5.32	113.09	107.87
3	J	310	MTE	C7-C6-N5	5.29	113.06	107.87
3	G	307	MTE	O3'-C7-N8	5.29	113.41	108.61
3	C	303	MTE	O3'-C7-N8	5.20	113.32	108.61
3	A	301	MTE	C7-C6-N5	5.18	112.95	107.87
3	I	309	MTE	C7-C6-N5	5.13	112.91	107.87
3	B	302	MTE	C7-C6-N5	5.05	112.83	107.87
3	A	301	MTE	O3'-C7-N8	4.91	113.06	108.61
3	E	305	MTE	C7-C6-N5	4.66	112.44	107.87
3	J	310	MTE	O3'-C7-N8	4.59	112.78	108.61
3	I	309	MTE	C4-C9-N5	4.58	128.76	116.27
3	F	306	MTE	C4-C9-N5	4.53	128.62	116.27

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	G	307	MTE	C4-C9-N5	4.53	128.61	116.27
3	J	310	MTE	C4-C9-N5	4.52	128.58	116.27
3	I	309	MTE	O3'-C7-N8	4.49	112.68	108.61
3	A	301	MTE	C4-C9-N5	4.45	128.40	116.27
3	F	306	MTE	C7-C6-N5	4.43	112.22	107.87
3	B	302	MTE	C4-C9-N5	4.40	128.25	116.27
3	D	304	MTE	C4-C9-N5	4.39	128.24	116.27
3	H	308	MTE	C4-C9-N5	4.39	128.23	116.27
3	I	309	MTE	C9-C4-N3	4.38	124.17	112.13
3	F	306	MTE	C9-C4-N3	4.36	124.09	112.13
3	E	305	MTE	C4-C9-N5	4.35	128.14	116.27
3	B	302	MTE	C9-C4-N3	4.34	124.05	112.13
3	J	310	MTE	C9-C4-N3	4.32	124.01	112.13
3	C	303	MTE	C4-C9-N5	4.28	127.94	116.27
3	A	301	MTE	C9-C4-N3	4.16	123.56	112.13
3	F	306	MTE	O3'-C7-N8	4.13	112.35	108.61
3	C	303	MTE	C9-C4-N3	4.12	123.45	112.13
3	H	308	MTE	C9-C4-N3	4.07	123.30	112.13
3	G	307	MTE	C9-C4-N3	4.03	123.19	112.13
3	E	305	MTE	C9-C4-N3	4.02	123.18	112.13
3	D	304	MTE	C9-C4-N3	4.01	123.14	112.13
3	B	302	MTE	O4'-P-O1P	3.77	116.63	106.44
3	H	308	MTE	O3'-C7-N8	3.61	111.88	108.61
3	H	308	MTE	O4'-P-O1P	3.52	115.94	106.44
3	C	303	MTE	O4'-P-O1P	3.42	115.68	106.44
3	B	302	MTE	O3P-P-O2P	3.29	120.16	107.80
3	E	305	MTE	O4'-P-O1P	3.28	115.30	106.44
3	F	306	MTE	O4'-P-O1P	3.26	115.26	106.44
3	I	309	MTE	O4'-P-O1P	3.25	115.22	106.44
3	I	309	MTE	N3-C2-N1	-3.24	117.39	123.32
3	B	302	MTE	O3'-C7-N8	3.20	111.51	108.61
3	G	307	MTE	N3-C2-N1	-3.16	117.54	123.32
3	I	309	MTE	O3P-P-O2P	3.12	119.52	107.80
3	G	307	MTE	O3P-P-O2P	3.06	119.29	107.80
3	B	302	MTE	O4-C4-N3	3.05	125.84	120.11
3	C	303	MTE	C2-N3-C4	-3.03	119.61	125.11
3	E	305	MTE	O3P-P-O2P	3.01	119.09	107.80
3	A	301	MTE	C2-N3-C4	-2.99	119.68	125.11
3	F	306	MTE	N3-C2-N1	-2.98	117.86	123.32
3	D	304	MTE	O4-C4-N3	2.95	125.66	120.11
3	F	306	MTE	C2-N3-C4	-2.94	119.77	125.11
3	H	308	MTE	O4-C4-N3	2.93	125.62	120.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	303	MTE	O4-C4-N3	2.93	125.61	120.11
3	F	306	MTE	O4-C4-N3	2.91	125.59	120.11
3	D	304	MTE	N3-C2-N1	-2.90	118.02	123.32
3	J	310	MTE	O4'-P-O1P	2.90	114.27	106.44
3	B	302	MTE	C2-N3-C4	-2.89	119.87	125.11
3	H	308	MTE	N3-C2-N1	-2.88	118.05	123.32
3	J	310	MTE	N3-C2-N1	-2.87	118.06	123.32
3	F	306	MTE	O3P-P-O2P	2.86	118.53	107.80
3	A	301	MTE	N3-C2-N1	-2.84	118.13	123.32
3	H	308	MTE	O3P-P-O2P	2.83	118.42	107.80
3	E	305	MTE	O3'-C7-N8	2.83	111.17	108.61
3	C	303	MTE	N3-C2-N1	-2.81	118.17	123.32
3	B	302	MTE	N3-C2-N1	-2.78	118.23	123.32
3	D	304	MTE	O4'-P-O1P	2.78	113.95	106.44
3	E	305	MTE	C2-N3-C4	-2.77	120.09	125.11
3	G	307	MTE	O4-C4-N3	2.75	125.29	120.11
3	D	304	MTE	O3P-P-O2P	2.74	118.07	107.80
3	J	310	MTE	C2-N3-C4	-2.72	120.17	125.11
3	J	310	MTE	O4-C4-N3	2.70	125.19	120.11
3	E	305	MTE	O4-C4-N3	2.70	125.19	120.11
3	E	305	MTE	N3-C2-N1	-2.67	118.43	123.32
3	D	304	MTE	C2-N3-C4	-2.61	120.37	125.11
3	A	301	MTE	O4'-P-O1P	2.60	113.47	106.44
3	H	308	MTE	C2-N3-C4	-2.58	120.43	125.11
3	H	308	MTE	O3P-P-O4'	-2.54	100.04	106.67
3	I	309	MTE	C2-N3-C4	-2.52	120.54	125.11
3	F	306	MTE	O3P-P-O4'	-2.49	100.19	106.67
3	I	309	MTE	O4-C4-N3	2.47	124.75	120.11
3	A	301	MTE	O4-C4-N3	2.42	124.67	120.11
3	C	303	MTE	O3P-P-O2P	2.34	116.58	107.80
3	H	308	MTE	O2P-P-O1P	-2.25	102.07	110.83
3	D	304	MTE	N2-C2-N3	2.25	121.50	116.76
3	A	301	MTE	O3P-P-O2P	2.24	116.19	107.80
3	J	310	MTE	O3P-P-O2P	2.23	116.15	107.80
3	F	306	MTE	N2-C2-N3	2.20	121.40	116.76
3	E	305	MTE	N2-C2-N3	2.17	121.35	116.76
3	F	306	MTE	O2P-P-O1P	-2.15	102.46	110.83
3	B	302	MTE	O4'-C4'-C3'	2.13	113.84	108.58
3	H	308	MTE	O4'-C4'-C3'	2.12	113.81	108.58
3	B	302	MTE	O3P-P-O4'	-2.11	101.16	106.67
3	D	304	MTE	O3'-C7-N8	2.10	110.51	108.61
3	C	303	MTE	N2-C2-N3	2.10	121.19	116.76

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	302	MTE	N2-C2-N3	2.08	121.16	116.76
3	D	304	MTE	O4'-C4'-C3'	2.08	113.72	108.58
3	G	307	MTE	C2-N3-C4	-2.08	121.34	125.11
3	H	308	MTE	N2-C2-N3	2.06	121.11	116.76
3	E	305	MTE	O4'-C4'-C3'	2.00	113.53	108.58

All (10) chirality outliers are listed below:

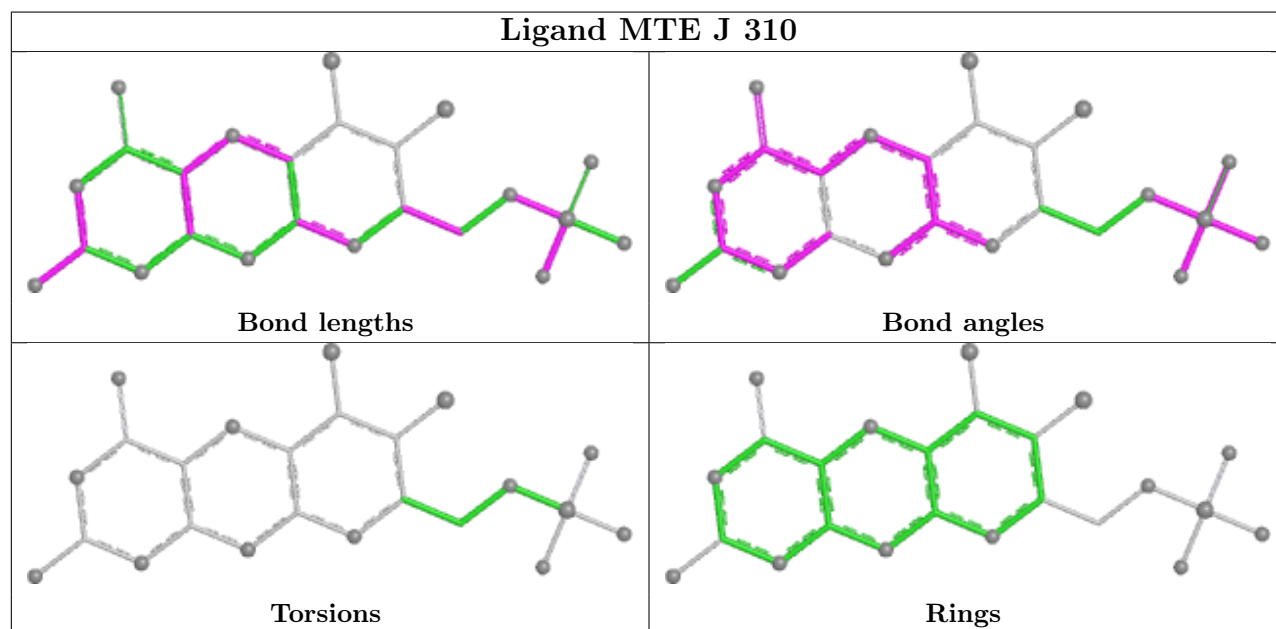
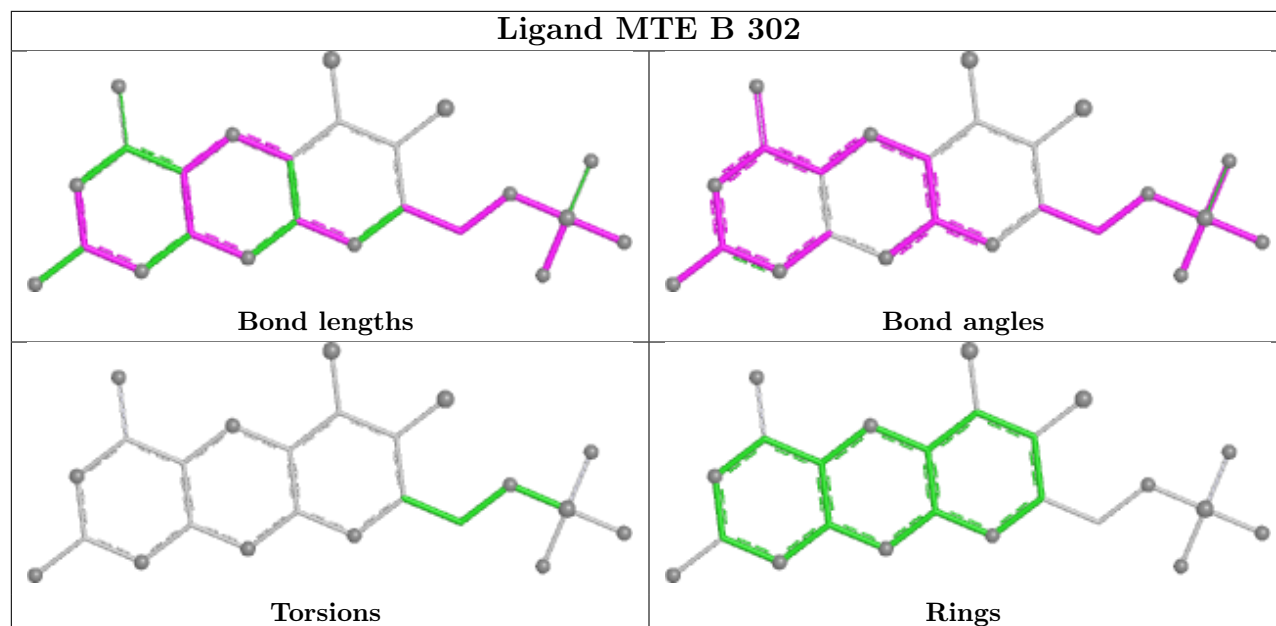
Mol	Chain	Res	Type	Atom
3	A	301	MTE	C6
3	B	302	MTE	C6
3	C	303	MTE	C6
3	D	304	MTE	C6
3	E	305	MTE	C6
3	F	306	MTE	C6
3	G	307	MTE	C6
3	H	308	MTE	C6
3	I	309	MTE	C6
3	J	310	MTE	C6

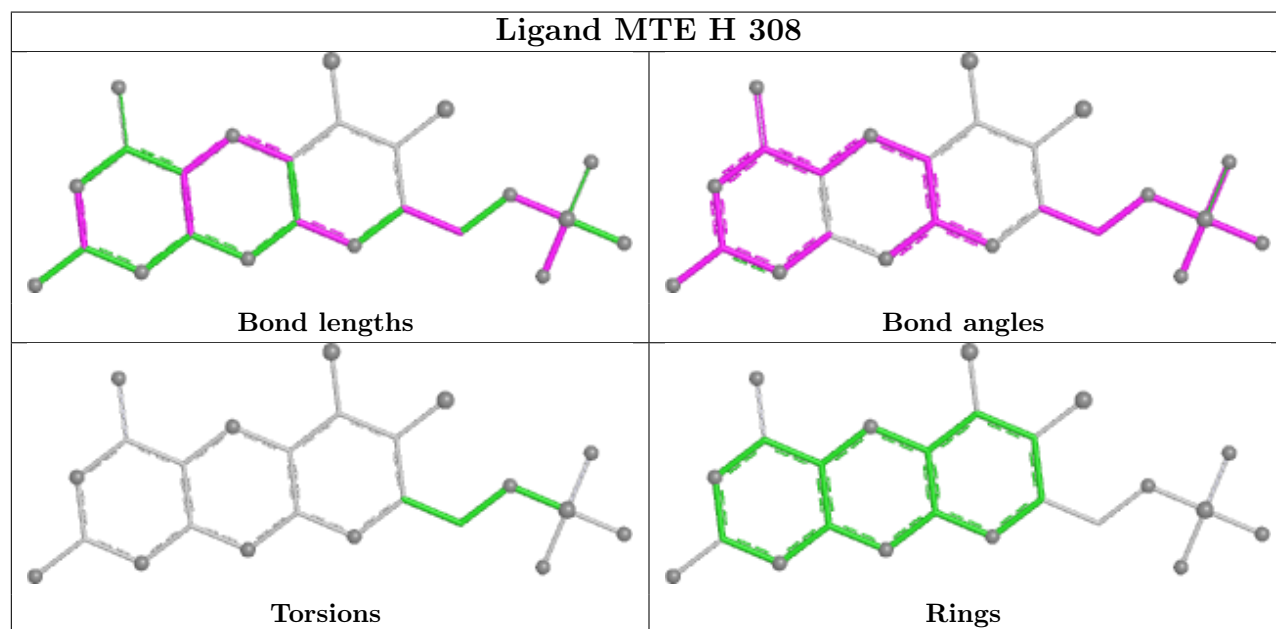
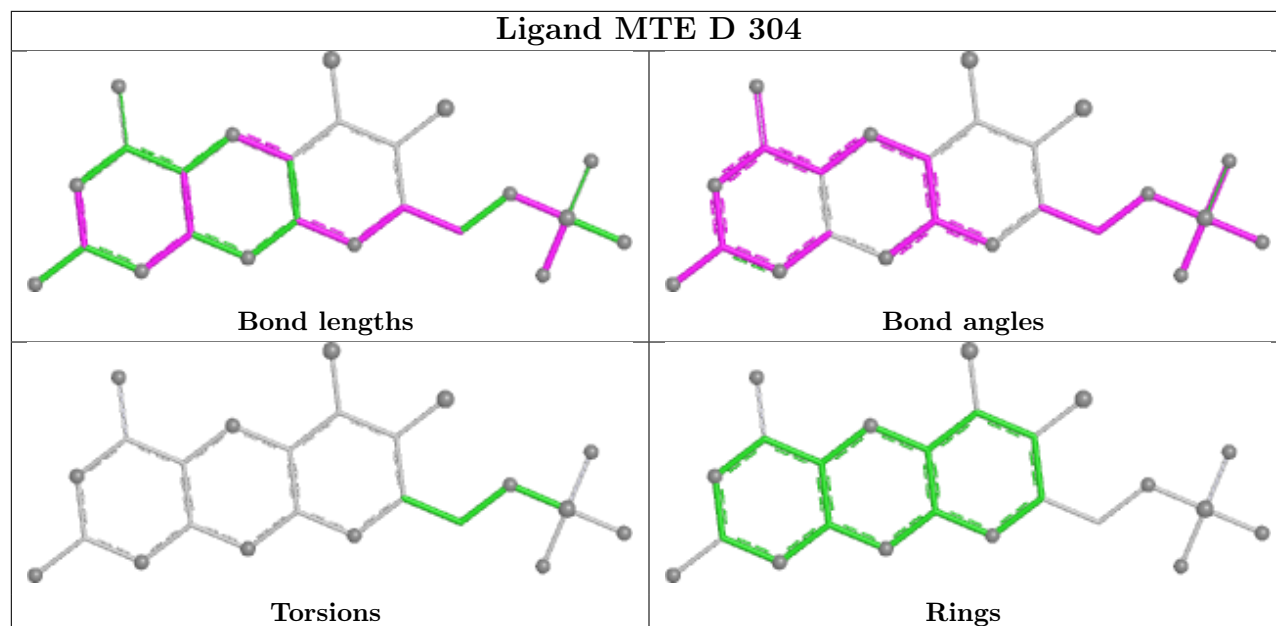
There are no torsion outliers.

There are no ring outliers.

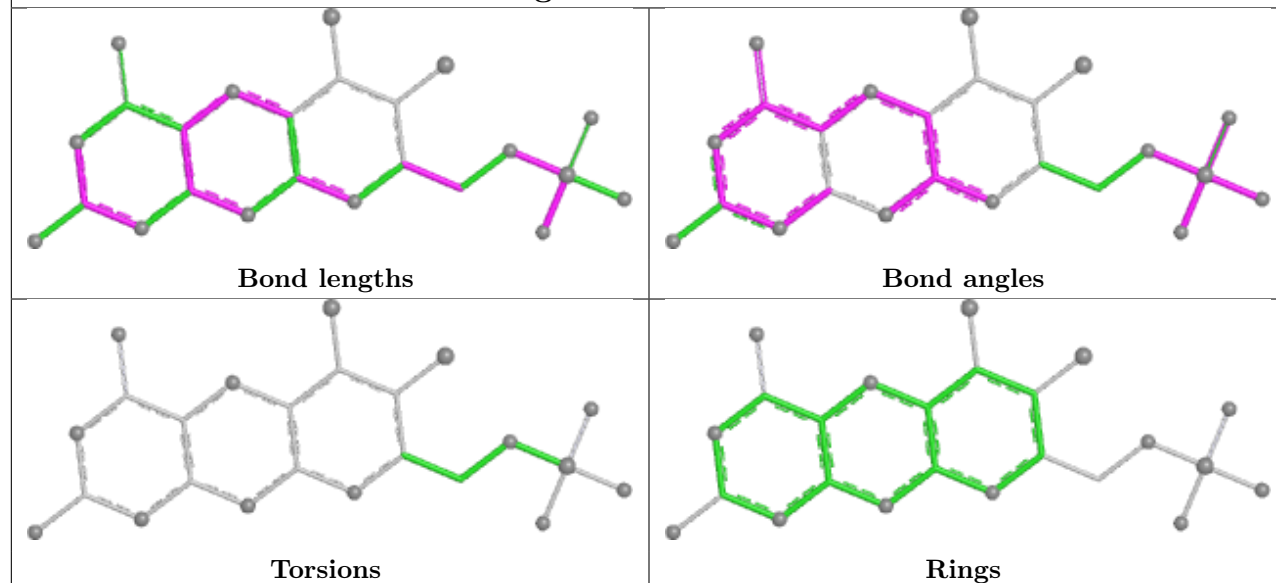
No monomer is involved in short contacts.

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.

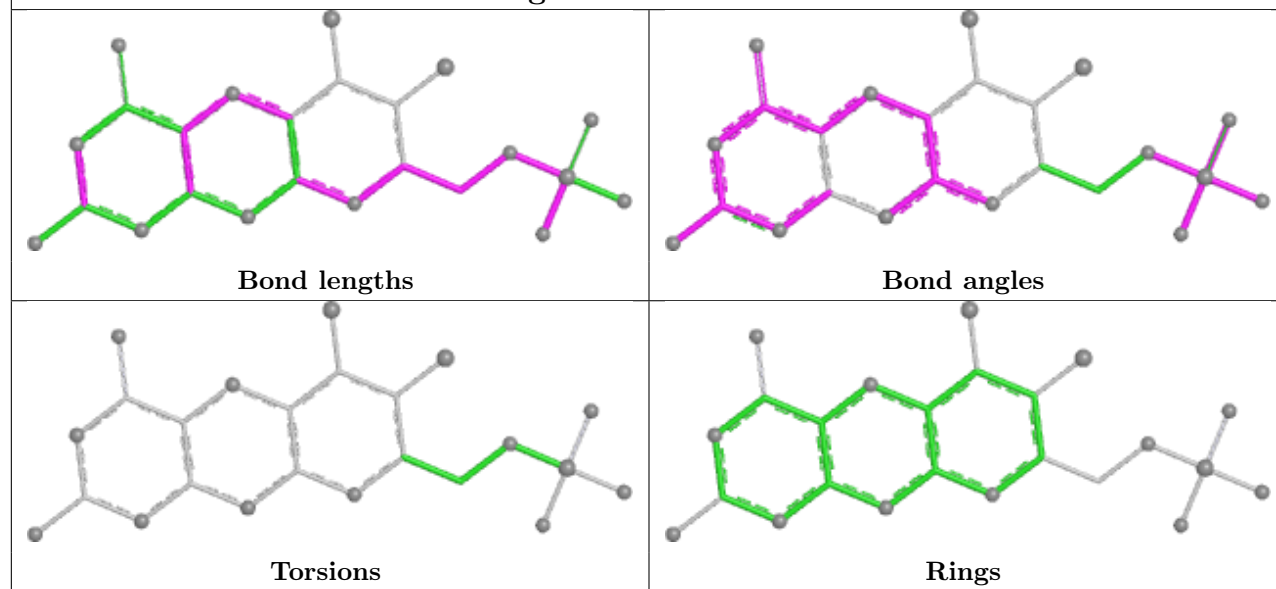




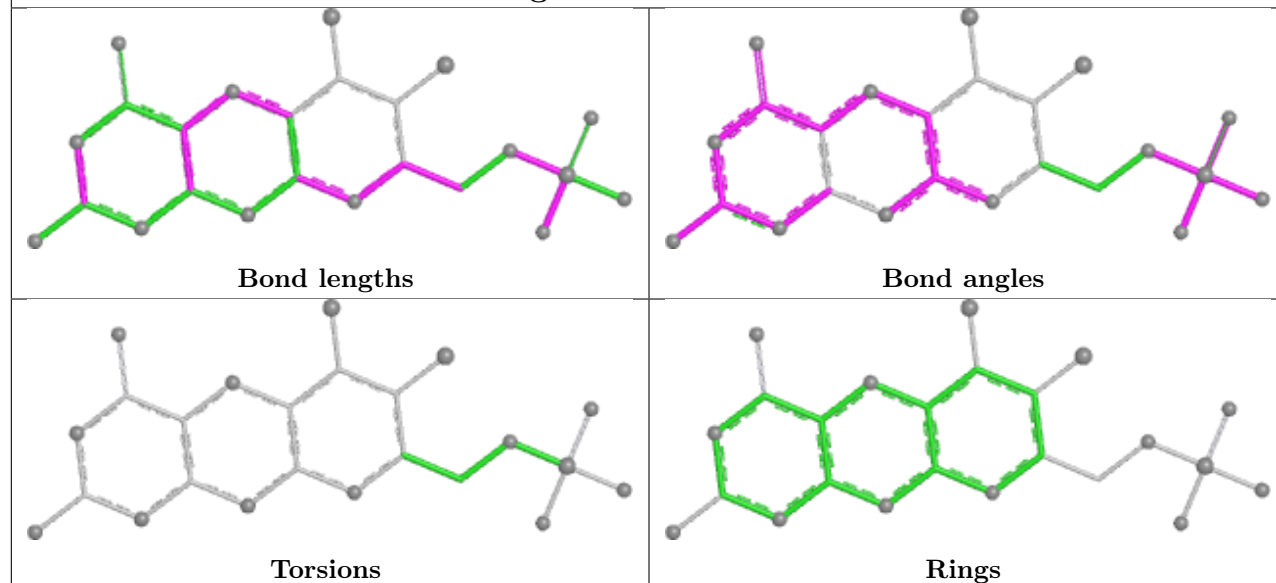
Ligand MTE I 309



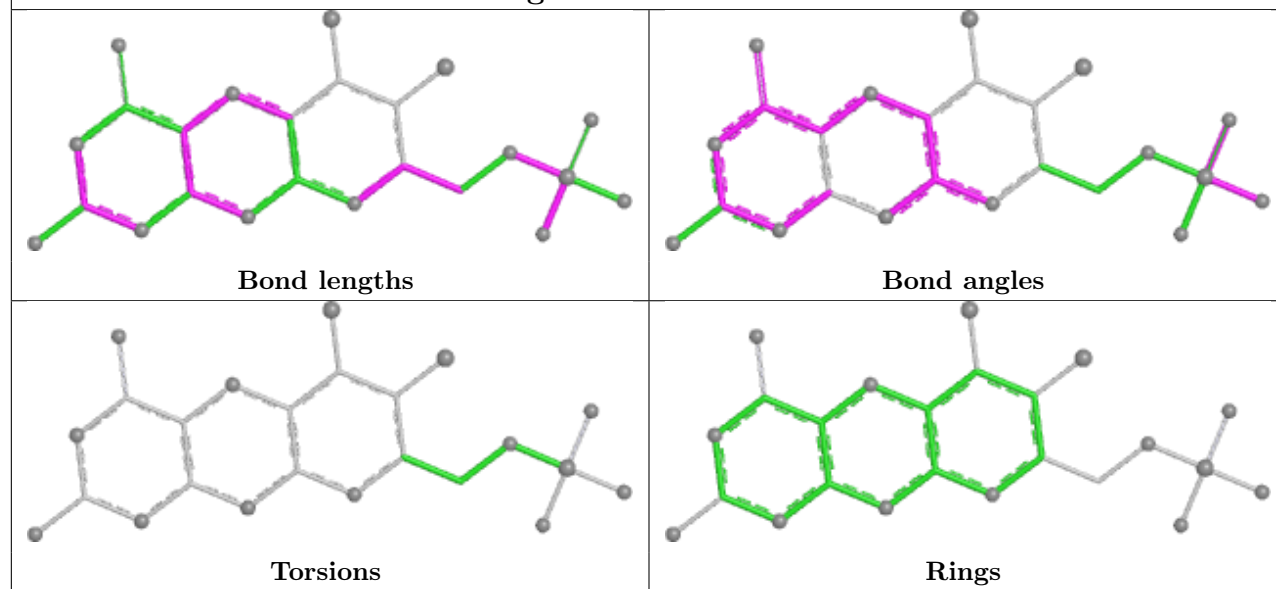
Ligand MTE C 303

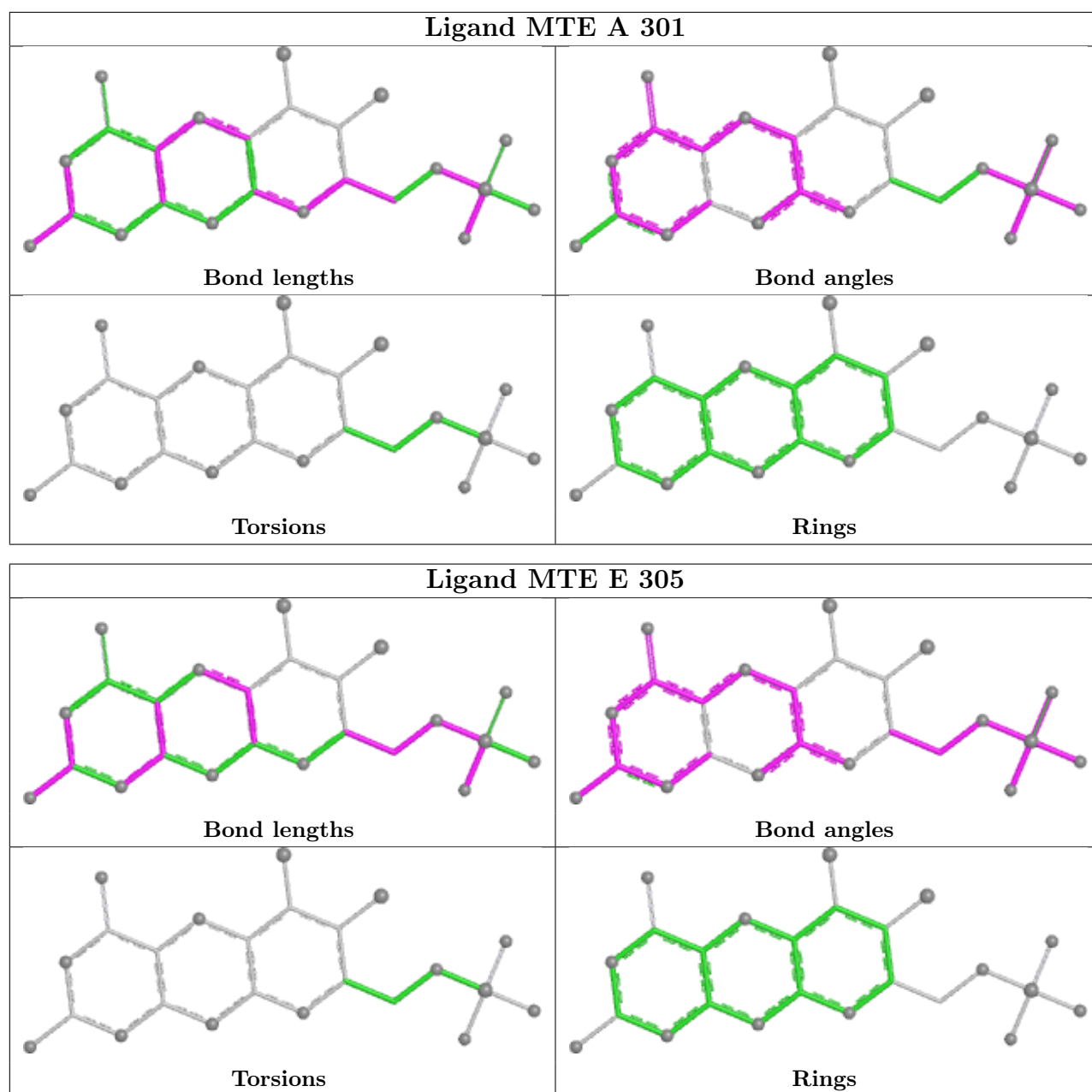


Ligand MTE F 306



Ligand MTE G 307





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	263/298 (88%)	0.32	8 (3%) 52 49	22, 46, 78, 95	0
1	B	262/298 (87%)	0.94	26 (9%) 13 10	33, 68, 94, 106	0
1	C	264/298 (88%)	0.34	9 (3%) 48 45	22, 49, 78, 90	0
1	D	263/298 (88%)	0.33	3 (1%) 78 76	24, 46, 76, 92	0
1	E	262/298 (87%)	0.34	10 (3%) 44 41	25, 46, 80, 97	0
1	F	262/298 (87%)	1.07	28 (10%) 11 8	34, 64, 88, 94	1 (0%)
1	G	261/298 (87%)	0.58	19 (7%) 21 18	21, 49, 84, 100	0
1	H	263/298 (88%)	1.06	28 (10%) 11 8	36, 68, 95, 102	0
1	I	262/298 (87%)	1.03	32 (12%) 8 6	34, 61, 90, 106	0
1	J	262/298 (87%)	0.30	17 (6%) 25 22	20, 44, 73, 91	0
All	All	2624/2980 (88%)	0.63	180 (6%) 23 20	20, 54, 88, 106	1 (0%)

All (180) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	J	151	PHE	5.6
1	C	257	GLY	4.9
1	A	286	LEU	4.8
1	E	257	GLY	4.8
1	I	151	PHE	4.8
1	B	257	GLY	4.5
1	F	257	GLY	4.4
1	I	257	GLY	4.3
1	H	257	GLY	4.2
1	F	19	ALA	4.2
1	D	151	PHE	4.1
1	F	102	CYS	4.0
1	B	20	LEU	4.0

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Mol	Chain	Res	Type	RSRZ
1	I	19	ALA	3.8
1	G	31	ASN	3.8
1	H	20	LEU	3.7
1	A	151	PHE	3.6
1	H	286	LEU	3.6
1	H	151	PHE	3.5
1	A	257	GLY	3.5
1	E	280	LEU	3.4
1	E	285	ASP	3.4
1	G	20	LEU	3.4
1	D	257	GLY	3.4
1	I	55	ASP	3.4
1	I	152	ILE	3.3
1	C	151	PHE	3.3
1	F	277	VAL	3.3
1	J	229	ASP	3.3
1	F	286	LEU	3.2
1	B	151	PHE	3.2
1	F	151	PHE	3.2
1	J	257	GLY	3.2
1	G	257	GLY	3.1
1	G	143	GLN	3.1
1	I	263	GLN	3.1
1	F	133	VAL	3.1
1	I	193	VAL	3.1
1	B	189	ASN	3.0
1	F	123	LEU	3.0
1	H	19	ALA	3.0
1	H	113	ILE	3.0
1	G	151	PHE	2.9
1	A	250	THR	2.9
1	B	184	ALA	2.9
1	I	20	LEU	2.9
1	F	166	LEU	2.9
1	G	67	PRO	2.9
1	G	32	LEU	2.9
1	B	71	LYS	2.8
1	G	87	LEU	2.8
1	I	285	ASP	2.8
1	G	28	TRP	2.8
1	F	77	ALA	2.8
1	I	150	ARG	2.8

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Mol	Chain	Res	Type	RSRZ
1	F	232	GLY	2.8
1	A	285	ASP	2.7
1	H	226	ALA	2.7
1	E	151	PHE	2.7
1	H	152	ILE	2.7
1	J	285	ASP	2.7
1	E	28	TRP	2.7
1	I	223	TRP	2.7
1	F	89	ARG	2.7
1	A	284	LEU	2.7
1	E	123	LEU	2.7
1	F	241	VAL	2.6
1	B	127	THR	2.6
1	F	51	LEU	2.6
1	G	286	LEU	2.6
1	I	225	LEU	2.6
1	B	77	ALA	2.6
1	I	54	ALA	2.6
1	J	51	LEU	2.6
1	I	21	GLU	2.6
1	H	143	GLN	2.6
1	H	81	THR	2.6
1	J	88	THR	2.6
1	B	178	VAL	2.6
1	B	26	ALA	2.5
1	I	191	ALA	2.5
1	C	263	GLN	2.5
1	H	252	ARG	2.5
1	F	120	LEU	2.5
1	I	176	MET	2.5
1	H	34	LEU	2.5
1	B	45	ASN	2.5
1	C	31	ASN	2.5
1	F	274	ALA	2.5
1	I	66	ASP	2.5
1	I	102	CYS	2.5
1	F	65	THR	2.5
1	J	250	THR	2.5
1	D	141	PRO	2.5
1	I	67	PRO	2.5
1	F	93	LEU	2.4
1	G	30	ASN	2.4

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Mol	Chain	Res	Type	RSRZ
1	F	130	ALA	2.4
1	H	153	GLY	2.4
1	I	89	ARG	2.4
1	J	263	GLN	2.4
1	B	67	PRO	2.4
1	F	174	THR	2.4
1	F	126	PRO	2.4
1	H	277	VAL	2.4
1	J	225	LEU	2.4
1	G	92	PRO	2.3
1	B	22	PHE	2.3
1	E	277	VAL	2.3
1	G	93	LEU	2.3
1	B	61	GLY	2.3
1	F	249	ALA	2.3
1	B	69	THR	2.3
1	J	152	ILE	2.3
1	I	70	LEU	2.3
1	I	58	ALA	2.3
1	G	88	THR	2.3
1	B	150	ARG	2.3
1	E	279	SER	2.3
1	H	91	PHE	2.3
1	B	286	LEU	2.3
1	G	80	LEU	2.3
1	H	214	LEU	2.3
1	B	98	TYR	2.3
1	F	124	ALA	2.3
1	B	87	LEU	2.2
1	E	284	LEU	2.2
1	F	80	LEU	2.2
1	I	187	PRO	2.2
1	I	42	GLY	2.2
1	G	27	ALA	2.2
1	J	284	LEU	2.2
1	B	193	VAL	2.2
1	H	99	ARG	2.2
1	C	230	GLU	2.2
1	H	88	THR	2.2
1	H	130	ALA	2.2
1	I	88	THR	2.2
1	B	152	ILE	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	51	LEU	2.2
1	C	150	ARG	2.2
1	F	150	ARG	2.2
1	J	241	VAL	2.2
1	G	64	LYS	2.2
1	A	26	ALA	2.2
1	H	32	LEU	2.1
1	H	75	GLU	2.1
1	E	102	CYS	2.1
1	B	181	TYR	2.1
1	B	269	LEU	2.1
1	H	80	LEU	2.1
1	I	138	ILE	2.1
1	I	175	LEU	2.1
1	H	238	ASN	2.1
1	I	44	ASN	2.1
1	H	27	ALA	2.1
1	B	185	LEU	2.1
1	C	175	LEU	2.1
1	A	31	ASN	2.1
1	J	275	ASP	2.1
1	H	180	VAL	2.1
1	I	210	VAL	2.1
1	B	255	GLY	2.1
1	J	155	GLY	2.1
1	H	140	ALA	2.1
1	G	175	LEU	2.1
1	J	277	VAL	2.1
1	F	250	THR	2.0
1	F	169	ALA	2.0
1	H	185	LEU	2.0
1	F	113	ILE	2.0
1	J	150	ARG	2.0
1	B	190	GLY	2.0
1	I	22	PHE	2.0
1	G	250	THR	2.0
1	I	256	SER	2.0
1	J	143	GLN	2.0
1	C	286	LEU	2.0
1	H	266	PRO	2.0
1	I	192	PRO	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

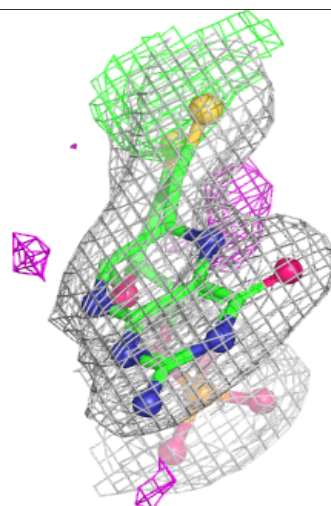
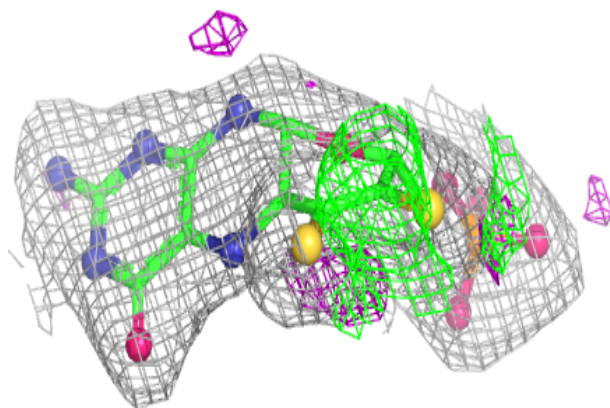
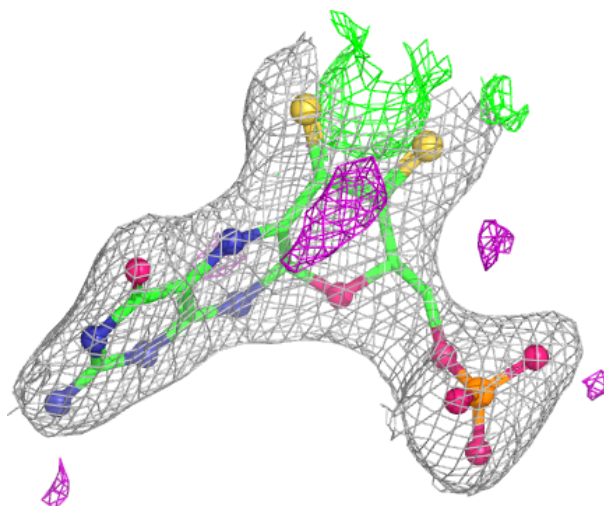
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	MTE	B	302	24/24	0.95	0.09	22,39,50,56	0
3	MTE	F	306	24/24	0.95	0.08	26,38,46,47	0
3	MTE	H	308	24/24	0.95	0.08	30,43,48,50	0
3	MTE	I	309	24/24	0.95	0.09	26,42,49,50	0
3	MTE	C	303	24/24	0.96	0.08	15,32,39,42	0
2	W	I	409	1/1	0.96	0.07	42,42,42,42	0
3	MTE	E	305	24/24	0.97	0.06	8,22,33,36	0
3	MTE	D	304	24/24	0.97	0.07	18,28,35,38	0
3	MTE	G	307	24/24	0.98	0.06	17,26,32,36	0
3	MTE	A	301	24/24	0.98	0.05	16,23,30,32	0
2	W	F	406	1/1	0.98	0.06	41,41,41,41	0
3	MTE	J	310	24/24	0.98	0.06	8,22,28,29	0
2	W	C	403	1/1	0.99	0.07	32,32,32,32	0
2	W	J	410	1/1	0.99	0.07	30,30,30,30	0
2	W	D	404	1/1	0.99	0.06	32,32,32,32	0
2	W	E	405	1/1	0.99	0.04	32,32,32,32	0
2	W	B	402	1/1	0.99	0.06	43,43,43,43	0
2	W	H	408	1/1	0.99	0.06	41,41,41,41	0
2	W	G	407	1/1	1.00	0.06	29,29,29,29	0
2	W	A	401	1/1	1.00	0.07	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.

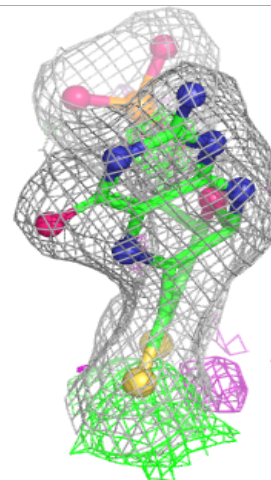
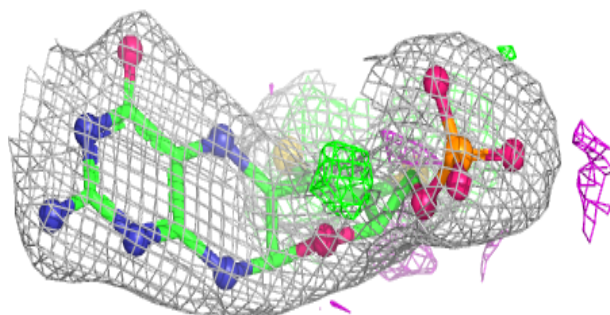
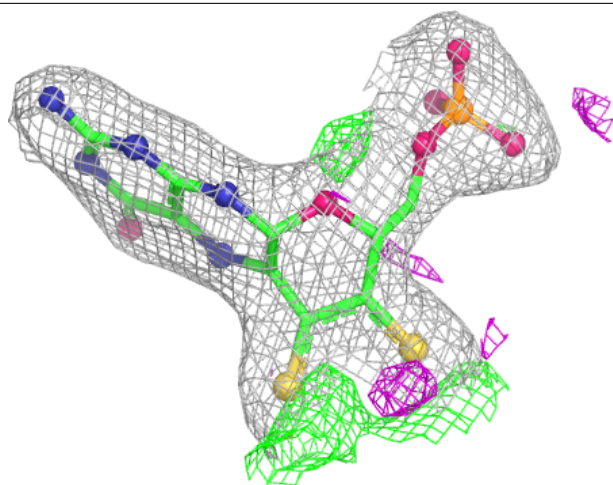
Electron density around MTE B 302:

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



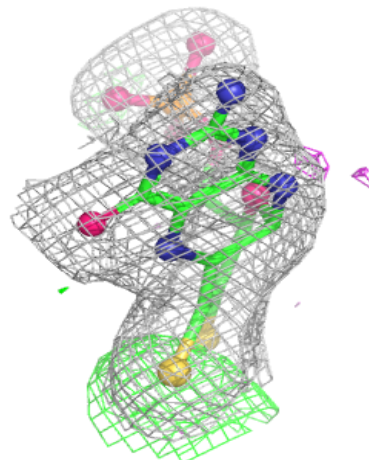
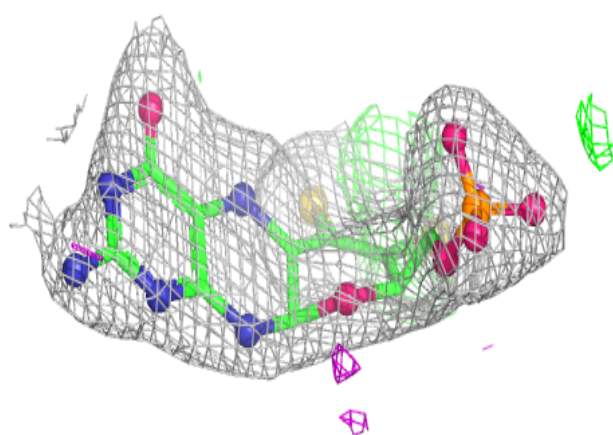
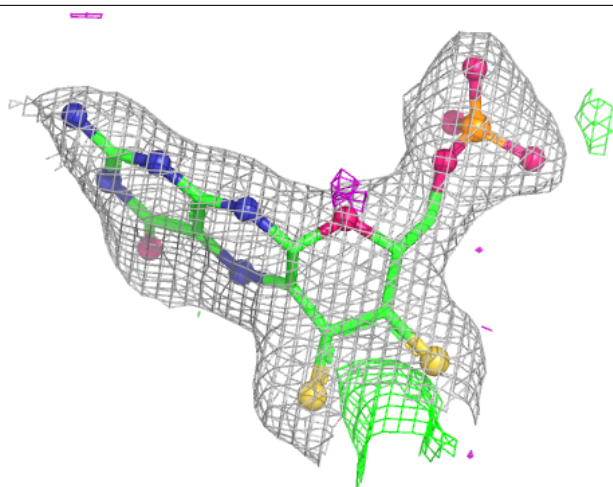
Electron density around MTE F 306:

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



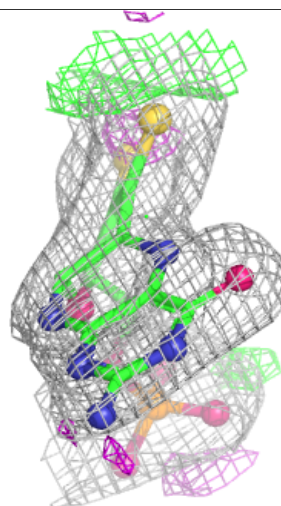
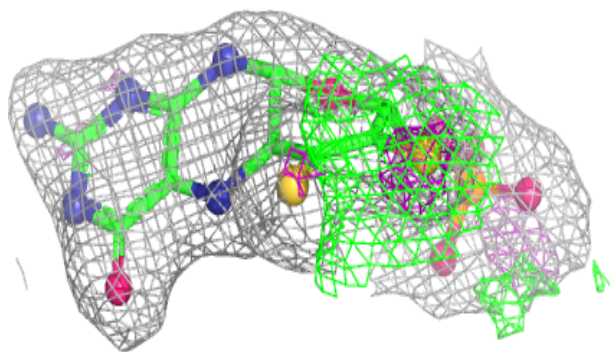
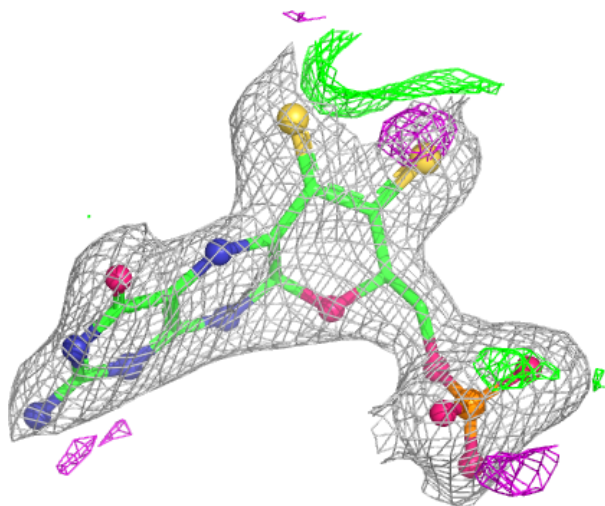
Electron density around MTE H 308:

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



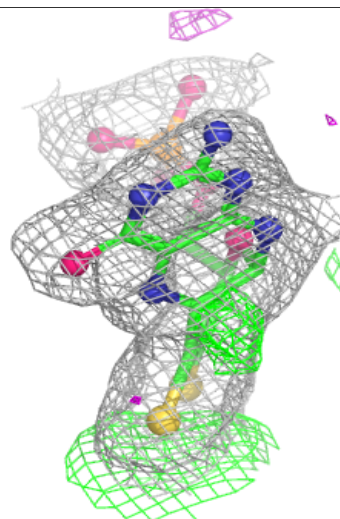
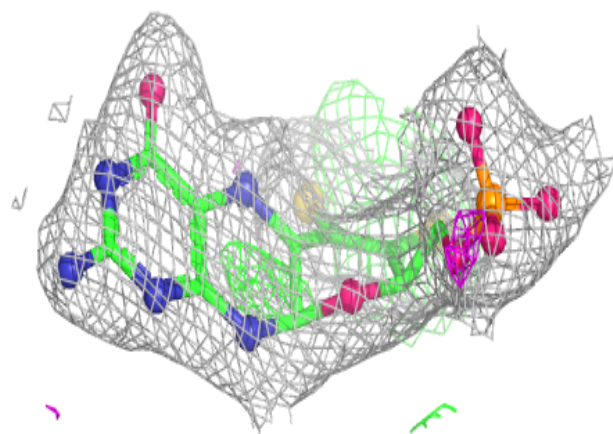
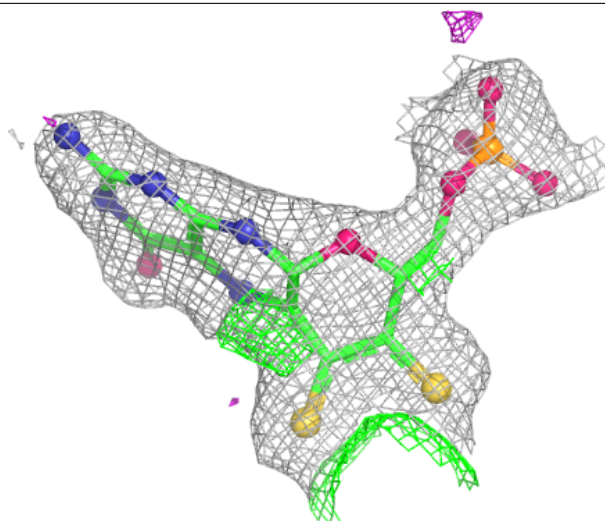
Electron density around MTE I 309:

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



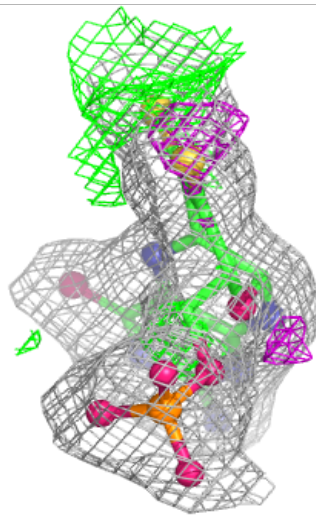
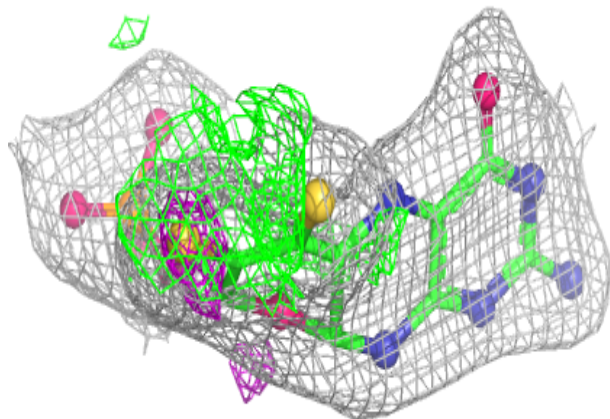
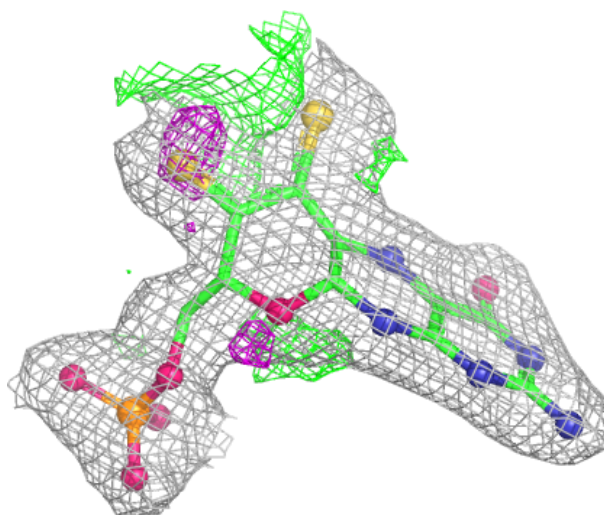
Electron density around MTE C 303:

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



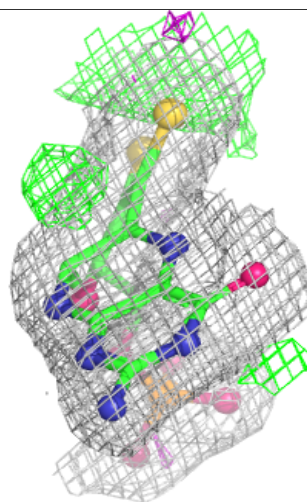
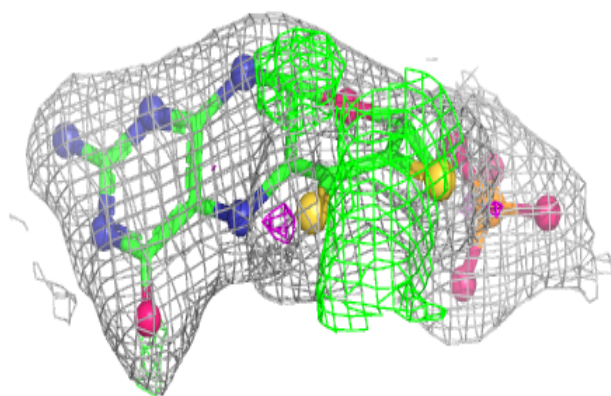
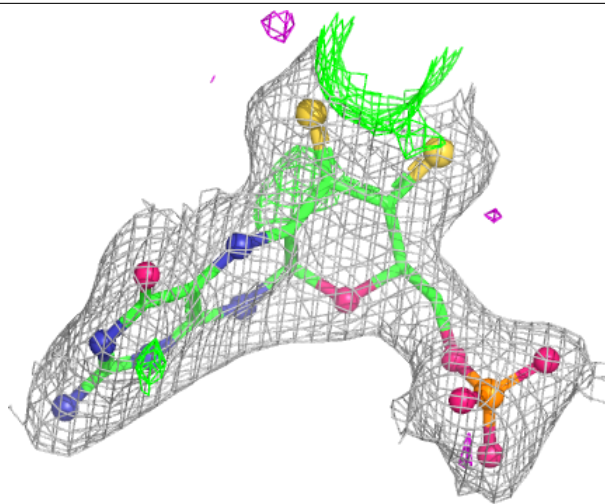
Electron density around MTE E 305:

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



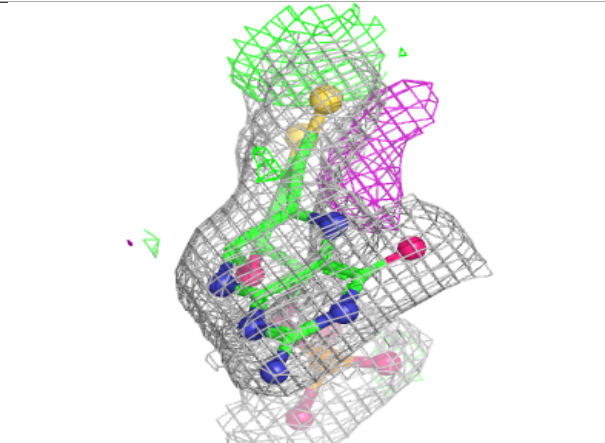
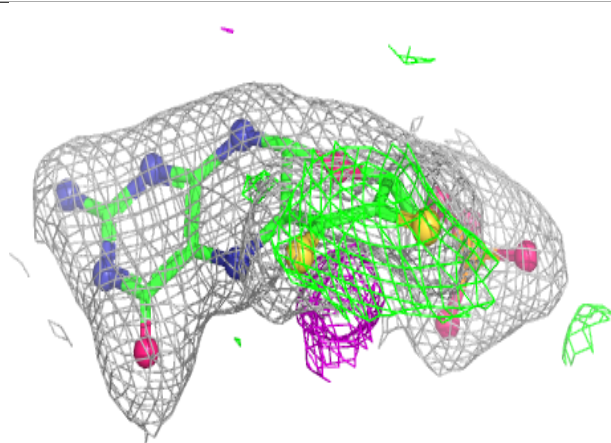
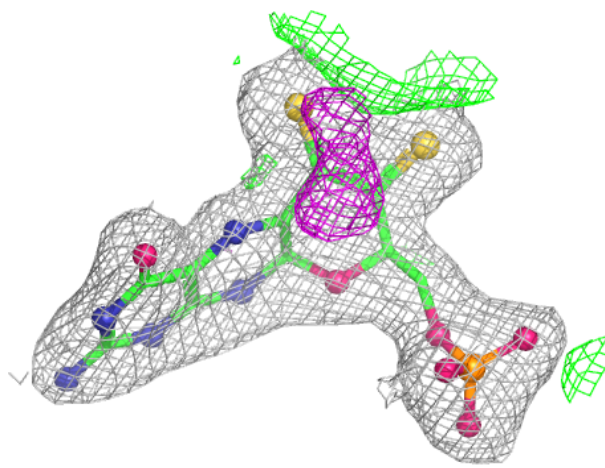
Electron density around MTE D 304:

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



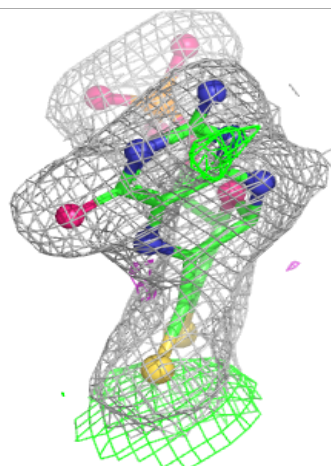
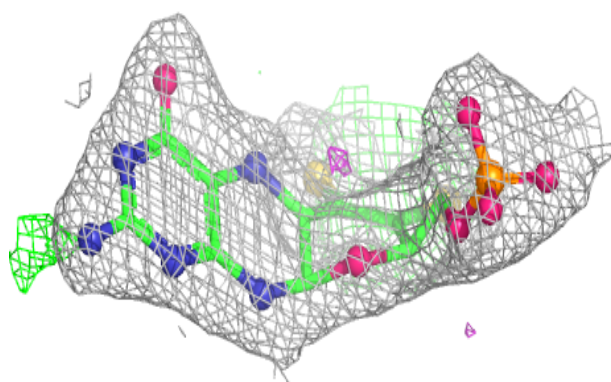
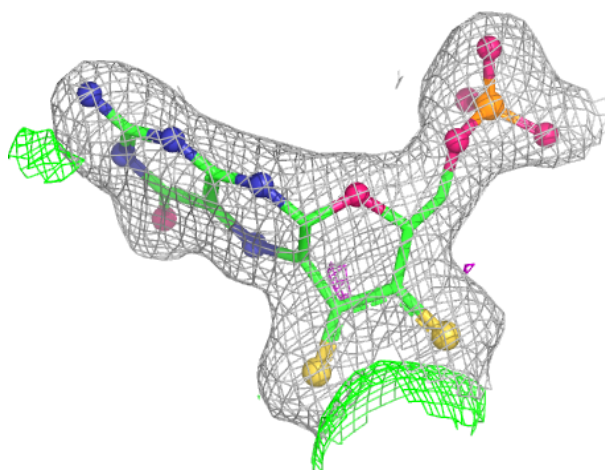
Electron density around MTE G 307:

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



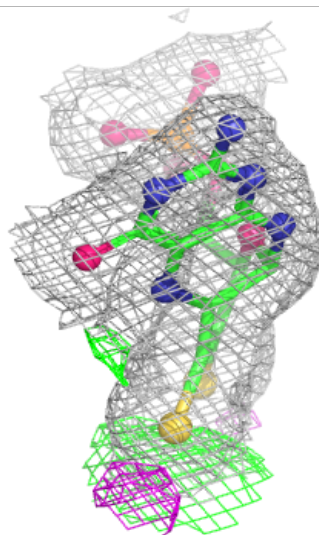
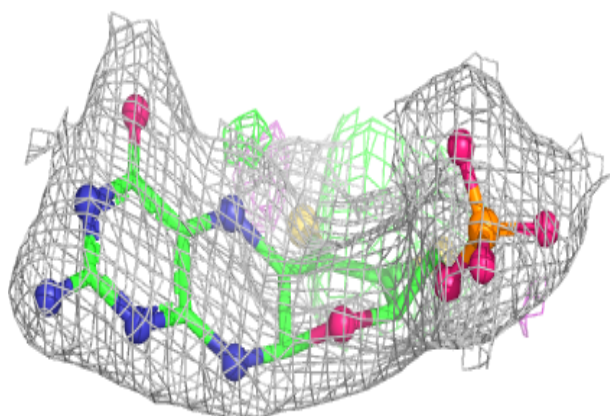
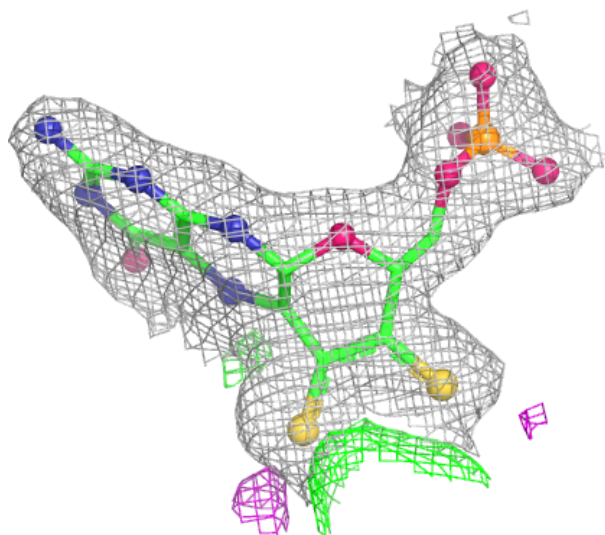
Electron density around MTE A 301:

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



Electron density around MTE J 310:

$2mF_o-DF_c$ (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.