



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 17, 2026 – 06:34 PM UTC

PDB ID : 3TT1 / pdb_00003tt1
Title : Crystal Structure of LeuT in the outward-open conformation in complex with Fab
Authors : Krishnamurthy, H.; Gouaux, E.
Deposited on : 2011-09-13
Resolution : 3.10 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

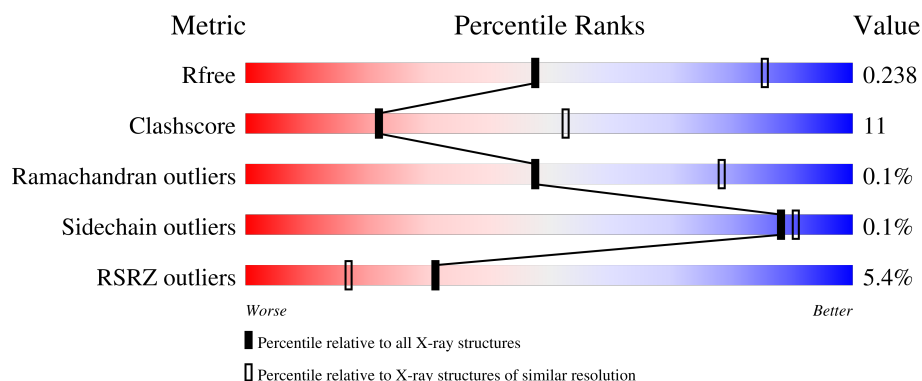
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.10 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	1456 (3.10-3.10)
Clashscore	190562	1539 (3.10-3.10)
Ramachandran outliers	187476	1467 (3.10-3.10)
Sidechain outliers	187428	1467 (3.10-3.10)
RSRZ outliers	180081	1456 (3.10-3.10)

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	519	7% 75% 20% .
1	B	519	6% 76% 20% .
2	L	218	% 82% 17% .
2	M	218	2% 82% 17% .
3	H	219	5% 71% 26% .

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Mol	Chain	Length	Quality of chain
3	I	219	

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
5	SOG	A	520	-	-	X	-
5	SOG	B	520	-	-	X	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 14274 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 Leucine transporter LeuT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	498	Total	C	N	O	S	0	0	0
			3835	2605	595	623	12			
1	B	498	Total	C	N	O	S	0	0	0
			3835	2605	595	623	12			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	108	PHE	TYR	engineered mutation	UNP O67854
A	288	ALA	LYS	engineered mutation	UNP O67854
A	514	GLY	-	expression tag	UNP O67854
A	515	THR	-	expression tag	UNP O67854
A	516	LEU	-	expression tag	UNP O67854
A	517	VAL	-	expression tag	UNP O67854
A	518	PRO	-	expression tag	UNP O67854
A	519	ARG	-	expression tag	UNP O67854
B	108	PHE	TYR	engineered mutation	UNP O67854
B	288	ALA	LYS	engineered mutation	UNP O67854
B	514	GLY	-	expression tag	UNP O67854
B	515	THR	-	expression tag	UNP O67854
B	516	LEU	-	expression tag	UNP O67854
B	517	VAL	-	expression tag	UNP O67854
B	518	PRO	-	expression tag	UNP O67854
B	519	ARG	-	expression tag	UNP O67854

- Molecule 2 is a protein called mouse monoclonal 1gG2a Fab fragment, heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	L	215	Total	C	N	O	S	0	0	0
			1656	1029	276	345	6			
2	M	215	Total	C	N	O	S	0	0	0
			1656	1029	276	345	6			

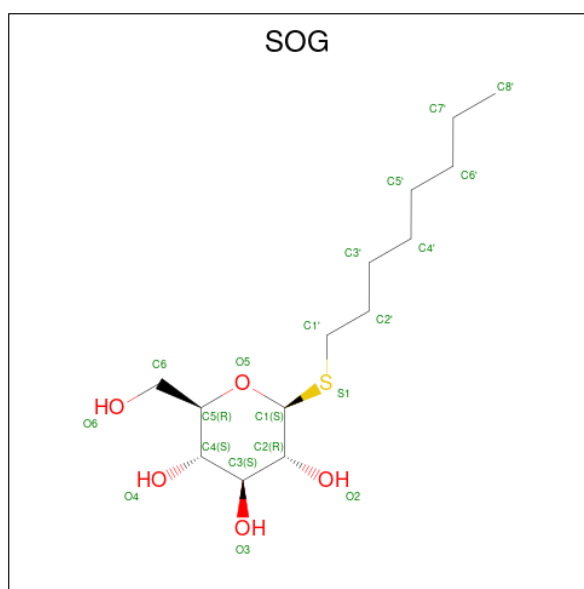
- Molecule 3 is a protein called mouse monoclonal 1gG2a Fab fragment, kappa light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	H	212	Total	C	N	O	S	0	0	0
			1598	1012	260	318	8			
3	I	212	Total	C	N	O	S	0	0	0
			1598	1012	260	318	8			

- Molecule 4 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total	Na	0	0
			2	2		
4	B	2	Total	Na	0	0
			2	2		

- Molecule 5 is octyl 1-thio-beta-D-glucopyranoside (CCD ID: SOG) (formula: C₁₄H₂₈O₅S).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	O	S	0	0
			20	14	5	1		
5	A	1	Total	C			0	0
			8	8				
5	A	1	Total	C			0	0
			8	8				
5	B	1	Total	C	O	S	0	0
			20	14	5	1		
5	B	1	Total	C			0	0
			8	8				

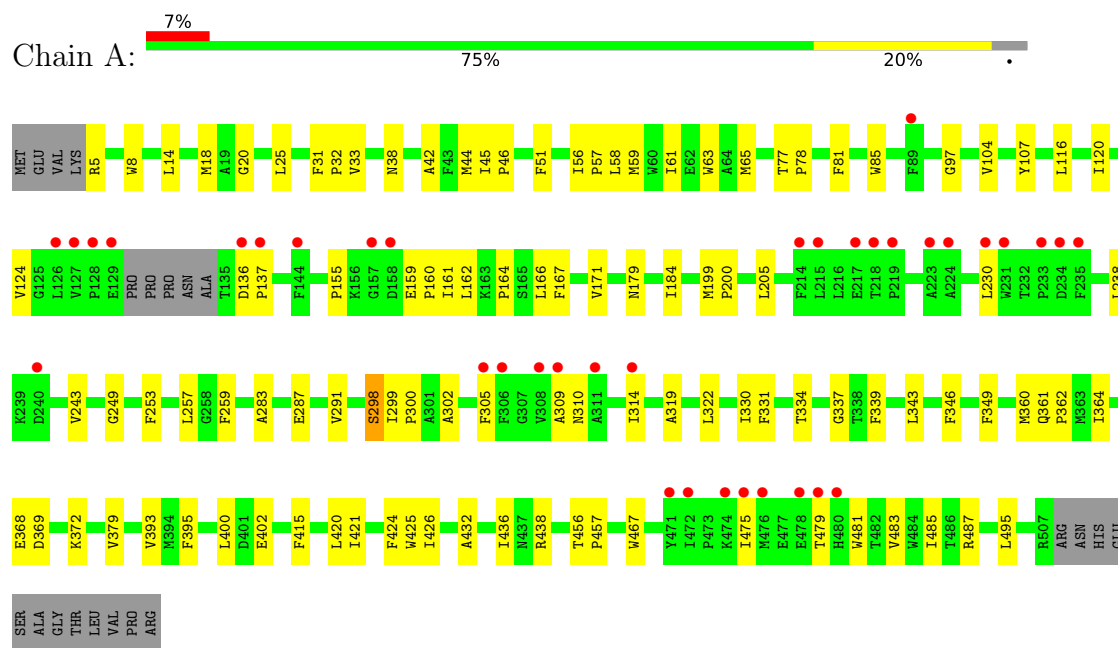
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	17	Total 17	O 17	0	0
6	B	10	Total 10	O 10	0	0
6	H	1	Total 1	O 1	0	0

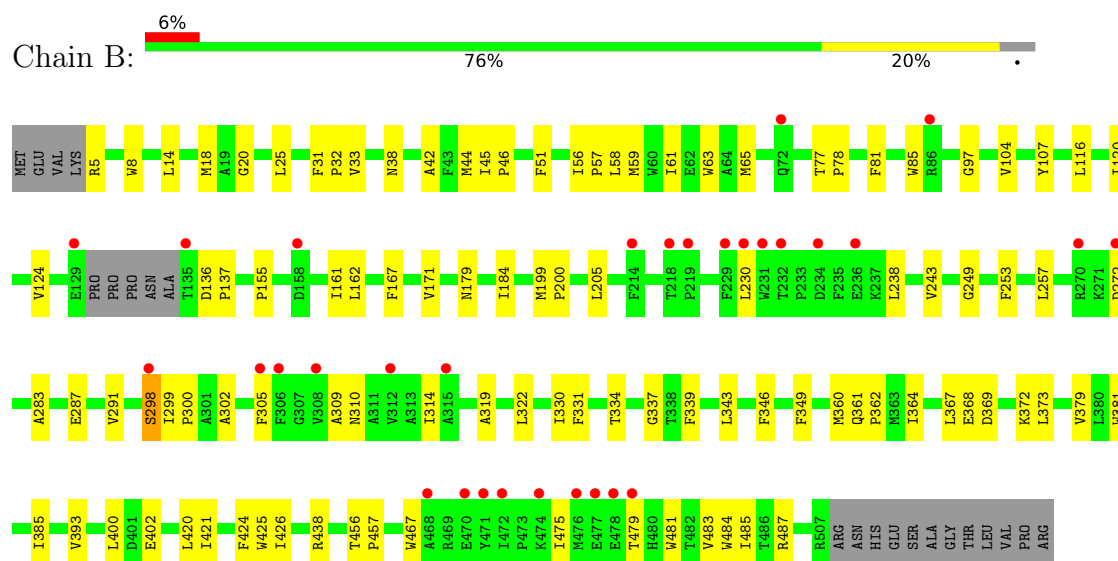
3 Residue-property plots [i](#)

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: Leucine transporter LeuT



• Molecule 1: Leucine transporter LeuT



- Chain I:
-
- | Residue | Count |
|---------|-------|
| Q1 | 1 |
| Q6 | 1 |
| L11 | 1 |
| Y27 | 1 |
| Y32 | 1 |
| Y33 | 1 |
| Y34 | 1 |
| N35 | 1 |
| K36 | 1 |
| M37 | 1 |
| G42 | 1 |
| W47 | 1 |
| E50 | 1 |
| F51 | 1 |
| Y52 | 1 |
| P53 | 1 |
| F57 | 1 |
| S78 | 1 |
| T87 | 1 |
| S88 | 1 |
| Y94 | 1 |
| F95 | 1 |
| C96 | 1 |
| A97 | 1 |
| R98 | 1 |
| A99 | 1 |
| V100 | 1 |
| V101 | 1 |
| M102 | 1 |
| M103 | 1 |
| G104 | 1 |
| R105 | 1 |
| M108 | 1 |
| G112 | 1 |
| Q113 | 1 |
| T118 | 1 |
| Y130 | 1 |
| P131 | 1 |
| V135 | 1 |
| CYS | 1 |
| GLY | 1 |
| ASP | 1 |
| THR | 1 |
| TRP | 1 |

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	102.86Å 162.76Å 210.08Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.17 – 3.10 40.17 – 3.10	Depositor EDS
% Data completeness (in resolution range)	89.1 (40.17-3.10) 89.0 (40.17-3.10)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.10 (at 3.12Å)	Xtrriage
Refinement program	PHENIX (phenix.refine: 1.7_650)	Depositor
R, R_{free}	0.219 , 0.240 0.219 , 0.238	Depositor DCC
R_{free} test set	2949 reflections (5.11%)	wwPDB-VP
Wilson B-factor (Å ²)	75.0	Xtrriage
Anisotropy	0.160	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 54.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	14274	wwPDB-VP
Average B, all atoms (Å ²)	82.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.17% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: SOG, NA

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.30	0/3956	0.69	0/5403
1	B	0.29	0/3956	0.68	0/5403
2	L	0.26	0/1694	0.69	0/2303
2	M	0.27	0/1694	0.69	0/2303
3	H	0.26	0/1636	0.67	0/2230
3	I	0.27	0/1636	0.68	0/2230
All	All	0.28	0/14572	0.68	0/19872

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	3835	0	3813	79	0
1	B	3835	0	3813	75	0
2	L	1656	0	1567	25	0
2	M	1656	0	1567	25	0
3	H	1598	0	1550	49	0
3	I	1598	0	1550	50	0
4	A	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	2	0	0	0	0
5	A	36	0	58	12	0
5	B	28	0	42	25	0
6	A	17	0	0	2	0
6	B	10	0	0	0	0
6	H	1	0	0	0	0
All	All	14274	0	13960	302	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 11.

All (302) 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)
5:B:520:SOG:C3	5:B:520:SOG:C2	1.81	1.56
5:A:520:SOG:H2'2	5:B:520:SOG:H3'1	1.33	1.08
3:I:218:ILE:HG22	3:I:219:GLU:H	1.34	0.91
3:H:218:ILE:HG22	3:H:219:GLU:H	1.34	0.90
5:A:520:SOG:H2'2	5:B:520:SOG:C3'	2.03	0.87
1:A:395:PHE:HZ	5:A:520:SOG:H1'1	1.44	0.83
5:B:520:SOG:C2	5:B:520:SOG:C4	2.52	0.83
5:B:520:SOG:C3	5:B:520:SOG:O2	2.30	0.80
1:A:395:PHE:CZ	5:A:520:SOG:H1'1	2.16	0.80
3:H:6:GLN:HE21	3:H:112:GLY:HA3	1.47	0.79
1:A:253:PHE:HE2	1:A:467:TRP:HH2	1.33	0.77
3:I:6:GLN:HE21	3:I:112:GLY:HA3	1.50	0.77
1:B:253:PHE:HE2	1:B:467:TRP:HH2	1.33	0.76
5:A:520:SOG:C2'	5:B:520:SOG:C3'	2.64	0.75
3:H:37:MET:HE3	3:H:108:MET:SD	2.27	0.74
3:I:37:MET:HE3	3:I:108:MET:SD	2.26	0.74
3:H:131:PRO:HA	3:H:217:LYS:HE3	1.70	0.74
3:I:131:PRO:HA	3:I:217:LYS:HE3	1.70	0.72
1:B:484:TRP:CE3	5:B:520:SOG:H5'1	2.25	0.72
1:B:302:ALA:HB1	1:B:314:ILE:HD13	1.72	0.71
1:B:479:THR:HB	1:B:483:VAL:HG21	1.72	0.71
1:A:479:THR:HB	1:A:483:VAL:HG21	1.72	0.71
5:B:520:SOG:O2	5:B:520:SOG:O4	2.01	0.70
3:H:11:LEU:HD22	3:H:155:PRO:HG3	1.72	0.70
5:A:520:SOG:O6	5:A:520:SOG:O3	2.09	0.70
3:H:6:GLN:H	3:H:113:GLN:HE22	1.40	0.70
1:A:302:ALA:HB1	1:A:314:ILE:HD13	1.73	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:45:ILE:HB	1:B:46:PRO:HD3	1.74	0.69
5:A:520:SOG:H1'2	5:B:520:SOG:H2'2	1.74	0.69
3:I:6:GLN:H	3:I:113:GLN:HE22	1.41	0.69
1:A:136:ASP:HB2	1:A:137:PRO:HD2	1.77	0.67
1:B:136:ASP:HB2	1:B:137:PRO:HD2	1.77	0.67
3:I:159:THR:HB	3:I:206:ALA:HB3	1.77	0.66
1:A:45:ILE:HB	1:A:46:PRO:HD3	1.76	0.66
3:H:159:THR:HB	3:H:206:ALA:HB3	1.79	0.65
5:B:520:SOG:HO2	5:B:520:SOG:HO4	1.30	0.64
5:A:520:SOG:H2'1	5:B:520:SOG:H3'2	1.79	0.64
5:B:520:SOG:C4	5:B:520:SOG:O2	2.47	0.63
2:M:199:GLU:HG2	2:M:208:PRO:HB2	1.81	0.63
3:I:11:LEU:HD22	3:I:155:PRO:HG3	1.81	0.62
2:L:41:GLN:HB2	2:L:51:LEU:HD11	1.81	0.62
2:M:46:GLN:HB2	2:M:47:PRO:HD2	1.80	0.62
2:L:124:PRO:HD3	2:L:136:VAL:HG22	1.82	0.62
2:M:124:PRO:HD3	2:M:136:VAL:HG22	1.81	0.62
1:B:309:ALA:O	1:B:310:ASN:HB2	1.98	0.62
2:L:46:GLN:HB2	2:L:47:PRO:HD2	1.80	0.62
2:M:41:GLN:HB2	2:M:51:LEU:HD11	1.80	0.61
2:L:199:GLU:HG2	2:L:208:PRO:HB2	1.81	0.61
1:B:484:TRP:CE3	5:B:520:SOG:C5'	2.82	0.61
3:H:37:MET:HE1	3:H:47:TRP:HD1	1.64	0.61
1:A:309:ALA:O	1:A:310:ASN:HB2	2.01	0.61
3:I:218:ILE:HG22	3:I:219:GLU:N	2.13	0.60
1:B:18:MET:HG3	1:B:63:TRP:CZ2	2.37	0.60
1:A:438:ARG:HD3	3:I:52:TYR:CE1	2.37	0.60
3:I:37:MET:HE1	3:I:47:TRP:HD1	1.65	0.60
5:B:520:SOG:C4	5:B:520:SOG:HO2	2.13	0.59
1:B:253:PHE:HE2	1:B:467:TRP:CH2	2.18	0.59
2:L:154:ILE:HD12	2:L:196:TYR:CD2	2.38	0.59
1:A:18:MET:HG3	1:A:63:TRP:CZ2	2.38	0.58
3:H:156:GLU:HG3	3:H:157:PRO:HA	1.84	0.58
2:M:154:ILE:HD12	2:M:196:TYR:CD2	2.37	0.58
5:A:520:SOG:C2'	5:B:520:SOG:H3'2	2.32	0.58
1:A:77:THR:OG1	1:A:97:GLY:HA3	2.03	0.58
1:B:364:ILE:HD13	1:B:379:VAL:HG22	1.85	0.58
3:I:27:TYR:CZ	3:I:98:ARG:HD3	2.39	0.57
1:A:364:ILE:HD13	1:A:379:VAL:HG22	1.85	0.57
2:L:53:TYR:O	2:L:57:ASN:HB2	2.04	0.57
1:B:77:THR:OG1	1:B:97:GLY:HA3	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:27:TYR:CZ	3:H:98:ARG:HD3	2.39	0.57
2:M:53:TYR:O	2:M:57:ASN:HB2	2.05	0.57
1:A:253:PHE:HE2	1:A:467:TRP:CH2	2.19	0.57
3:I:11:LEU:HD12	3:I:118:THR:O	2.06	0.56
2:M:100:TYR:HE2	3:I:99:ARG:HD3	1.71	0.56
3:H:218:ILE:HG22	3:H:219:GLU:N	2.13	0.56
2:M:128:GLN:HG2	2:M:133:GLY:O	2.06	0.56
1:A:58:LEU:HD23	1:A:257:LEU:HD13	1.88	0.56
3:H:207:HIS:NE2	3:H:209:ALA:HB3	2.21	0.56
1:A:438:ARG:HD3	3:I:52:TYR:CZ	2.41	0.55
2:M:213:PHE:HD1	2:M:215:ARG:H	1.54	0.55
1:A:104:VAL:HA	1:A:107:TYR:CE2	2.41	0.55
2:L:13:VAL:HG21	2:L:82:VAL:HG21	1.88	0.55
1:B:58:LEU:HD23	1:B:257:LEU:HD13	1.89	0.55
3:I:207:HIS:NE2	3:I:209:ALA:HB3	2.22	0.55
1:B:51:PHE:CZ	1:B:56:ILE:HD11	2.42	0.55
3:H:35:ASN:HB3	3:H:50:GLU:HB3	1.88	0.54
1:A:51:PHE:CZ	1:A:56:ILE:HD11	2.43	0.54
2:L:128:GLN:HG2	2:L:133:GLY:O	2.06	0.54
3:H:156:GLU:HG3	3:H:157:PRO:CA	2.37	0.54
2:M:13:VAL:HG21	2:M:82:VAL:HG21	1.90	0.54
1:B:438:ARG:HD3	3:H:52:TYR:CE1	2.43	0.54
1:B:331:PHE:O	1:B:337:GLY:HA3	2.08	0.54
3:I:35:ASN:HB3	3:I:50:GLU:HB3	1.88	0.54
1:B:104:VAL:HA	1:B:107:TYR:CE2	2.41	0.54
2:L:213:PHE:HD1	2:L:215:ARG:H	1.54	0.54
1:B:179:ASN:HB3	1:B:360:MET:HE1	1.89	0.53
3:I:205:VAL:O	3:I:212:THR:HG23	2.08	0.53
1:A:402:GLU:CD	1:A:483:VAL:HG13	2.33	0.53
1:A:56:ILE:N	1:A:57:PRO:HD2	2.23	0.53
2:M:43:LYS:HD2	2:M:46:GLN:NE2	2.23	0.53
3:I:156:GLU:HG3	3:I:157:PRO:HA	1.89	0.53
1:A:287:GLU:O	1:A:291:VAL:HG22	2.09	0.53
1:B:56:ILE:N	1:B:57:PRO:HD2	2.24	0.53
1:A:331:PHE:O	1:A:337:GLY:HA3	2.08	0.53
2:L:43:LYS:HD2	2:L:46:GLN:NE2	2.24	0.53
1:A:485:ILE:HD13	1:B:485:ILE:HD13	1.90	0.53
1:A:179:ASN:HB3	1:A:360:MET:HE1	1.90	0.53
1:B:298:SER:HB3	1:B:322:LEU:HD21	1.90	0.52
1:B:402:GLU:CD	1:B:483:VAL:HG13	2.34	0.52
3:H:205:VAL:O	3:H:212:THR:HG23	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:218:ILE:CG2	3:H:219:GLU:H	2.17	0.52
2:M:3:VAL:HG22	2:M:26:SER:HB3	1.91	0.52
1:A:298:SER:HB3	1:A:322:LEU:HD21	1.91	0.52
1:B:5:ARG:HH22	1:B:369:ASP:CG	2.18	0.52
1:B:438:ARG:HD3	3:H:52:TYR:CZ	2.44	0.52
3:I:87:THR:HG22	3:I:88:SER:N	2.25	0.52
1:B:199:MET:HB2	1:B:200:PRO:HD3	1.92	0.52
1:B:155:PRO:HD3	1:B:162:LEU:HD23	1.92	0.51
1:A:230:LEU:HD11	1:A:299:ILE:HB	1.92	0.51
1:A:25:LEU:HD12	1:A:25:LEU:H	1.75	0.51
1:B:249:GLY:O	1:B:253:PHE:HD2	1.93	0.51
1:B:456:THR:N	1:B:457:PRO:HD2	2.26	0.51
1:B:287:GLU:O	1:B:291:VAL:HG22	2.10	0.51
1:A:456:THR:N	1:A:457:PRO:HD2	2.26	0.51
2:M:140:LEU:HD22	2:M:179:MET:HE3	1.92	0.51
3:H:156:GLU:CG	3:H:157:PRO:HA	2.40	0.51
1:A:199:MET:HB2	1:A:200:PRO:HD3	1.92	0.51
1:B:438:ARG:HD2	3:H:33:TYR:OH	2.10	0.51
3:H:192:THR:HG22	3:H:194:SER:H	1.76	0.51
1:A:44:MET:HA	1:A:44:MET:HE2	1.93	0.50
1:A:299:ILE:HB	1:A:300:PRO:HD3	1.92	0.50
1:A:438:ARG:HD2	3:I:33:TYR:OH	2.11	0.50
3:H:87:THR:HG22	3:H:88:SER:N	2.26	0.50
3:I:1:GLN:N	3:I:1:GLN:CD	2.70	0.50
1:A:495:LEU:HD22	5:A:521:SOG:H8'1	1.94	0.50
1:B:44:MET:HE2	1:B:44:MET:HA	1.93	0.50
3:I:156:GLU:HG3	3:I:157:PRO:CA	2.42	0.50
3:I:218:ILE:CG2	3:I:219:GLU:H	2.16	0.50
1:B:230:LEU:HD11	1:B:299:ILE:HB	1.92	0.50
2:L:140:LEU:HD22	2:L:179:MET:HE3	1.93	0.50
1:B:299:ILE:HB	1:B:300:PRO:HD3	1.93	0.50
3:H:11:LEU:HD12	3:H:118:THR:O	2.12	0.50
2:M:100:TYR:HE2	3:I:99:ARG:CD	2.24	0.49
1:A:155:PRO:HD3	1:A:162:LEU:HD23	1.93	0.49
1:B:25:LEU:HD12	1:B:25:LEU:H	1.76	0.49
3:I:192:THR:HG22	3:I:194:SER:H	1.76	0.49
1:A:5:ARG:HH22	1:A:369:ASP:CG	2.20	0.49
1:A:124:VAL:HG12	1:A:124:VAL:O	2.12	0.49
2:L:3:VAL:HG22	2:L:26:SER:HB3	1.93	0.49
1:A:249:GLY:O	1:A:253:PHE:HD2	1.94	0.49
2:L:128:GLN:HG3	3:H:130:TYR:CE2	2.48	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:M:129:LEU:O	2:M:187:LYS:HD2	2.12	0.49
1:B:59:MET:SD	1:B:283:ALA:HB2	2.53	0.49
3:H:6:GLN:HE21	3:H:112:GLY:CA	2.21	0.49
5:A:520:SOG:O3	5:A:520:SOG:C6	2.56	0.49
1:B:484:TRP:CZ3	5:B:520:SOG:H4'1	2.47	0.49
3:I:103:MET:HA	3:I:105:ARG:H	1.77	0.49
2:L:129:LEU:O	2:L:187:LYS:HD2	2.13	0.49
3:I:87:THR:HG22	3:I:88:SER:H	1.78	0.49
1:B:167:PHE:O	1:B:171:VAL:HG23	2.14	0.48
1:A:59:MET:SD	1:A:283:ALA:HB2	2.53	0.48
1:A:167:PHE:O	1:A:171:VAL:HG23	2.13	0.48
1:B:124:VAL:HG12	1:B:124:VAL:O	2.12	0.48
2:L:199:GLU:HG3	2:L:210:VAL:HG22	1.95	0.48
3:I:156:GLU:CG	3:I:157:PRO:HA	2.43	0.48
1:B:305:PHE:HZ	1:B:334:THR:CG2	2.28	0.47
1:B:475:ILE:HG22	1:B:487:ARG:HH21	1.79	0.47
3:H:87:THR:HG22	3:H:88:SER:H	1.79	0.47
1:A:305:PHE:HZ	1:A:334:THR:CG2	2.27	0.47
3:H:151:LYS:HG2	3:H:152:GLY:N	2.29	0.47
3:H:191:VAL:HG12	3:H:192:THR:H	1.79	0.47
1:A:395:PHE:CD1	5:B:520:SOG:C6	2.97	0.47
3:H:196:TRP:CG	3:H:197:PRO:HA	2.50	0.47
3:I:196:TRP:CG	3:I:197:PRO:HA	2.50	0.47
1:A:38:ASN:HB3	1:A:243:VAL:HG13	1.97	0.47
3:H:6:GLN:NE2	3:H:96:CYS:H	2.12	0.47
3:I:51:ILE:O	3:I:53:PRO:HD3	2.15	0.47
3:H:51:ILE:O	3:H:53:PRO:HD3	2.14	0.47
1:B:484:TRP:CH2	5:B:520:SOG:H4'1	2.51	0.47
3:H:103:MET:HA	3:H:105:ARG:H	1.77	0.47
3:I:6:GLN:NE2	3:I:96:CYS:H	2.13	0.47
1:A:57:PRO:O	1:A:61:ILE:HG13	2.15	0.46
1:B:31:PHE:HB3	1:B:32:PRO:HD3	1.98	0.46
2:M:43:LYS:NZ	2:M:85:GLU:O	2.48	0.46
1:B:421:ILE:O	1:B:425:TRP:HB2	2.15	0.46
2:M:199:GLU:HG3	2:M:210:VAL:HG22	1.96	0.46
1:B:57:PRO:O	1:B:61:ILE:HG13	2.15	0.46
1:A:31:PHE:HB3	1:A:32:PRO:HD3	1.97	0.46
1:A:65:MET:HE1	1:A:81:PHE:HE1	1.79	0.46
3:H:32:TYR:CE1	3:H:98:ARG:NH1	2.84	0.46
2:M:144:TYR:HA	2:M:145:PRO:C	2.41	0.46
3:I:151:LYS:HG2	3:I:152:GLY:N	2.30	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:421:ILE:O	1:A:425:TRP:HB2	2.16	0.46
2:L:43:LYS:NZ	2:L:85:GLU:O	2.48	0.46
1:B:38:ASN:HB3	1:B:243:VAL:HG13	1.97	0.46
1:A:25:LEU:HD12	1:A:25:LEU:N	2.30	0.46
1:B:65:MET:HE1	1:B:81:PHE:HE1	1.81	0.45
3:I:191:VAL:HG12	3:I:192:THR:H	1.80	0.45
1:B:25:LEU:HD12	1:B:25:LEU:N	2.31	0.45
2:L:144:TYR:HA	2:L:145:PRO:C	2.40	0.45
3:I:32:TYR:CE1	3:I:98:ARG:NH1	2.84	0.45
1:A:481:TRP:CZ3	1:B:161:ILE:HD11	2.51	0.45
1:A:475:ILE:HG22	1:A:487:ARG:HH21	1.81	0.45
3:H:207:HIS:HE2	3:H:209:ALA:HB3	1.82	0.45
3:I:50:GLU:CD	3:I:99:ARG:HH21	2.24	0.45
1:A:159:GLU:HA	1:A:160:PRO:HD3	1.82	0.45
2:L:48:PRO:HD2	3:H:111:TRP:CE3	2.52	0.45
3:H:191:VAL:HG12	3:H:192:THR:N	2.32	0.45
2:M:128:GLN:HG3	3:I:130:TYR:CE2	2.52	0.45
3:I:207:HIS:HE2	3:I:209:ALA:HB3	1.82	0.45
1:A:166:LEU:HB2	5:A:520:SOG:O3	2.17	0.44
2:M:119:VAL:HG22	2:M:140:LEU:HG	1.99	0.44
1:A:395:PHE:CD1	5:B:520:SOG:H62	2.52	0.44
1:B:481:TRP:HB2	5:B:521:SOG:H5'2	2.00	0.44
3:I:191:VAL:HG12	3:I:192:THR:N	2.33	0.44
2:L:36:TYR:CE2	3:H:106:TYR:CD2	3.06	0.43
3:H:50:GLU:CD	3:H:99:ARG:HH21	2.26	0.43
2:L:119:VAL:HG22	2:L:140:LEU:HG	2.00	0.43
3:H:1:GLN:CD	3:H:1:GLN:N	2.75	0.43
3:H:37:MET:O	3:H:94:TYR:HA	2.19	0.43
1:A:184:ILE:HA	1:A:379:VAL:HG21	2.01	0.43
1:A:42:ALA:HB1	1:A:238:LEU:HD11	2.01	0.43
1:A:420:LEU:C	1:A:420:LEU:HD23	2.43	0.43
1:A:420:LEU:HD23	1:A:420:LEU:O	2.19	0.43
1:B:184:ILE:HA	1:B:379:VAL:HG21	2.00	0.43
2:L:187:LYS:O	2:L:191:GLU:HG2	2.18	0.43
1:B:116:LEU:O	1:B:120:ILE:HG13	2.19	0.43
3:H:147:GLY:O	3:H:217:LYS:HE2	2.19	0.43
1:A:85:TRP:CG	1:A:426:ILE:HG21	2.54	0.42
1:A:259:PHE:HE2	6:A:608:HOH:O	2.01	0.42
1:B:85:TRP:CG	1:B:426:ILE:HG21	2.54	0.42
1:B:339:PHE:CE2	1:B:343:LEU:HD11	2.54	0.42
1:B:420:LEU:HD23	1:B:420:LEU:C	2.44	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:77:THR:N	1:A:78:PRO:CD	2.83	0.42
1:A:393:VAL:HG13	1:A:400:LEU:HB2	2.01	0.42
1:B:42:ALA:HB1	1:B:238:LEU:HD11	2.01	0.42
1:B:367:LEU:HB3	1:B:373:LEU:HD22	2.02	0.42
1:A:116:LEU:O	1:A:120:ILE:HG13	2.20	0.42
1:B:20:GLY:HA2	1:B:199:MET:SD	2.60	0.42
1:B:481:TRP:CB	5:B:521:SOG:H5'2	2.49	0.42
2:M:187:LYS:O	2:M:191:GLU:HG2	2.18	0.42
1:B:77:THR:N	1:B:78:PRO:CD	2.83	0.42
3:I:37:MET:O	3:I:94:TYR:HA	2.19	0.42
1:B:305:PHE:CE2	1:B:330:ILE:HG23	2.54	0.42
1:A:161:ILE:HD11	1:B:481:TRP:CE3	2.55	0.42
2:L:192:ARG:O	2:L:192:ARG:HG2	2.19	0.42
3:H:131:PRO:HG3	3:H:215:ASP:OD2	2.20	0.42
1:A:432:ALA:O	1:A:436:ILE:HG13	2.20	0.42
1:A:481:TRP:CE3	1:B:161:ILE:HD11	2.55	0.42
1:B:361:GLN:N	1:B:362:PRO:CD	2.83	0.42
1:B:393:VAL:HG13	1:B:400:LEU:HB2	2.01	0.42
3:I:147:GLY:O	3:I:217:LYS:HE2	2.20	0.42
1:A:116:LEU:HB2	1:A:349:PHE:CD1	2.55	0.41
2:M:192:ARG:O	2:M:192:ARG:HG2	2.19	0.41
3:I:6:GLN:HE21	3:I:112:GLY:CA	2.26	0.41
3:I:6:GLN:HE22	3:I:96:CYS:H	1.68	0.41
1:A:361:GLN:N	1:A:362:PRO:CD	2.83	0.41
1:B:381:TRP:O	1:B:385:ILE:HG13	2.20	0.41
2:L:140:LEU:HD23	2:L:148:ILE:HD11	2.02	0.41
3:H:102:THR:HG23	3:H:105:ARG:HG3	2.03	0.41
1:B:8:TRP:CD1	1:B:14:LEU:HD13	2.55	0.41
1:A:33:VAL:HG21	1:A:319:ALA:HA	2.03	0.41
1:A:339:PHE:CE2	1:A:343:LEU:HD11	2.55	0.41
3:H:160:LEU:C	3:H:160:LEU:HD23	2.46	0.41
3:I:207:HIS:CD2	3:I:210:SER:OG	2.73	0.41
1:A:305:PHE:CE2	1:A:330:ILE:HG23	2.55	0.41
3:I:130:TYR:HA	3:I:131:PRO:HD3	1.92	0.41
3:I:207:HIS:HD2	3:I:210:SER:OG	2.04	0.41
1:A:205:LEU:HD11	1:A:346:PHE:HD2	1.86	0.41
1:A:485:ILE:HD13	1:B:485:ILE:CD1	2.51	0.41
1:B:272:ASP:OD1	1:B:272:ASP:N	2.53	0.41
2:L:139:PHE:C	2:L:140:LEU:HD12	2.46	0.41
3:I:131:PRO:HG3	3:I:215:ASP:OD2	2.20	0.41
1:A:395:PHE:CD1	5:B:520:SOG:H61	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:400:LEU:HD23	1:A:400:LEU:C	2.46	0.41
1:B:116:LEU:HB2	1:B:349:PHE:CD1	2.56	0.41
3:H:105:ARG:HG3	3:H:105:ARG:O	2.21	0.41
3:H:207:HIS:CD2	3:H:210:SER:OG	2.74	0.41
1:A:8:TRP:CD1	1:A:14:LEU:HD13	2.55	0.41
1:A:77:THR:HG21	1:A:415:PHE:HZ	1.86	0.41
1:A:162:LEU:C	1:A:164:PRO:HD3	2.46	0.41
1:A:395:PHE:HD1	5:B:520:SOG:H61	1.86	0.41
1:B:400:LEU:HD23	1:B:400:LEU:C	2.46	0.41
3:H:189:VAL:HG13	3:H:189:VAL:O	2.21	0.41
2:M:140:LEU:HD23	2:M:148:ILE:HD11	2.02	0.41
1:B:33:VAL:HG21	1:B:319:ALA:HA	2.03	0.41
1:B:205:LEU:HD11	1:B:346:PHE:HD2	1.86	0.41
1:B:484:TRP:CD1	5:B:521:SOG:H7'2	2.55	0.41
2:M:139:PHE:C	2:M:140:LEU:HD12	2.46	0.41
1:B:420:LEU:HD23	1:B:420:LEU:O	2.21	0.40
2:L:214:ASN:O	2:L:215:ARG:C	2.64	0.40
3:I:189:VAL:O	3:I:189:VAL:HG13	2.21	0.40
3:H:154:PHE:HA	3:H:155:PRO:HA	1.80	0.40
2:M:65:ARG:HD2	2:M:81:PRO:O	2.21	0.40
3:I:105:ARG:O	3:I:105:ARG:HG3	2.21	0.40
1:A:253:PHE:HA	6:A:608:HOH:O	2.21	0.40
1:B:368:GLU:O	1:B:372:LYS:HA	2.22	0.40
3:I:219:GLU:HG2	3:I:219:GLU:OXT	2.21	0.40
1:A:20:GLY:HA2	1:A:199:MET:SD	2.61	0.40
1:A:368:GLU:O	1:A:372:LYS:HA	2.21	0.40
5:B:520:SOG:H3'1	5:B:520:SOG:H6'2	1.66	0.40
3:H:158:VAL:HG21	3:H:185:LEU:HD22	2.03	0.40
3:I:102:THR:HG23	3:I:105:ARG:HG3	2.03	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	494/519 (95%)	479 (97%)	14 (3%)	1 (0%)	43	73
1	B	494/519 (95%)	477 (97%)	16 (3%)	1 (0%)	43	73
2	L	213/218 (98%)	204 (96%)	9 (4%)	0	100	100
2	M	213/218 (98%)	205 (96%)	8 (4%)	0	100	100
3	H	208/219 (95%)	195 (94%)	13 (6%)	0	100	100
3	I	208/219 (95%)	196 (94%)	12 (6%)	0	100	100
All	All	1830/1912 (96%)	1756 (96%)	72 (4%)	2 (0%)	48	78

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	298	SER
1	B	298	SER

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	370/419 (88%)	369 (100%)	1 (0%)	86	87
1	B	370/419 (88%)	369 (100%)	1 (0%)	86	87
2	L	187/190 (98%)	187 (100%)	0	100	100
2	M	187/190 (98%)	187 (100%)	0	100	100
3	H	176/185 (95%)	176 (100%)	0	100	100
3	I	176/185 (95%)	176 (100%)	0	100	100
All	All	1466/1588 (92%)	1464 (100%)	2 (0%)	88	90

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	424	PHE
1	B	424	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	38	ASN
1	A	179	ASN
1	A	361	GLN
1	B	38	ASN
1	B	179	ASN
1	B	361	GLN
2	L	17	GLN
2	L	42	GLN
2	L	46	GLN
2	L	202	HIS
3	H	6	GLN
3	H	39	GLN
3	H	172	HIS
3	H	199	GLN
3	H	207	HIS
2	M	17	GLN
2	M	46	GLN
2	M	202	HIS
3	I	6	GLN
3	I	199	GLN
3	I	207	HIS

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 9 ligands modelled in this entry, 4 are monoatomic - leaving 5 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$
5	SOG	A	522	-	7,7,20	0.33	0	6,6,25	0.54	0
5	SOG	B	520	-	20,20,20	4.11	5 (25%)	24,25,25	9.15	10 (41%)
5	SOG	B	521	-	7,7,20	0.32	0	6,6,25	0.57	0
5	SOG	A	521	-	7,7,20	0.32	0	6,6,25	0.61	0
5	SOG	A	520	-	20,20,20	6.10	6 (30%)	24,25,25	7.35	15 (62%)

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
5	SOG	A	522	-	-	1/5/5/31	-
5	SOG	B	520	-	-	2/11/31/31	0/1/1/1
5	SOG	B	521	-	-	3/5/5/31	-
5	SOG	A	521	-	-	2/5/5/31	-
5	SOG	A	520	-	-	6/11/31/31	0/1/1/1

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	520	SOG	O5-C1	-21.74	1.09	1.42
5	B	520	SOG	O5-C1	-12.29	1.23	1.42
5	B	520	SOG	C3-C2	11.27	1.81	1.52
5	A	520	SOG	C1-C2	10.81	1.72	1.53
5	A	520	SOG	C4-C3	-7.10	1.33	1.52
5	A	520	SOG	C4-C5	6.97	1.67	1.53
5	A	520	SOG	C3-C2	6.38	1.68	1.52
5	B	520	SOG	O5-C5	-6.30	1.29	1.44
5	A	520	SOG	C6-C5	2.39	1.59	1.51
5	B	520	SOG	C1-C2	2.37	1.57	1.53
5	B	520	SOG	O4-C4	-2.05	1.37	1.43

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	520	SOG	C1-O5-C5	31.53	169.13	112.56
5	A	520	SOG	O5-C1-C2	-24.50	76.54	110.32
5	B	520	SOG	O5-C1-C2	-20.04	82.69	110.32
5	A	520	SOG	O2-C2-C1	-17.33	79.49	110.30
5	B	520	SOG	O2-C2-C1	-12.05	88.88	110.30
5	A	520	SOG	C4-C3-C2	-10.94	91.63	110.83
5	B	520	SOG	O5-C5-C6	10.38	132.17	106.44
5	B	520	SOG	O5-C1-S1	-9.51	83.82	109.94
5	A	520	SOG	C3-C4-C5	-9.50	93.00	110.23
5	B	520	SOG	O2-C2-C3	-9.11	88.89	110.38
5	B	520	SOG	O5-C5-C4	-8.74	93.94	109.70
5	B	520	SOG	C4-C3-C2	-7.47	97.71	110.83
5	B	520	SOG	C1-C2-C3	6.36	122.98	110.55
5	A	520	SOG	C1'-S1-C1	5.81	112.99	100.45
5	A	520	SOG	O5-C5-C6	5.71	120.59	106.44
5	A	520	SOG	C1-O5-C5	-5.36	102.95	112.56
5	A	520	SOG	O5-C1-S1	4.77	123.03	109.94
5	A	520	SOG	C1-C2-C3	4.20	118.76	110.55
5	A	520	SOG	O3-C3-C4	3.30	118.14	110.38
5	A	520	SOG	O2-C2-C3	2.99	117.43	110.38
5	A	520	SOG	C3'-C2'-C1'	2.83	123.82	113.10
5	A	520	SOG	O5-C5-C4	2.65	114.48	109.70
5	A	520	SOG	O4-C4-C5	2.33	115.07	109.32
5	A	520	SOG	C2'-C1'-S1	2.18	119.71	112.36
5	B	520	SOG	C3-C4-C5	-2.06	106.50	110.23

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	520	SOG	O5-C1-S1-C1'
5	A	520	SOG	C3'-C4'-C5'-C6'
5	B	520	SOG	C3'-C4'-C5'-C6'
5	A	520	SOG	C4-C5-C6-O6
5	B	521	SOG	C4'-C5'-C6'-C7'
5	B	520	SOG	C2'-C3'-C4'-C5'
5	A	521	SOG	C3'-C4'-C5'-C6'
5	B	521	SOG	C1'-C2'-C3'-C4'
5	A	521	SOG	C2'-C3'-C4'-C5'
5	B	521	SOG	C5'-C6'-C7'-C8'
5	A	520	SOG	C2'-C3'-C4'-C5'

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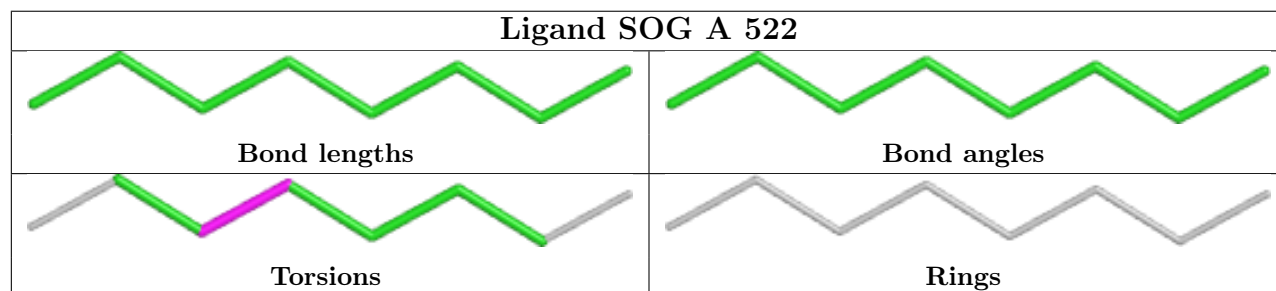
Mol	Chain	Res	Type	Atoms
5	A	522	SOG	C4'-C5'-C6'-C7'
5	A	520	SOG	C2'-C1'-S1-C1
5	A	520	SOG	C4'-C5'-C6'-C7'

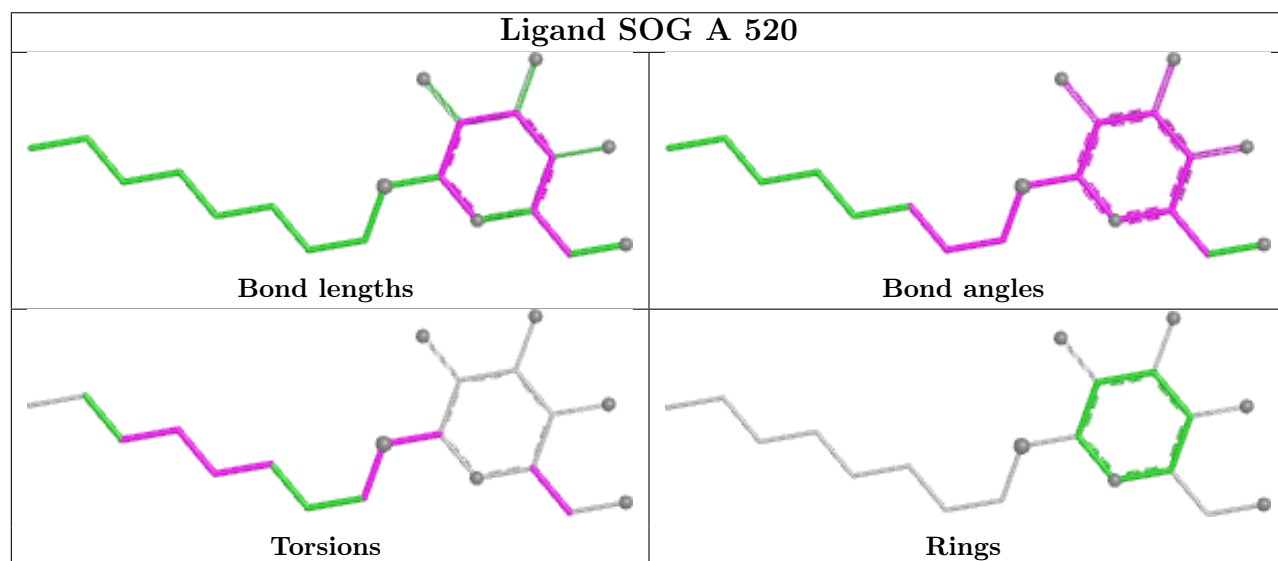
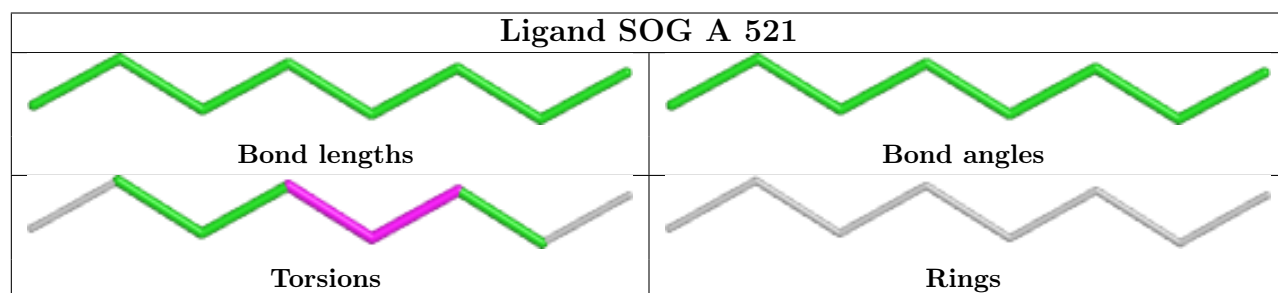
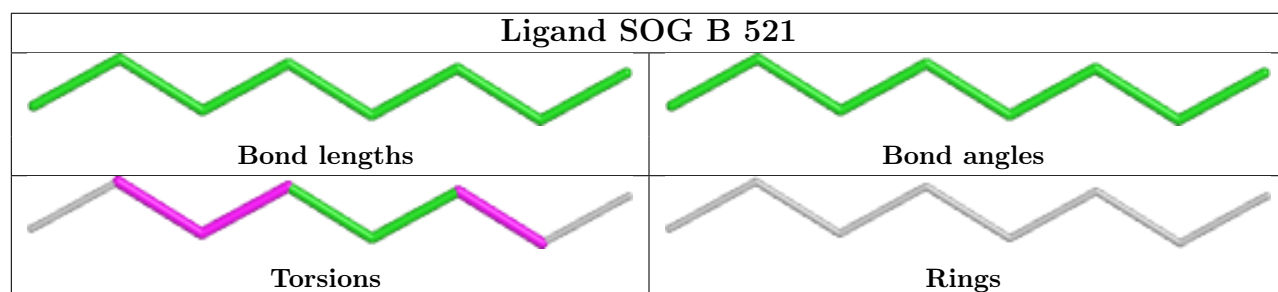
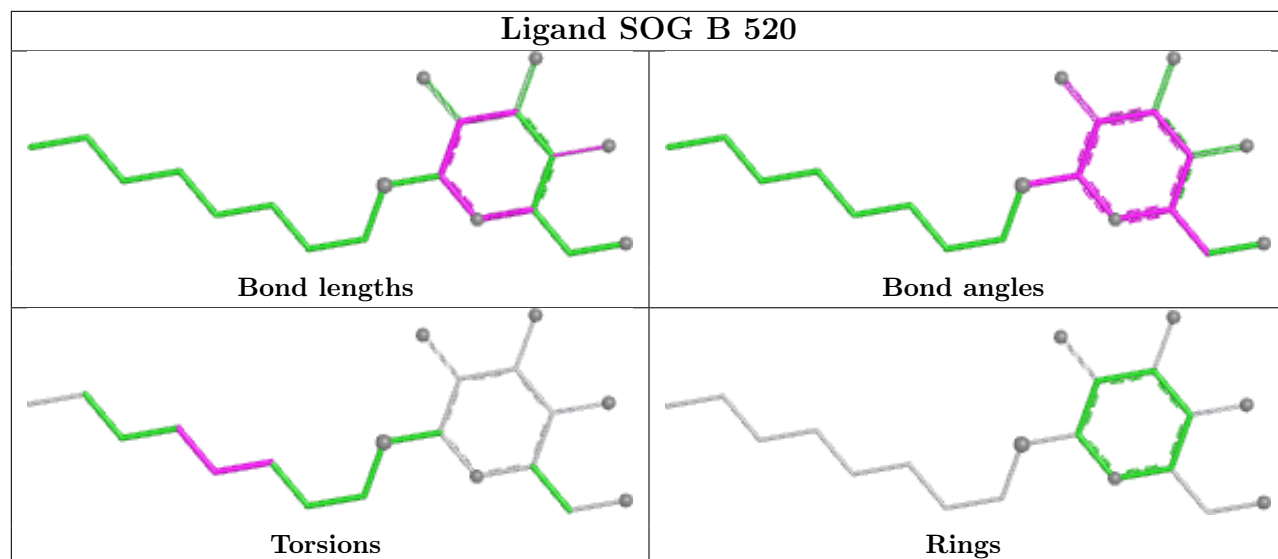
There are no ring outliers.

4 monomers are involved in 31 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	520	SOG	22	0
5	B	521	SOG	3	0
5	A	521	SOG	1	0
5	A	520	SOG	11	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

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	498/519 (95%)	0.19	37 (7%)	20 11	46, 71, 143, 194	0
1	B	498/519 (95%)	0.14	31 (6%)	26 14	48, 72, 145, 197	0
2	L	215/218 (98%)	0.10	3 (1%)	73 53	53, 79, 142, 189	0
2	M	215/218 (98%)	0.25	5 (2%)	61 39	51, 80, 142, 187	0
3	H	212/219 (96%)	0.37	12 (5%)	29 16	46, 76, 106, 129	0
3	I	212/219 (96%)	0.30	11 (5%)	33 17	54, 78, 107, 128	0
All	All	1850/1912 (96%)	0.21	99 (5%)	31 17	46, 75, 139, 197	0

All (99) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	479	THR	6.9
1	A	475	ILE	5.5
1	B	158	ASP	5.4
3	I	135	VAL	5.2
1	B	305	PHE	5.0
1	A	305	PHE	4.9
1	A	234	ASP	4.7
1	A	219	PRO	4.4
1	A	126	LEU	4.3
1	A	129	GLU	4.2
3	H	103	MET	4.1
1	A	472	ILE	4.1
1	B	236	GLU	4.0
3	H	215	ASP	3.8
1	B	231	TRP	3.7
1	B	308	VAL	3.7
3	I	103	MET	3.6
1	B	234	ASP	3.6
1	A	309	ALA	3.5

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Mol	Chain	Res	Type	RSRZ
1	B	472	ILE	3.5
3	I	212	THR	3.4
1	A	127	VAL	3.4
1	A	214	PHE	3.3
1	B	477	GLU	3.3
1	A	235	PHE	3.2
3	I	1	GLN	3.2
1	A	478	GLU	3.2
1	A	231	TRP	3.1
1	A	233	PRO	3.1
1	B	474	LYS	3.1
1	B	229	PHE	3.1
1	B	471	TYR	3.1
1	B	129	GLU	3.0
3	H	212	THR	3.0
2	L	154	ILE	3.0
1	A	474	LYS	2.9
1	A	128	PRO	2.9
1	B	470	GLU	2.9
1	A	311	ALA	2.8
1	B	230	LEU	2.8
1	A	218	THR	2.8
1	B	306	PHE	2.7
1	B	312	VAL	2.7
1	B	219	PRO	2.7
1	B	476	MET	2.7
1	B	298	SER	2.7
3	H	135	VAL	2.7
1	A	476	MET	2.6
1	B	272	ASP	2.6
3	I	215	ASP	2.6
1	A	308	VAL	2.5
3	H	101	VAL	2.5
1	A	158	ASP	2.5
1	B	214	PHE	2.5
1	A	157	GLY	2.5
1	B	478	GLU	2.5
1	B	135	THR	2.5
3	I	42	GLY	2.4
1	B	232	THR	2.4
1	B	468	ALA	2.4
2	M	134	ALA	2.4

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Mol	Chain	Res	Type	RSRZ
3	H	9	ALA	2.4
2	L	35	SER	2.4
3	I	143	SER	2.4
1	A	137	PRO	2.4
3	I	101	VAL	2.4
1	B	218	THR	2.4
1	A	480	HIS	2.4
1	B	270	ARG	2.3
1	B	86	ARG	2.3
3	I	57	THR	2.3
1	B	72	GLN	2.3
3	H	1	GLN	2.3
1	A	224	ALA	2.2
1	B	479	THR	2.2
1	A	136	ASP	2.2
1	A	217	GLU	2.2
1	A	471	TYR	2.2
1	A	223	ALA	2.2
3	H	192	THR	2.2
3	H	17	SER	2.2
1	A	314	ILE	2.2
2	M	197	THR	2.1
3	I	210	SER	2.1
1	A	89	PHE	2.1
1	A	144	PHE	2.1
2	L	158	GLU	2.1
1	A	240	ASP	2.1
1	A	306	PHE	2.1
3	H	193	SER	2.1
3	I	78	SER	2.1
2	M	194	ASN	2.1
2	M	13	VAL	2.1
2	M	15	LEU	2.1
1	B	315	ALA	2.1
1	A	215	LEU	2.0
1	A	230	LEU	2.0
3	H	102	THR	2.0
3	H	214	VAL	2.0

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

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

6.3 Carbohydrates

There are no oligosaccharides in this entry.

6.4 Ligands

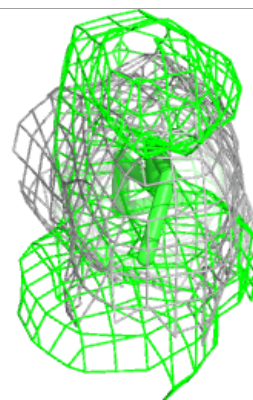
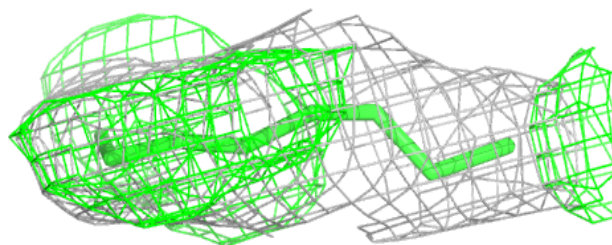
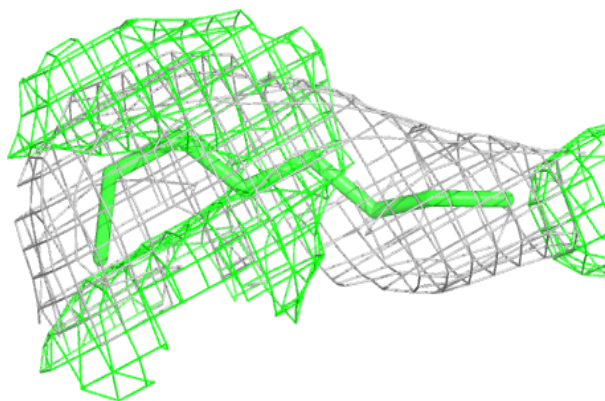
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	NA	B	601	1/1	0.48	0.16	89,89,89,89	0
5	SOG	B	521	8/20	0.67	0.34	51,59,64,64	0
5	SOG	B	520	20/20	0.75	0.18	56,106,142,154	0
5	SOG	A	522	8/20	0.80	0.26	93,105,110,113	0
5	SOG	A	520	20/20	0.83	0.16	69,111,130,142	0
5	SOG	A	521	8/20	0.88	0.19	49,59,74,75	0
4	NA	A	602	1/1	0.90	0.08	82,82,82,82	0
4	NA	B	602	1/1	0.91	0.11	122,122,122,122	0
4	NA	A	601	1/1	0.95	0.07	77,77,77,77	0

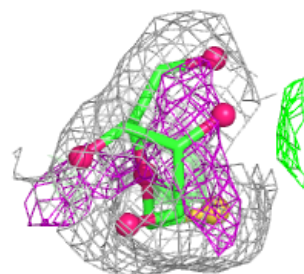
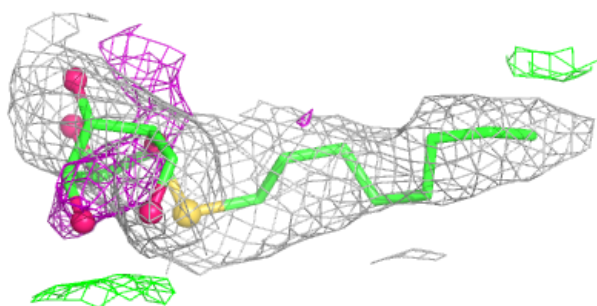
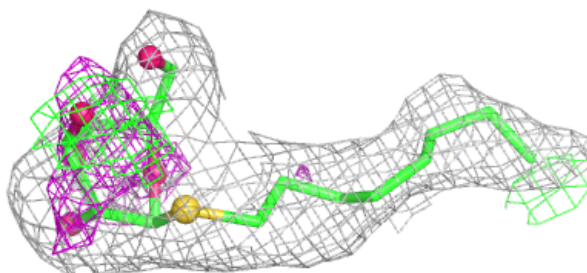
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around SOG B 521:

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

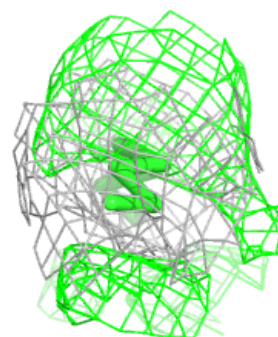
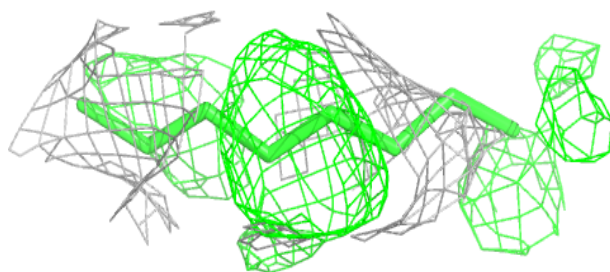
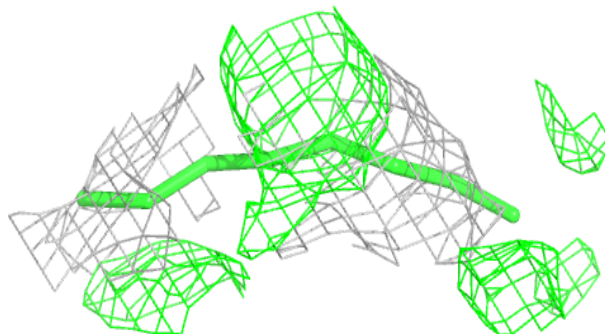
**Electron density around SOG B 520:**

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

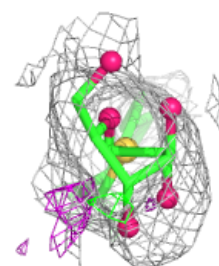
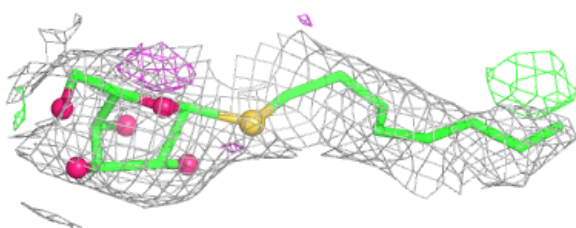
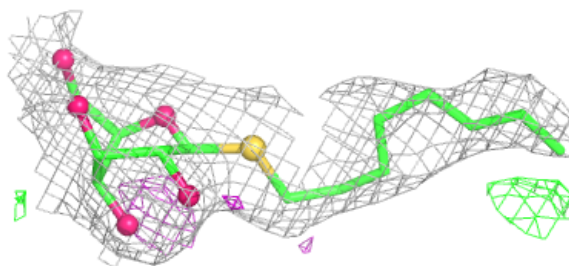


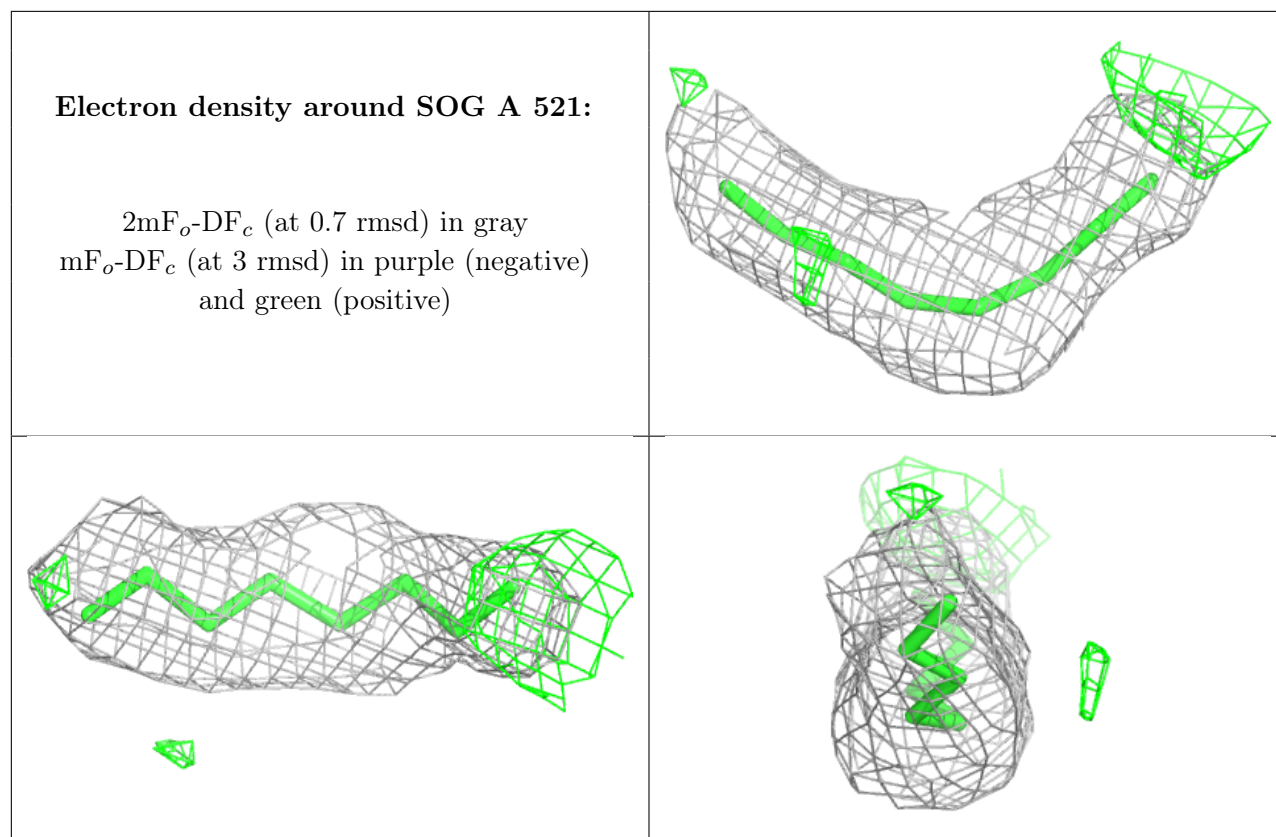
Electron density around SOG A 522:

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

**Electron density around SOG A 520:**

$2mF_o-DF_c$ (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.