



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 02:03 AM UTC

PDB ID : 2VT4 / pdb_00002vt4
Title : TURKEY BETA1 ADRENERGIC RECEPTOR WITH STABILISING MUTATIONS AND BOUND CYANOPINDOLOL
Authors : Warne, A.; Serrano-Vega, M.J.; Baker, J.G.; Moukhametzianov, R.; Edwards, P.C.; Henderson, R.; Leslie, A.G.W.; Tate, C.G.; Schertler, G.F.X.
Deposited on : 2008-05-09
Resolution : 2.70 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

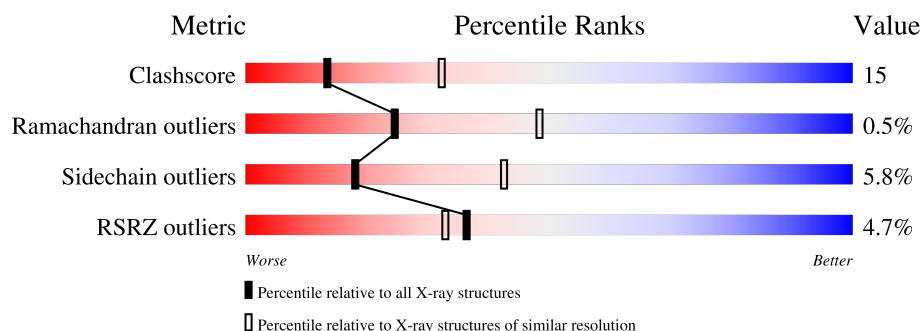
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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))
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	313	2% 61% 23% • 12%
1	B	313	6% 59% 27% • 12%
1	C	313	4% 60% 22% • 16%
1	D	313	4% 57% 27% • 13%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 8976 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 BETA1 ADRENERGIC RECEPTOR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	274	Total	C	N	O	S	0	0	0
			2175	1441	355	361	18			
1	B	276	Total	C	N	O	S	0	0	0
			2195	1455	358	364	18			
1	C	263	Total	C	N	O	S	0	0	0
			2081	1377	337	349	18			
1	D	272	Total	C	N	O	S	0	0	0
			2156	1427	352	359	18			

There are 32 discrepancies between the modelled and reference sequences:

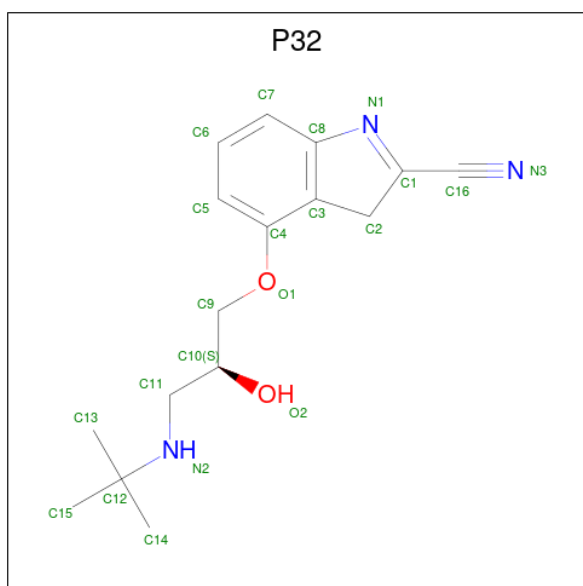
Chain	Residue	Modelled	Actual	Comment	Reference
A	68	SER	ARG	engineered mutation	UNP P07700
A	90	VAL	MET	engineered mutation	UNP P07700
A	116	LEU	CYS	engineered mutation	UNP P07700
A	227	ALA	TYR	engineered mutation	UNP P07700
A	282	LEU	ALA	engineered mutation	UNP P07700
A	327	ALA	PHE	engineered mutation	UNP P07700
A	338	MET	PHE	engineered mutation	UNP P07700
A	358	ALA	CYS	engineered mutation	UNP P07700
B	68	SER	ARG	engineered mutation	UNP P07700
B	90	VAL	MET	engineered mutation	UNP P07700
B	116	LEU	CYS	engineered mutation	UNP P07700
B	227	ALA	TYR	engineered mutation	UNP P07700
B	282	LEU	ALA	engineered mutation	UNP P07700
B	327	ALA	PHE	engineered mutation	UNP P07700
B	338	MET	PHE	engineered mutation	UNP P07700
B	358	ALA	CYS	engineered mutation	UNP P07700
C	68	SER	ARG	engineered mutation	UNP P07700
C	90	VAL	MET	engineered mutation	UNP P07700
C	116	LEU	CYS	engineered mutation	UNP P07700
C	227	ALA	TYR	engineered mutation	UNP P07700
C	282	LEU	ALA	engineered mutation	UNP P07700

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Chain	Residue	Modelled	Actual	Comment	Reference
C	327	ALA	PHE	engineered mutation	UNP P07700
C	338	MET	PHE	engineered mutation	UNP P07700
C	358	ALA	CYS	engineered mutation	UNP P07700
D	68	SER	ARG	engineered mutation	UNP P07700
D	90	VAL	MET	engineered mutation	UNP P07700
D	116	LEU	CYS	engineered mutation	UNP P07700
D	227	ALA	TYR	engineered mutation	UNP P07700
D	282	LEU	ALA	engineered mutation	UNP P07700
D	327	ALA	PHE	engineered mutation	UNP P07700
D	338	MET	PHE	engineered mutation	UNP P07700
D	358	ALA	CYS	engineered mutation	UNP P07700

- Molecule 2 is Cyanopindolol (CCD ID: P32) (formula: $C_{16}H_{21}N_3O_2$).

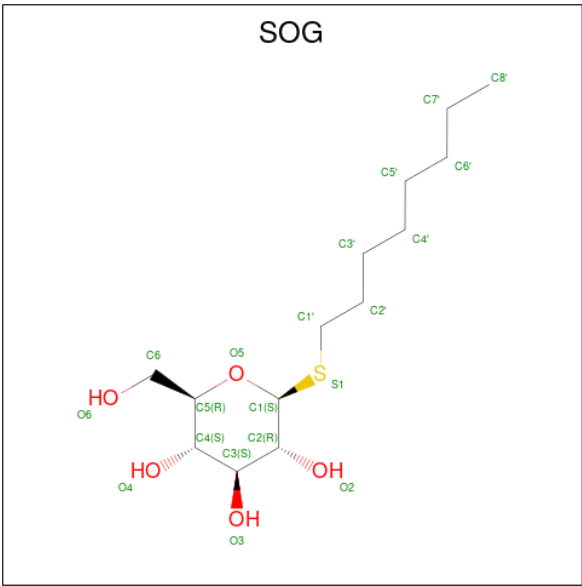


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			21	16	3	2		
2	B	1	Total	C	N	O	0	0
			21	16	3	2		
2	C	1	Total	C	N	O	0	0
			21	16	3	2		
2	D	1	Total	C	N	O	0	0
			21	16	3	2		

- Molecule 3 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Na 1 1	0	0
3	B	1	Total Na 1 1	0	0
3	C	1	Total Na 1 1	0	0
3	D	1	Total Na 1 1	0	0

- Molecule 4 is octyl 1-thio-beta-D-glucopyranoside (CCD ID: SOG) (formula: C₁₄H₂₈O₅S).



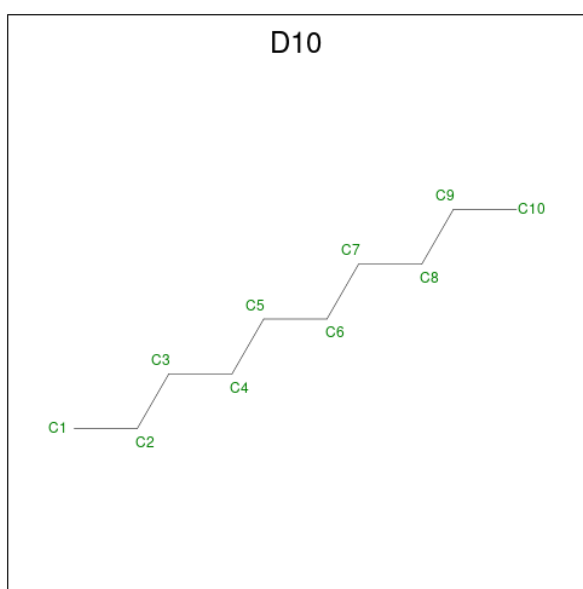
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O S 20 14 5 1	0	0
4	A	1	Total C O S 20 14 5 1	0	0
4	A	1	Total C O S 20 14 5 1	0	0
4	B	1	Total C O S 20 14 5 1	0	0
4	B	1	Total C S 10 9 1	0	0
4	B	1	Total C O S 20 14 5 1	0	0
4	C	1	Total C O S 20 14 5 1	0	0
4	C	1	Total C O S 20 14 5 1	0	0

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	C	1	Total	C	O	S	0	0
			20	14	5	1		
4	D	1	Total	C	O	S	0	0
			20	14	5	1		
4	D	1	Total	C	O	S	0	0
			20	14	5	1		
4	D	1	Total	C	O	S	0	0
			20	14	5	1		

- Molecule 5 is DECANE (CCD ID: D10) (formula: C₁₀H₂₂).



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

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	9	Total	O	0	0
			9	9		
6	B	7	Total	O	0	0
			7	7		
6	C	7	Total	O	0	0
			7	7		

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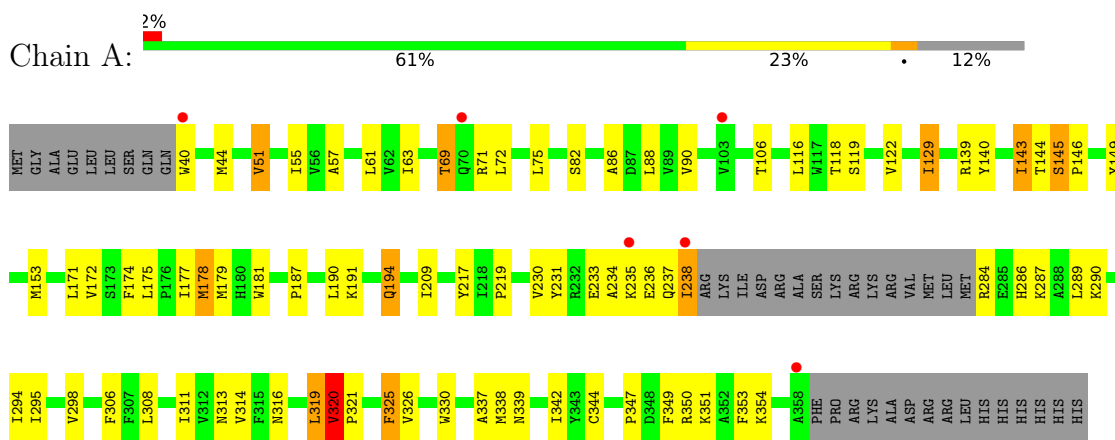
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	D	8	Total	O	0	0
			8	8		

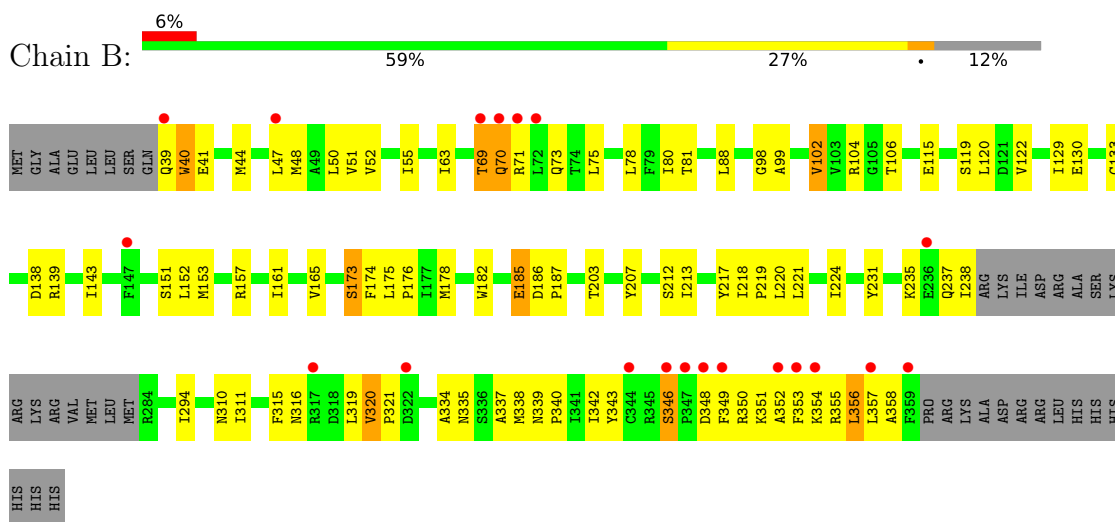
3 Residue-property plots [i](#)

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.

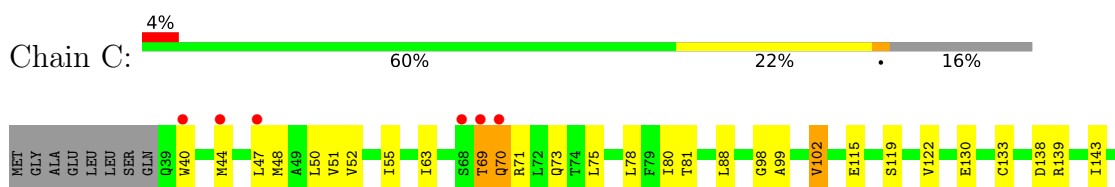
• Molecule 1: BETA1 ADRENERGIC RECEPTOR

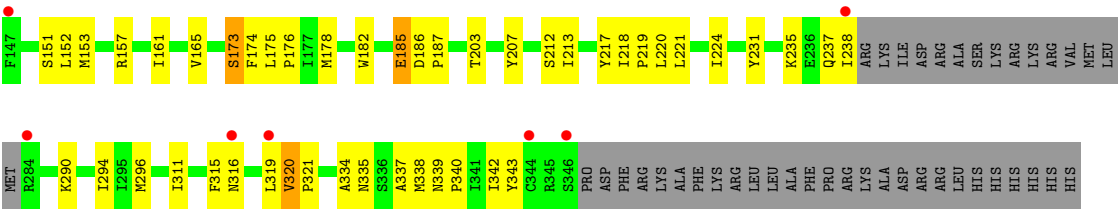


• Molecule 1: BETA1 ADRENERGIC RECEPTOR

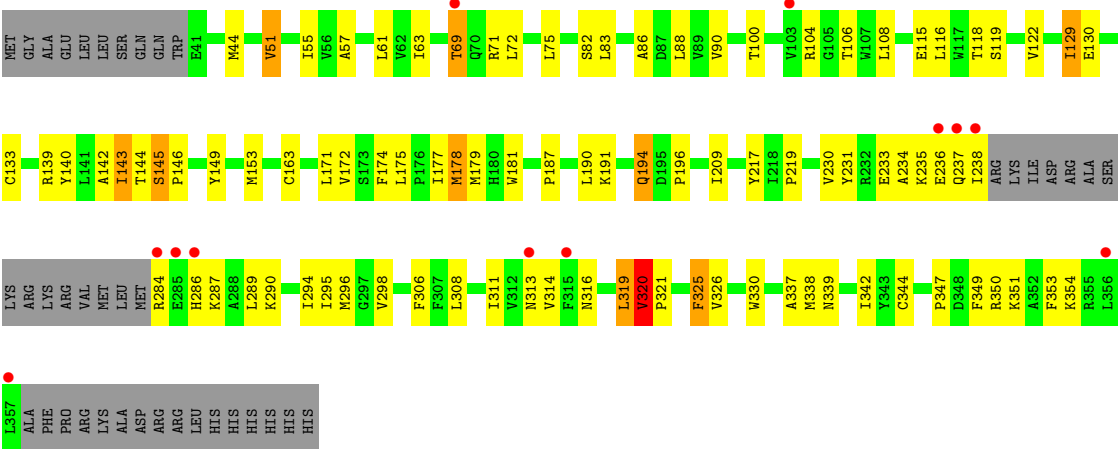


• Molecule 1: BETA1 ADRENERGIC RECEPTOR





● Molecule 1: BETA1 ADRENERGIC RECEPTOR



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	55.50Å 86.80Å 95.50Å 67.60° 73.30° 85.80°	Depositor
Resolution (Å)	45.10 – 2.70 45.10 – 2.70	Depositor EDS
% Data completeness (in resolution range)	95.8 (45.10-2.70) 95.8 (45.10-2.70)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.72 (at 2.69Å)	Xtriage
Refinement program	PHENIX (PHENIX.PHASER)	Depositor
R, R_{free}	0.212 , 0.268 0.205 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	39.3	Xtriage
Anisotropy	0.322	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 65.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	8976	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.11% 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: SOG, NA, D10, P32

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	0.60	0/2229	0.93	6/3042 (0.2%)
1	B	0.58	0/2250	0.88	4/3070 (0.1%)
1	C	0.57	0/2132	0.87	2/2913 (0.1%)
1	D	0.60	0/2208	0.93	7/3012 (0.2%)
All	All	0.59	0/8819	0.90	19/12037 (0.2%)

There are no bond length outliers.

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	82	SER	N-CA-C	-7.01	103.72	111.36
1	D	82	SER	N-CA-C	-6.94	103.80	111.36
1	A	51	VAL	N-CA-C	6.72	116.87	110.42
1	D	51	VAL	N-CA-C	6.61	116.76	110.42
1	D	320	VAL	CA-C-N	6.39	126.89	119.99
1	D	320	VAL	C-N-CA	6.39	126.89	119.99
1	C	320	VAL	CA-C-N	6.37	126.59	119.90
1	C	320	VAL	C-N-CA	6.37	126.59	119.90
1	B	320	VAL	CA-C-N	6.35	126.57	119.90
1	B	320	VAL	C-N-CA	6.35	126.57	119.90
1	A	320	VAL	CA-C-N	6.28	126.78	119.99
1	A	320	VAL	C-N-CA	6.28	126.78	119.99
1	D	338	MET	N-CA-C	6.13	118.04	111.36
1	D	51	VAL	CB-CA-C	-5.85	104.49	111.97
1	A	338	MET	N-CA-C	5.56	117.42	111.36
1	A	51	VAL	CB-CA-C	-5.47	104.96	111.97
1	D	196	PRO	N-CA-C	-5.08	107.12	114.18
1	B	346	SER	CA-C-N	5.04	126.14	119.84
1	B	346	SER	C-N-CA	5.04	126.14	119.84

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2175	0	2250	65	0
1	B	2195	0	2267	68	0
1	C	2081	0	2145	60	0
1	D	2156	0	2235	74	0
2	A	21	0	21	0	0
2	B	21	0	21	2	0
2	C	21	0	21	1	0
2	D	21	0	21	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	60	0	83	7	0
4	B	50	0	73	5	0
4	C	60	0	84	11	0
4	D	60	0	84	10	0
5	A	10	0	22	0	0
5	C	10	0	22	0	0
6	A	9	0	0	0	0
6	B	7	0	0	1	0
6	C	7	0	0	0	0
6	D	8	0	0	0	0
All	All	8976	0	9349	268	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 15.

All (268) 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:C:290:LYS:HA	4:C:405:SOG:H2'1	1.47	0.95

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:104:ARG:HG2	1:D:108:LEU:HD11	1.58	0.84
1:C:115:GLU:HB3	1:C:178:MET:HE1	1.62	0.82
1:B:115:GLU:HB3	1:B:178:MET:HE1	1.63	0.81
1:C:122:VAL:HG12	1:C:173:SER:HB2	1.64	0.80
1:D:238:ILE:HB	1:D:286:HIS:CE1	2.16	0.80
1:B:122:VAL:HG12	1:B:173:SER:HB2	1.65	0.78
1:D:295:ILE:HD12	1:D:339:ASN:ND2	2.00	0.77
1:D:313:ASN:ND2	1:D:325:PHE:HE1	1.81	0.77
1:C:290:LYS:HG3	4:C:405:SOG:H1'2	1.66	0.76
1:A:295:ILE:HD12	1:A:339:ASN:ND2	2.02	0.75
1:A:313:ASN:ND2	1:A:325:PHE:HE1	1.84	0.74
1:D:235:LYS:HD3	1:D:236:GLU:OE2	1.88	0.74
1:B:346:SER:O	1:B:350:ARG:HB2	1.89	0.72
1:A:235:LYS:O	1:A:238:ILE:HG12	1.90	0.71
1:C:98:GLY:O	1:C:102:VAL:HG23	1.93	0.69
1:B:98:GLY:O	1:B:102:VAL:HG23	1.95	0.67
1:A:69:THR:HG22	1:A:71:ARG:HB3	1.75	0.67
1:A:233:GLU:O	1:A:237:GLN:HB2	1.95	0.67
1:D:69:THR:HG22	1:D:71:ARG:HB3	1.75	0.67
1:D:233:GLU:O	1:D:237:GLN:HB2	1.95	0.67
1:B:106:THR:HG21	4:D:403:SOG:H1	1.78	0.66
1:B:70:GLN:HA	1:B:73:GLN:HB2	1.77	0.65
1:A:71:ARG:HH22	1:A:351:LYS:NZ	1.96	0.64
1:D:129:ILE:HG13	1:D:219:PRO:HB2	1.78	0.64
1:D:71:ARG:HH22	1:D:351:LYS:NZ	1.96	0.64
1:C:70:GLN:HA	1:C:73:GLN:HB2	1.77	0.64
1:A:235:LYS:C	1:A:237:GLN:H	2.04	0.63
1:B:231:TYR:CD1	4:B:405:SOG:H3'1	2.33	0.63
1:A:129:ILE:HG13	1:A:219:PRO:HB2	1.79	0.63
1:B:157:ARG:HG2	4:B:404:SOG:H2'1	1.80	0.62
1:D:313:ASN:ND2	1:D:325:PHE:CE1	2.65	0.62
1:B:316:ASN:ND2	1:B:319:LEU:HB2	2.14	0.62
1:D:178:MET:HA	4:D:403:SOG:C1'	2.28	0.62
4:A:402:SOG:O5	4:A:402:SOG:H2'2	2.00	0.62
1:C:316:ASN:ND2	1:C:319:LEU:HB2	2.15	0.62
1:D:296:MET:SD	4:D:405:SOG:H8'3	2.40	0.62
1:C:48:MET:HG2	1:C:99:ALA:HB1	1.81	0.61
1:B:48:MET:HG2	1:B:99:ALA:HB1	1.80	0.61
1:D:234:ALA:O	1:D:286:HIS:HE1	1.84	0.61
1:C:217:TYR:CE1	1:C:311:ILE:HD11	2.36	0.61
1:A:313:ASN:ND2	1:A:325:PHE:CE1	2.67	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:186:ASP:OD1	1:B:187:PRO:HD2	2.02	0.60
1:B:217:TYR:CE1	1:B:311:ILE:HD11	2.36	0.60
1:C:186:ASP:OD1	1:C:187:PRO:HD2	2.02	0.60
1:B:294:ILE:HG22	1:B:342:ILE:HD13	1.83	0.60
1:D:55:ILE:HG13	1:D:337:ALA:HB2	1.83	0.60
1:A:55:ILE:HG13	1:A:337:ALA:HB2	1.83	0.60
1:C:40:TRP:CE2	1:C:44:MET:HE3	2.36	0.60
1:C:218:ILE:HB	1:C:219:PRO:HD3	1.83	0.59
1:B:346:SER:HB3	1:B:349:PHE:HB2	1.85	0.59
1:B:40:TRP:CE2	1:B:44:MET:HE3	2.36	0.59
1:B:161:ILE:HG12	4:B:404:SOG:H6'1	1.84	0.59
1:B:218:ILE:HB	1:B:219:PRO:HD3	1.83	0.59
1:C:161:ILE:HG12	4:C:404:SOG:H6'1	1.85	0.59
1:B:213:ILE:HA	1:B:217:TYR:HB2	1.84	0.58
1:C:294:ILE:HG22	1:C:342:ILE:HD13	1.85	0.58
1:A:234:ALA:O	1:A:286:HIS:HE1	1.87	0.57
1:C:213:ILE:HA	1:C:217:TYR:HB2	1.84	0.57
1:A:289:LEU:O	4:A:405:SOG:H3'2	2.05	0.57
1:B:231:TYR:HD1	4:B:405:SOG:H3'1	1.69	0.57
1:C:231:TYR:HD1	4:C:405:SOG:H3'1	1.68	0.57
1:B:185:GLU:OE1	1:B:185:GLU:HA	2.05	0.56
1:B:350:ARG:HG2	1:B:354:LYS:NZ	2.20	0.56
1:D:290:LYS:HG3	4:D:405:SOG:S1	2.45	0.56
1:A:116:LEU:O	1:A:119:SER:HB2	2.05	0.56
1:D:349:PHE:HB3	1:D:353:PHE:CE2	2.41	0.56
1:B:335:ASN:ND2	6:B:2007:HOH:O	2.22	0.56
1:A:294:ILE:HG22	1:A:342:ILE:HD13	1.87	0.55
1:A:143:ILE:O	1:A:143:ILE:HG23	2.06	0.55
1:D:116:LEU:O	1:D:119:SER:HB2	2.07	0.55
1:B:353:PHE:O	1:B:357:LEU:HD12	2.07	0.55
1:D:235:LYS:C	1:D:237:GLN:H	2.13	0.55
1:C:185:GLU:OE1	1:C:185:GLU:HA	2.07	0.55
1:B:338:MET:O	1:B:342:ILE:HG13	2.07	0.54
1:A:349:PHE:HB3	1:A:353:PHE:CE2	2.42	0.54
1:D:238:ILE:HB	1:D:286:HIS:HE1	1.70	0.54
1:A:181:TRP:CD1	4:A:402:SOG:O6	2.61	0.54
1:D:209:ILE:HD11	1:D:314:VAL:HG11	1.88	0.54
1:D:234:ALA:O	1:D:286:HIS:CE1	2.60	0.54
1:A:140:TYR:HB2	1:A:230:VAL:CG2	2.38	0.54
1:A:295:ILE:HD12	1:A:339:ASN:CG	2.33	0.54
1:C:122:VAL:HG12	1:C:173:SER:CB	2.37	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:143:ILE:HG23	1:D:143:ILE:O	2.07	0.54
1:D:295:ILE:HD12	1:D:339:ASN:CG	2.33	0.54
1:C:207:TYR:C	1:C:207:TYR:CD2	2.86	0.53
1:D:140:TYR:HB2	1:D:230:VAL:CG2	2.39	0.53
1:A:71:ARG:O	1:A:71:ARG:HG2	2.07	0.53
1:A:209:ILE:HD11	1:A:314:VAL:HG11	1.89	0.53
1:B:207:TYR:C	1:B:207:TYR:CD2	2.86	0.53
1:B:104:ARG:HH22	1:D:115:GLU:CD	2.17	0.53
1:D:294:ILE:HG22	1:D:342:ILE:HD13	1.89	0.53
1:A:40:TRP:O	1:A:44:MET:HG2	2.09	0.53
1:C:235:LYS:O	1:C:238:ILE:HG22	2.09	0.52
1:C:338:MET:O	1:C:342:ILE:HG13	2.09	0.52
1:C:339:ASN:HB2	1:C:340:PRO:HD3	1.90	0.52
1:C:48:MET:O	1:C:51:VAL:HG22	2.09	0.52
1:B:122:VAL:HG12	1:B:173:SER:CB	2.38	0.52
1:A:284:ARG:HD3	1:A:287:LYS:HD2	1.91	0.52
1:C:296:MET:HE1	4:C:405:SOG:C8'	2.40	0.52
1:A:234:ALA:O	1:A:286:HIS:CE1	2.63	0.52
1:A:347:PRO:O	1:A:351:LYS:HG3	2.10	0.52
1:C:75:LEU:O	1:C:78:LEU:HB2	2.10	0.52
1:D:284:ARG:HD3	1:D:287:LYS:HD2	1.91	0.52
1:A:326:VAL:O	1:A:330:TRP:HD1	1.93	0.52
1:B:119:SER:HB3	1:B:174:PHE:CD2	2.45	0.52
1:D:326:VAL:O	1:D:330:TRP:HD1	1.93	0.52
1:A:75:LEU:HD13	1:A:153:MET:CG	2.40	0.51
1:B:48:MET:HG2	1:B:99:ALA:CB	2.40	0.51
1:C:119:SER:HB3	1:C:174:PHE:CD2	2.45	0.51
1:A:231:TYR:HD1	4:A:405:SOG:H1'1	1.75	0.51
1:A:171:LEU:HD12	1:A:175:LEU:HD12	1.92	0.51
1:B:235:LYS:O	1:B:238:ILE:HG22	2.10	0.51
1:B:339:ASN:HB2	1:B:340:PRO:HD3	1.91	0.51
1:D:71:ARG:O	1:D:71:ARG:HG2	2.09	0.51
1:D:44:MET:HA	1:D:44:MET:HE2	1.92	0.51
1:D:75:LEU:HD13	1:D:153:MET:CG	2.40	0.51
1:B:119:SER:HB3	1:B:174:PHE:CE2	2.46	0.51
1:D:171:LEU:HD12	1:D:175:LEU:HD12	1.92	0.51
1:C:48:MET:HG2	1:C:99:ALA:CB	2.41	0.50
1:B:182:TRP:O	1:B:203:THR:HA	2.11	0.50
1:C:182:TRP:O	1:C:203:THR:HA	2.12	0.50
1:A:118:THR:O	1:A:122:VAL:HG23	2.10	0.50
1:B:48:MET:O	1:B:51:VAL:HG22	2.10	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:178:MET:HE2	1:B:178:MET:HA	1.92	0.50
1:A:44:MET:HA	1:A:44:MET:HE2	1.92	0.50
1:D:347:PRO:O	1:D:351:LYS:HG3	2.12	0.50
1:A:316:ASN:HB3	1:A:319:LEU:HB2	1.92	0.50
1:C:119:SER:HB3	1:C:174:PHE:CE2	2.46	0.50
1:D:316:ASN:HB3	1:D:319:LEU:HB2	1.92	0.50
1:C:231:TYR:CD1	4:C:405:SOG:H3'1	2.47	0.50
1:D:118:THR:O	1:D:122:VAL:HG23	2.11	0.49
1:C:178:MET:HE2	1:C:178:MET:HA	1.93	0.49
1:C:319:LEU:O	1:C:319:LEU:HG	2.13	0.49
1:B:152:LEU:O	1:B:157:ARG:NH1	2.40	0.49
1:D:71:ARG:HH22	1:D:351:LYS:HZ2	1.58	0.49
4:A:405:SOG:H3'2	4:A:405:SOG:H6'1	1.53	0.49
1:B:319:LEU:HG	1:B:319:LEU:O	2.13	0.49
1:B:351:LYS:O	1:B:355:ARG:HG3	2.13	0.49
1:B:75:LEU:O	1:B:78:LEU:HB2	2.13	0.49
1:B:75:LEU:HD13	1:B:153:MET:HG3	1.94	0.49
1:C:152:LEU:O	1:C:157:ARG:NH1	2.42	0.48
1:D:178:MET:HB2	4:D:403:SOG:O6	2.14	0.48
1:C:75:LEU:HD13	1:C:153:MET:HG3	1.94	0.48
1:D:235:LYS:C	1:D:237:GLN:N	2.70	0.48
1:A:174:PHE:O	1:A:178:MET:HG2	2.12	0.48
1:A:290:LYS:HA	4:A:405:SOG:H2'1	1.95	0.48
1:D:145:SER:N	1:D:146:PRO:CD	2.76	0.48
1:C:55:ILE:HG12	1:C:337:ALA:HB2	1.96	0.48
1:C:130:GLU:HB3	1:C:165:VAL:HG13	1.95	0.48
1:D:86:ALA:O	1:D:90:VAL:HG23	2.13	0.48
1:A:139:ARG:HA	1:A:139:ARG:NE	2.28	0.48
1:D:306:PHE:C	1:D:306:PHE:CD2	2.92	0.48
1:D:174:PHE:O	1:D:178:MET:HG2	2.13	0.48
1:A:306:PHE:C	1:A:306:PHE:CD2	2.92	0.47
1:B:55:ILE:HG12	1:B:337:ALA:HB2	1.96	0.47
1:C:122:VAL:HG13	2:C:400:P32:C6	2.44	0.47
1:A:71:ARG:HH22	1:A:351:LYS:HZ2	1.59	0.47
1:B:39:GLN:C	1:B:41:GLU:N	2.72	0.47
1:D:71:ARG:HH12	1:D:351:LYS:NZ	2.11	0.47
1:A:71:ARG:HH12	1:A:351:LYS:NZ	2.12	0.47
1:A:86:ALA:O	1:A:90:VAL:HG23	2.14	0.47
1:D:57:ALA:O	1:D:61:LEU:HB2	2.15	0.47
1:C:235:LYS:C	1:C:237:GLN:H	2.23	0.47
1:D:75:LEU:HD13	1:D:153:MET:HG2	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:145:SER:N	1:A:146:PRO:CD	2.78	0.47
1:B:130:GLU:HB3	1:B:165:VAL:HG13	1.96	0.47
1:D:217:TYR:CE1	1:D:311:ILE:HD11	2.50	0.47
1:D:231:TYR:HD1	4:D:405:SOG:H3'1	1.79	0.47
1:B:69:THR:HG22	1:B:71:ARG:HB3	1.96	0.47
1:D:139:ARG:HA	1:D:139:ARG:NE	2.29	0.47
1:B:235:LYS:C	1:B:237:GLN:H	2.23	0.47
1:B:217:TYR:O	1:B:221:LEU:HG	2.15	0.46
1:C:51:VAL:HG23	1:C:52:VAL:N	2.31	0.46
1:B:63:ILE:HG12	1:B:81:THR:HG23	1.97	0.46
1:B:122:VAL:HG13	2:B:400:P32:C6	2.44	0.46
1:C:217:TYR:O	1:C:221:LEU:HG	2.15	0.46
1:C:44:MET:O	1:C:48:MET:HB2	2.16	0.46
1:A:143:ILE:HG22	1:A:144:THR:HG23	1.96	0.46
1:A:187:PRO:O	1:A:191:LYS:HG3	2.15	0.46
1:B:51:VAL:HG23	1:B:52:VAL:N	2.31	0.46
1:D:289:LEU:O	4:D:405:SOG:H4'1	2.15	0.46
1:A:57:ALA:O	1:A:61:LEU:HB2	2.15	0.46
1:A:75:LEU:HD13	1:A:153:MET:HG2	1.97	0.46
1:C:69:THR:HG22	1:C:71:ARG:HB3	1.97	0.46
1:A:238:ILE:HG23	1:A:286:HIS:HE1	1.82	0.45
1:B:120:LEU:HD23	1:B:120:LEU:HA	1.79	0.45
1:D:143:ILE:HG22	1:D:144:THR:HG23	1.98	0.45
1:C:63:ILE:HG12	1:C:81:THR:HG23	1.98	0.45
4:C:402:SOG:O6	1:D:163:CYS:HB3	2.16	0.45
1:A:238:ILE:HD13	4:A:405:SOG:O3	2.17	0.45
1:B:44:MET:O	1:B:48:MET:HB2	2.17	0.45
1:A:217:TYR:CE1	1:A:311:ILE:HD11	2.52	0.44
1:C:290:LYS:CA	4:C:405:SOG:H2'1	2.33	0.44
1:D:178:MET:HA	4:D:403:SOG:H1'1	1.96	0.44
1:A:235:LYS:O	1:A:237:GLN:N	2.51	0.44
1:D:179:MET:HE2	1:D:181:TRP:NE1	2.31	0.44
1:C:157:ARG:HG2	4:C:404:SOG:H2'1	1.99	0.44
1:A:344:CYS:O	1:A:350:ARG:HD3	2.18	0.44
1:D:69:THR:HB	1:D:72:LEU:HB2	2.00	0.44
1:B:352:ALA:O	1:B:356:LEU:HD12	2.18	0.44
1:C:290:LYS:HG3	4:C:405:SOG:C1'	2.44	0.44
1:D:100:THR:O	1:D:104:ARG:HB3	2.18	0.44
1:D:235:LYS:O	1:D:237:GLN:N	2.51	0.44
1:A:308:LEU:HA	1:A:308:LEU:HD23	1.73	0.44
1:A:320:VAL:HA	1:A:321:PRO:HD3	1.78	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:69:THR:CG2	1:B:71:ARG:HB3	2.48	0.44
1:A:179:MET:HE2	1:A:181:TRP:NE1	2.32	0.43
1:A:149:TYR:CD2	1:A:149:TYR:C	2.97	0.43
1:A:294:ILE:O	1:A:298:VAL:HG23	2.19	0.43
1:D:344:CYS:O	1:D:350:ARG:HD3	2.19	0.43
1:C:55:ILE:CG1	1:C:337:ALA:HB2	2.48	0.43
1:D:88:LEU:HD12	1:D:88:LEU:HA	1.69	0.43
1:D:143:ILE:CG2	1:D:144:THR:HG23	2.49	0.43
1:D:179:MET:HE2	1:D:181:TRP:HE1	1.83	0.43
1:A:69:THR:HB	1:A:72:LEU:HB2	2.01	0.43
1:B:315:PHE:CD1	1:B:315:PHE:N	2.86	0.43
1:C:69:THR:CG2	1:C:71:ARG:HB3	2.49	0.43
1:B:122:VAL:CG1	1:B:173:SER:HB2	2.44	0.42
1:D:149:TYR:CD2	1:D:149:TYR:C	2.97	0.42
1:D:187:PRO:O	1:D:191:LYS:HG3	2.17	0.42
1:A:129:ILE:HD12	1:A:129:ILE:O	2.19	0.42
1:B:55:ILE:CG1	1:B:337:ALA:HB2	2.49	0.42
1:B:75:LEU:CD1	1:B:153:MET:HG3	2.49	0.42
1:B:310:ASN:HD21	2:B:400:P32:C16	2.31	0.42
1:B:80:ILE:HD12	1:B:343:TYR:OH	2.20	0.42
1:B:152:LEU:HD23	1:B:152:LEU:HA	1.83	0.42
1:C:296:MET:CE	4:C:405:SOG:C8'	2.97	0.42
1:D:294:ILE:O	1:D:298:VAL:HG23	2.20	0.42
1:B:220:LEU:HG	1:B:224:ILE:HD12	2.00	0.42
1:C:80:ILE:HD12	1:C:343:TYR:OH	2.20	0.42
1:D:296:MET:HE1	4:D:405:SOG:C8'	2.50	0.42
1:A:88:LEU:HD12	1:A:88:LEU:HA	1.69	0.42
1:B:320:VAL:HG13	1:B:321:PRO:HD2	2.00	0.42
1:C:152:LEU:HD23	1:C:152:LEU:HA	1.83	0.42
1:A:75:LEU:HD13	1:A:153:MET:HG3	2.01	0.42
1:A:143:ILE:CG2	1:A:144:THR:HG23	2.49	0.42
1:B:175:LEU:HB2	1:B:176:PRO:HD3	2.01	0.42
1:D:320:VAL:HA	1:D:321:PRO:HD3	1.79	0.42
1:D:350:ARG:O	1:D:354:LYS:HG3	2.19	0.42
1:A:179:MET:HE2	1:A:181:TRP:HE1	1.84	0.41
1:A:350:ARG:O	1:A:354:LYS:HG3	2.19	0.41
1:C:315:PHE:N	1:C:315:PHE:CD1	2.87	0.41
1:C:75:LEU:CD1	1:C:153:MET:HG3	2.50	0.41
1:D:308:LEU:HA	1:D:308:LEU:HD23	1.75	0.41
1:C:220:LEU:HG	1:C:224:ILE:HD12	2.02	0.41
1:A:119:SER:HB3	1:A:174:PHE:CE2	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:139:ARG:NE	1:C:139:ARG:HA	2.35	0.41
1:D:190:LEU:O	1:D:194:GLN:HG2	2.21	0.41
1:B:231:TYR:CE1	4:B:405:SOG:H3'1	2.55	0.41
1:A:140:TYR:HB2	1:A:230:VAL:HG22	2.02	0.41
1:B:88:LEU:HD12	1:B:88:LEU:HA	1.93	0.41
1:C:334:ALA:O	1:C:335:ASN:C	2.64	0.41
1:B:334:ALA:O	1:B:335:ASN:C	2.64	0.41
1:C:175:LEU:HB2	1:C:176:PRO:HD3	2.02	0.41
1:C:339:ASN:CB	1:C:340:PRO:HD3	2.51	0.41
1:C:88:LEU:HD12	1:C:88:LEU:HA	1.93	0.41
1:D:238:ILE:HD12	1:D:286:HIS:ND1	2.36	0.41
1:D:296:MET:CE	4:D:405:SOG:C8'	2.98	0.41
1:A:190:LEU:O	1:A:194:GLN:HG2	2.21	0.40
1:A:238:ILE:HG23	1:A:286:HIS:CE1	2.56	0.40
1:B:139:ARG:NE	1:B:139:ARG:HA	2.36	0.40
1:D:129:ILE:HD12	1:D:129:ILE:O	2.20	0.40
1:D:142:ALA:HB2	1:D:149:TYR:CE1	2.56	0.40
1:D:130:GLU:O	1:D:133:CYS:HB2	2.21	0.40
1:D:83:LEU:HD12	1:D:83:LEU:HA	1.85	0.40
1:B:339:ASN:CB	1:B:340:PRO:HD3	2.52	0.40
1:C:320:VAL:HG13	1:C:321:PRO:HD2	2.02	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	270/313 (86%)	258 (96%)	10 (4%)	2 (1%)	18	41
1	B	272/313 (87%)	249 (92%)	21 (8%)	2 (1%)	18	41
1	C	259/313 (83%)	241 (93%)	18 (7%)	0	100	100
1	D	268/313 (86%)	254 (95%)	13 (5%)	1 (0%)	30	54

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	1069/1252 (85%)	1002 (94%)	62 (6%)	5 (0%)	24	48

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	236	GLU
1	B	40	TRP
1	A	69	THR
1	D	69	THR
1	B	358	ALA

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	236/271 (87%)	222 (94%)	14 (6%)	18	42
1	B	238/271 (88%)	223 (94%)	15 (6%)	16	39
1	C	227/271 (84%)	215 (95%)	12 (5%)	20	46
1	D	235/271 (87%)	222 (94%)	13 (6%)	19	45
All	All	936/1084 (86%)	882 (94%)	54 (6%)	18	42

All (54) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	51	VAL
1	A	63	ILE
1	A	106	THR
1	A	129	ILE
1	A	143	ILE
1	A	145	SER
1	A	172	VAL
1	A	177	ILE
1	A	178	MET
1	A	194	GLN

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Mol	Chain	Res	Type
1	A	238	ILE
1	A	319	LEU
1	A	320	VAL
1	A	325	PHE
1	B	47	LEU
1	B	50	LEU
1	B	69	THR
1	B	70	GLN
1	B	102	VAL
1	B	129	ILE
1	B	133	CYS
1	B	138	ASP
1	B	143	ILE
1	B	151	SER
1	B	173	SER
1	B	185	GLU
1	B	212	SER
1	B	348	ASP
1	B	356	LEU
1	C	47	LEU
1	C	50	LEU
1	C	69	THR
1	C	70	GLN
1	C	102	VAL
1	C	133	CYS
1	C	138	ASP
1	C	143	ILE
1	C	151	SER
1	C	173	SER
1	C	185	GLU
1	C	212	SER
1	D	51	VAL
1	D	63	ILE
1	D	106	THR
1	D	129	ILE
1	D	143	ILE
1	D	145	SER
1	D	172	VAL
1	D	177	ILE
1	D	178	MET
1	D	194	GLN
1	D	319	LEU

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Mol	Chain	Res	Type
1	D	320	VAL
1	D	325	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	286	HIS
1	A	313	ASN
1	B	310	ASN
1	C	310	ASN
1	D	237	GLN
1	D	286	HIS
1	D	313	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 22 ligands modelled in this entry, 4 are monoatomic - leaving 18 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	SOG	D	405	-	20,20,20	0.53	0	24,25,25	0.99	0
4	SOG	D	403	-	20,20,20	0.63	0	24,25,25	2.33	7 (29%)
4	SOG	A	405	-	20,20,20	1.02	2 (10%)	24,25,25	1.12	3 (12%)
4	SOG	D	402	-	20,20,20	1.23	3 (15%)	24,25,25	1.49	2 (8%)
4	SOG	A	402	-	20,20,20	1.92	6 (30%)	24,25,25	2.27	7 (29%)
2	P32	C	400	-	21,22,22	2.70	3 (14%)	25,31,31	1.97	5 (20%)
5	D10	C	406	-	9,9,9	0.63	0	8,8,8	0.64	0
2	P32	A	400	-	21,22,22	2.72	3 (14%)	25,31,31	1.56	5 (20%)
4	SOG	B	402	-	20,20,20	0.91	1 (5%)	24,25,25	1.88	4 (16%)
4	SOG	A	403	-	20,20,20	0.56	0	24,25,25	0.89	0
4	SOG	B	405	-	20,20,20	0.78	1 (5%)	24,25,25	0.81	1 (4%)
5	D10	A	406	-	9,9,9	0.69	0	8,8,8	0.57	0
2	P32	B	400	-	21,22,22	2.72	3 (14%)	25,31,31	1.84	5 (20%)
4	SOG	C	405	-	20,20,20	0.64	0	24,25,25	1.29	3 (12%)
4	SOG	B	404	-	9,9,20	0.63	0	8,8,25	1.17	1 (12%)
4	SOG	C	404	-	20,20,20	0.91	1 (5%)	24,25,25	1.06	2 (8%)
2	P32	D	400	-	21,22,22	2.80	3 (14%)	25,31,31	1.64	6 (24%)
4	SOG	C	402	-	20,20,20	0.72	1 (5%)	24,25,25	1.61	3 (12%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	SOG	D	405	-	-	0/11/31/31	0/1/1/1
4	SOG	D	403	-	-	2/11/31/31	0/1/1/1
4	SOG	A	405	-	-	0/11/31/31	0/1/1/1
4	SOG	D	402	-	-	2/11/31/31	0/1/1/1
4	SOG	A	402	-	-	2/11/31/31	0/1/1/1
2	P32	C	400	-	-	0/11/21/21	0/2/2/2
5	D10	C	406	-	-	3/7/7/7	-
2	P32	A	400	-	-	4/11/21/21	0/2/2/2
4	SOG	B	402	-	-	2/11/31/31	0/1/1/1
4	SOG	A	403	-	-	0/11/31/31	0/1/1/1
4	SOG	B	405	-	-	0/11/31/31	0/1/1/1
5	D10	A	406	-	-	6/7/7/7	-
2	P32	B	400	-	-	0/11/21/21	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	SOG	C	405	-	-	0/11/31/31	0/1/1/1
4	SOG	B	404	-	-	0/7/7/31	-
4	SOG	C	404	-	-	0/11/31/31	0/1/1/1
2	P32	D	400	-	-	4/11/21/21	0/2/2/2
4	SOG	C	402	-	-	2/11/31/31	0/1/1/1

All (27) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	400	P32	C2-C3	-11.40	1.32	1.50
2	B	400	P32	C2-C3	-11.17	1.32	1.50
2	A	400	P32	C2-C3	-11.04	1.33	1.50
2	C	400	P32	C2-C3	-10.75	1.33	1.50
4	A	402	SOG	O5-C1	5.55	1.51	1.42
2	D	400	P32	C2-C1	-3.72	1.32	1.50
4	A	402	SOG	C1-S1	3.52	1.86	1.80
2	C	400	P32	C2-C1	-3.50	1.33	1.50
4	A	402	SOG	O5-C5	3.33	1.52	1.44
2	A	400	P32	C2-C1	-3.31	1.34	1.50
2	A	400	P32	C8-N1	-3.24	1.33	1.39
2	B	400	P32	C2-C1	-3.22	1.34	1.50
2	D	400	P32	C8-N1	-3.09	1.33	1.39
2	C	400	P32	C8-N1	-3.00	1.33	1.39
4	B	402	SOG	O5-C1	2.97	1.47	1.42
2	B	400	P32	C8-N1	-2.85	1.34	1.39
4	D	402	SOG	O5-C1	2.80	1.46	1.42
4	A	405	SOG	O5-C1	2.54	1.46	1.42
4	A	405	SOG	C1-S1	2.38	1.84	1.80
4	C	402	SOG	O5-C1	2.35	1.46	1.42
4	A	402	SOG	C6-C5	2.33	1.59	1.51
4	C	404	SOG	O5-C1	2.32	1.46	1.42
4	A	402	SOG	C4-C5	2.11	1.57	1.53
4	D	402	SOG	C6-C5	2.05	1.58	1.51
4	B	405	SOG	O5-C1	2.05	1.45	1.42
4	A	402	SOG	C1'-S1	2.03	1.84	1.81
4	D	402	SOG	C1-S1	2.01	1.84	1.80

All (54) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	400	P32	C11-N2-C12	-7.52	108.76	116.52
4	A	402	SOG	O5-C5-C6	6.95	123.67	106.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	400	P32	C11-N2-C12	-6.95	109.35	116.52
4	D	403	SOG	O5-C5-C6	-6.56	90.18	106.44
4	B	402	SOG	C6-C5-C4	-6.11	98.03	113.02
4	D	403	SOG	C6-C5-C4	-5.06	100.58	113.02
4	D	402	SOG	C6-C5-C4	4.43	123.89	113.02
2	D	400	P32	C11-N2-C12	-4.38	112.00	116.52
4	C	402	SOG	C6-C5-C4	-4.15	102.83	113.02
4	C	402	SOG	C1'-S1-C1	3.92	108.92	100.45
2	A	400	P32	C11-N2-C12	-3.91	112.48	116.52
4	C	405	SOG	C1'-S1-C1	-3.83	92.17	100.45
4	D	403	SOG	C1-C2-C3	-3.60	103.51	110.55
4	B	402	SOG	C1'-S1-C1	3.50	108.00	100.45
4	D	403	SOG	C1-O5-C5	-3.41	106.44	112.56
4	A	402	SOG	C6-C5-C4	3.38	121.32	113.02
4	A	405	SOG	O5-C5-C6	3.35	114.75	106.44
4	A	402	SOG	C4'-C3'-C2'	3.23	130.69	114.37
4	A	402	SOG	C1'-S1-C1	3.11	107.17	100.45
4	A	402	SOG	O5-C1-S1	3.06	118.33	109.94
2	D	400	P32	C2-C3-C8	-2.97	108.06	109.52
2	C	400	P32	C2-C1-N1	-2.93	111.00	114.45
2	A	400	P32	C11-C10-C9	-2.87	106.15	110.54
2	C	400	P32	O1-C4-C3	2.84	119.71	115.75
4	C	402	SOG	C2'-C1'-S1	-2.81	102.90	112.36
2	D	400	P32	C2-C1-N1	-2.81	111.14	114.45
2	B	400	P32	O1-C4-C3	2.80	119.65	115.75
4	A	402	SOG	O2-C2-C1	2.77	115.22	110.30
4	D	402	SOG	O5-C5-C6	2.73	113.20	106.44
4	B	402	SOG	C2'-C1'-S1	-2.64	103.46	112.36
4	D	403	SOG	O5-C1-C2	2.64	113.96	110.32
2	B	400	P32	C2-C1-N1	-2.63	111.35	114.45
4	B	405	SOG	O5-C5-C6	2.58	112.82	106.44
2	A	400	P32	O1-C4-C3	2.58	119.34	115.75
2	A	400	P32	C2-C1-N1	-2.56	111.43	114.45
2	D	400	P32	C11-C10-C9	-2.53	106.66	110.54
4	C	404	SOG	O5-C5-C6	2.47	112.57	106.44
2	D	400	P32	O1-C4-C3	2.43	119.14	115.75
2	A	400	P32	C2-C3-C8	-2.40	108.34	109.52
4	B	404	SOG	C3'-C2'-C1'	-2.38	104.08	113.10
4	C	405	SOG	C6'-C5'-C4'	-2.37	102.39	114.37
4	D	403	SOG	O4-C4-C3	-2.26	105.04	110.38
4	A	402	SOG	O4-C4-C5	2.24	114.84	109.32
4	B	402	SOG	O5-C5-C6	2.23	111.97	106.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	400	P32	C11-C10-C9	-2.18	107.20	110.54
4	C	405	SOG	C4'-C3'-C2'	-2.15	103.50	114.37
4	A	405	SOG	C3'-C2'-C1'	2.15	121.24	113.10
2	D	400	P32	C2-C3-C4	2.13	131.49	127.78
4	D	403	SOG	C2'-C1'-S1	-2.12	105.21	112.36
4	A	405	SOG	C5'-C4'-C3'	-2.11	103.69	114.37
2	C	400	P32	O1-C4-C5	-2.09	119.38	123.95
4	C	404	SOG	O5-C5-C4	-2.04	106.02	109.70
2	C	400	P32	C2-C3-C4	2.03	131.33	127.78
2	B	400	P32	O1-C4-C5	-2.01	119.57	123.95

There are no chirality outliers.

All (27) torsion outliers are listed below:

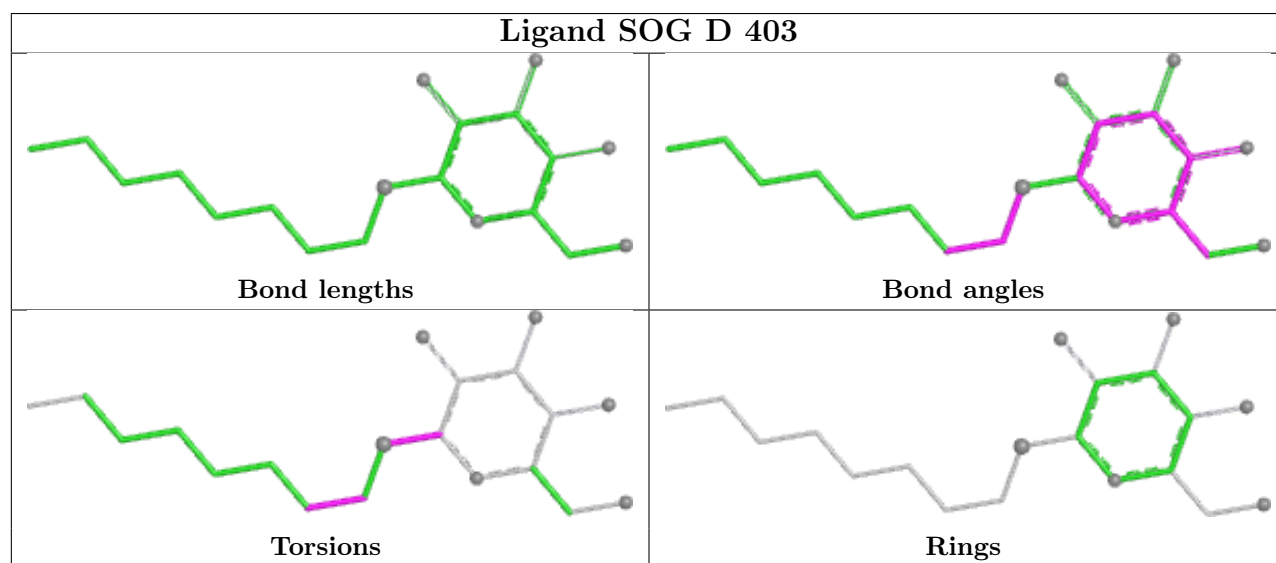
Mol	Chain	Res	Type	Atoms
2	A	400	P32	C11-C10-C9-O1
2	D	400	P32	C11-C10-C9-O1
4	A	402	SOG	O5-C1-S1-C1'
4	B	402	SOG	C2-C1-S1-C1'
4	B	402	SOG	O5-C1-S1-C1'
4	C	402	SOG	O5-C1-S1-C1'
4	D	403	SOG	O5-C1-S1-C1'
2	A	400	P32	O2-C10-C9-O1
2	D	400	P32	O2-C10-C9-O1
4	A	402	SOG	C4-C5-C6-O6
5	A	406	D10	C3-C4-C5-C6
5	A	406	D10	C2-C3-C4-C5
5	A	406	D10	C1-C2-C3-C4
5	A	406	D10	C4-C5-C6-C7
5	A	406	D10	C7-C8-C9-C10
5	C	406	D10	C3-C4-C5-C6
4	C	402	SOG	C2-C1-S1-C1'
4	D	402	SOG	C4-C5-C6-O6
5	A	406	D10	C6-C7-C8-C9
5	C	406	D10	C7-C8-C9-C10
5	C	406	D10	C1-C2-C3-C4
2	A	400	P32	C5-C4-O1-C9
2	D	400	P32	C5-C4-O1-C9
4	D	403	SOG	S1-C1'-C2'-C3'
4	D	402	SOG	C2-C1-S1-C1'
2	A	400	P32	C3-C4-O1-C9
2	D	400	P32	C3-C4-O1-C9

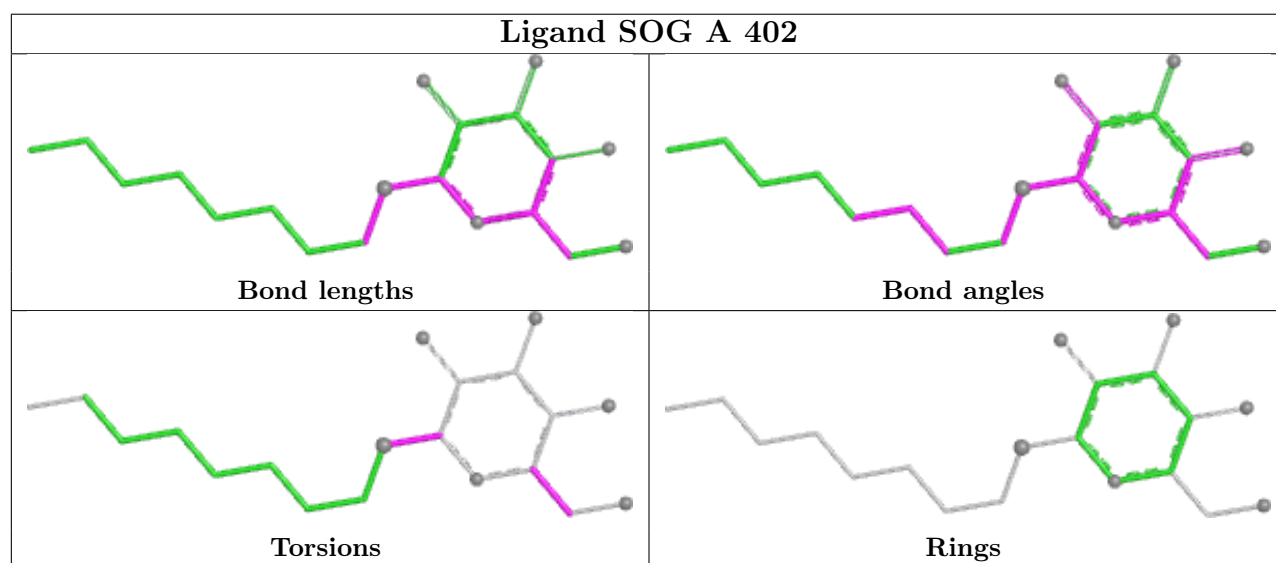
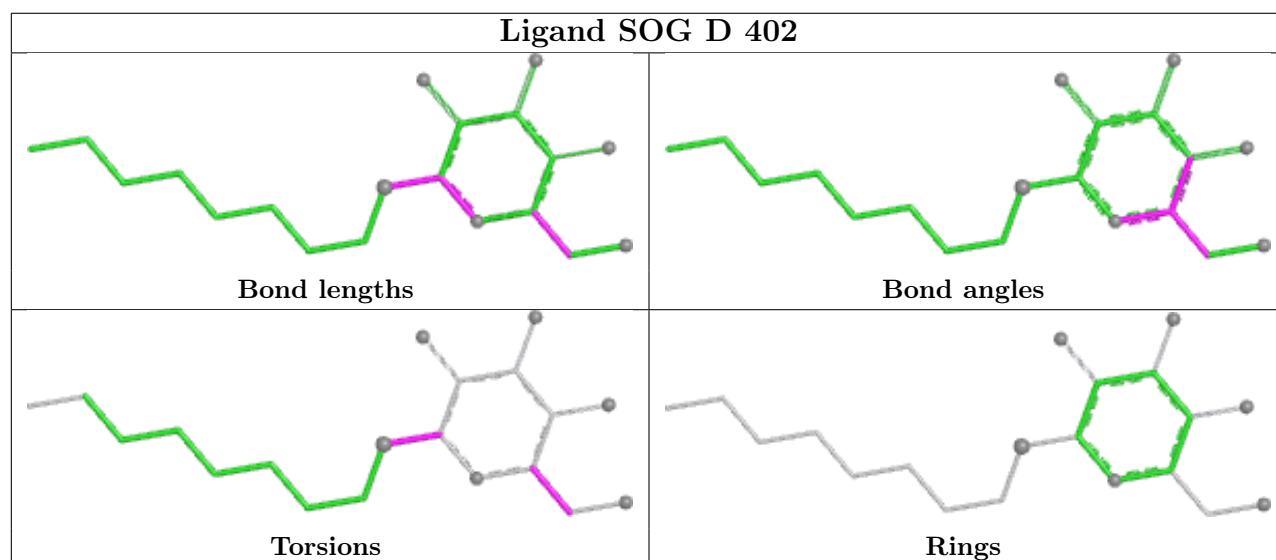
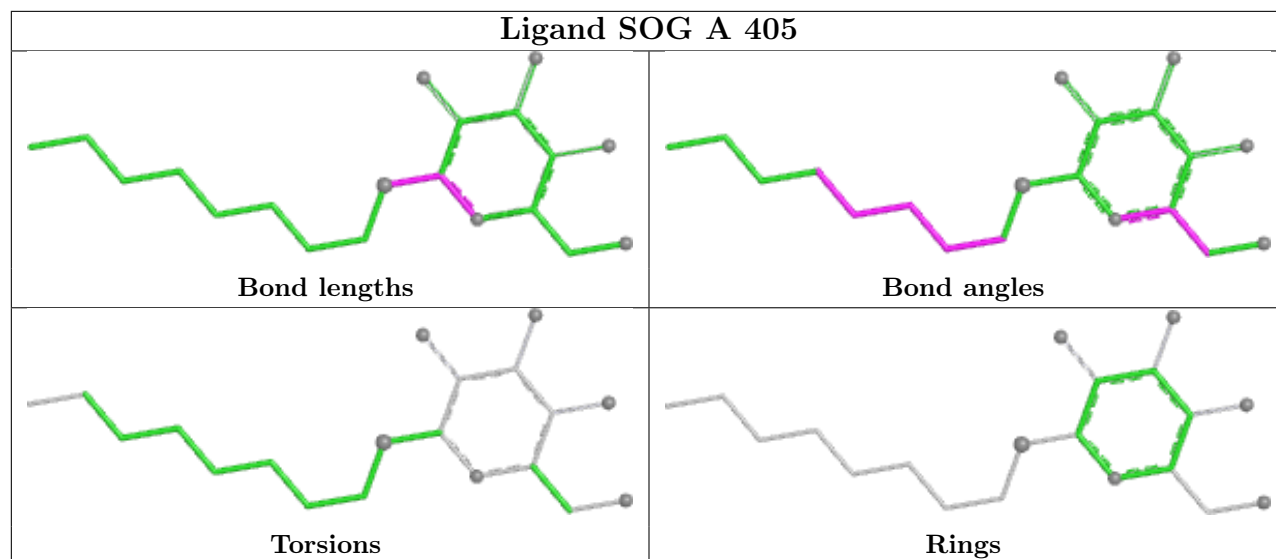
There are no ring outliers.

11 monomers are involved in 36 short contacts:

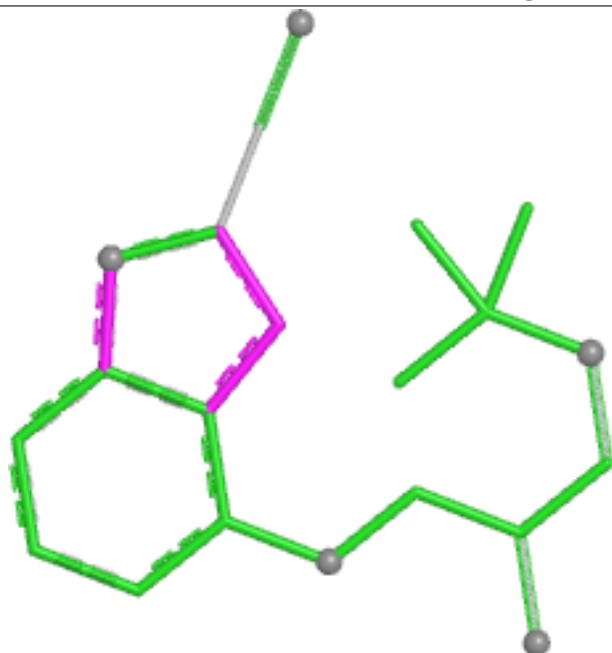
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	D	405	SOG	6	0
4	D	403	SOG	4	0
4	A	405	SOG	5	0
4	A	402	SOG	2	0
2	C	400	P32	1	0
4	B	405	SOG	3	0
2	B	400	P32	2	0
4	C	405	SOG	8	0
4	B	404	SOG	2	0
4	C	404	SOG	2	0
4	C	402	SOG	1	0

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

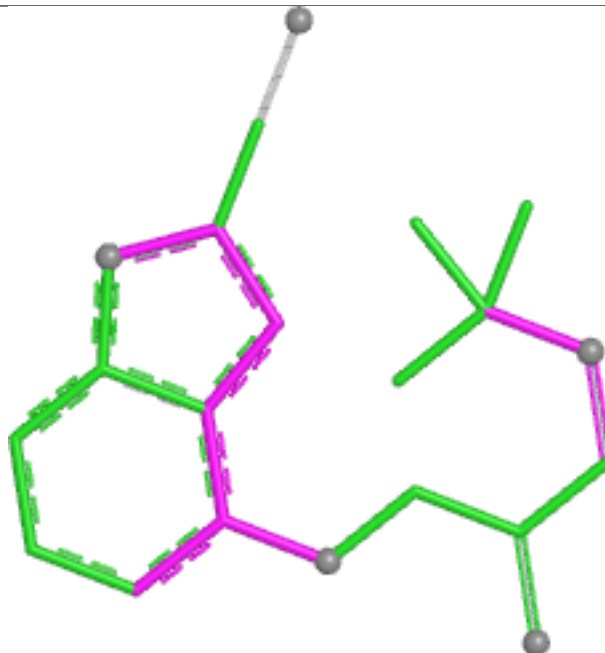




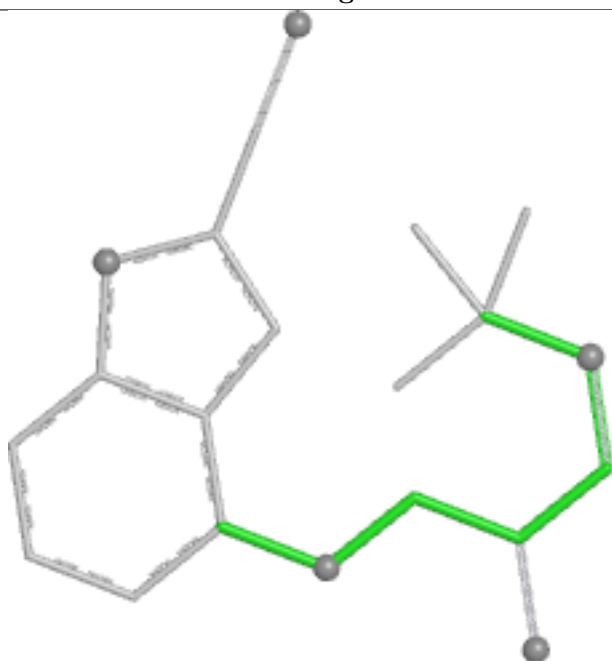
Ligand P32 C 400



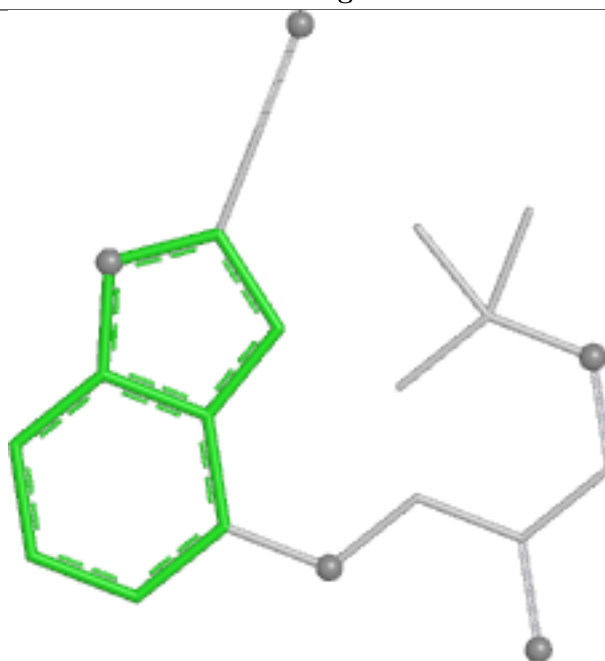
Bond lengths



Bond angles

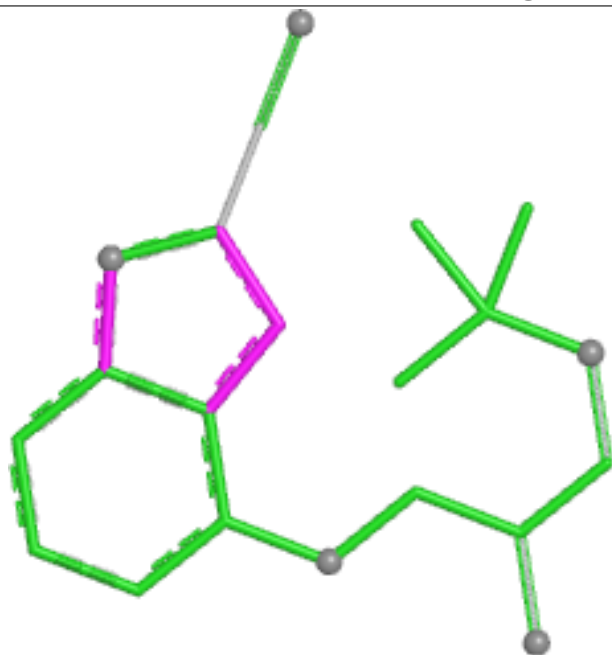


Torsions

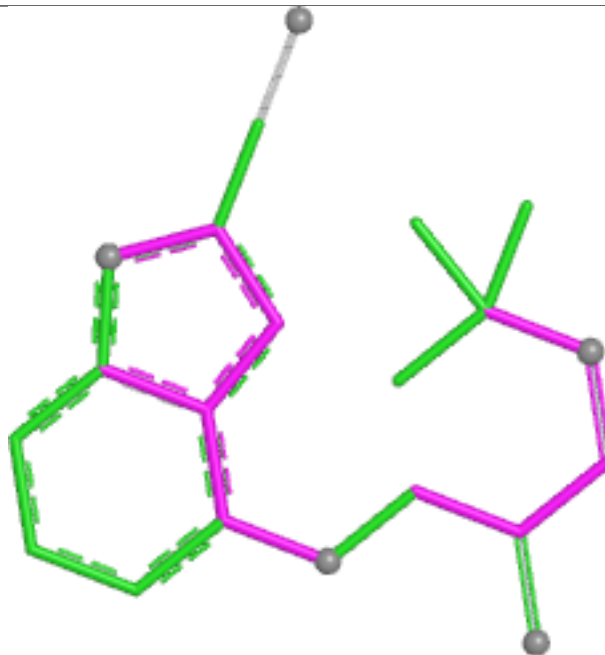


Rings

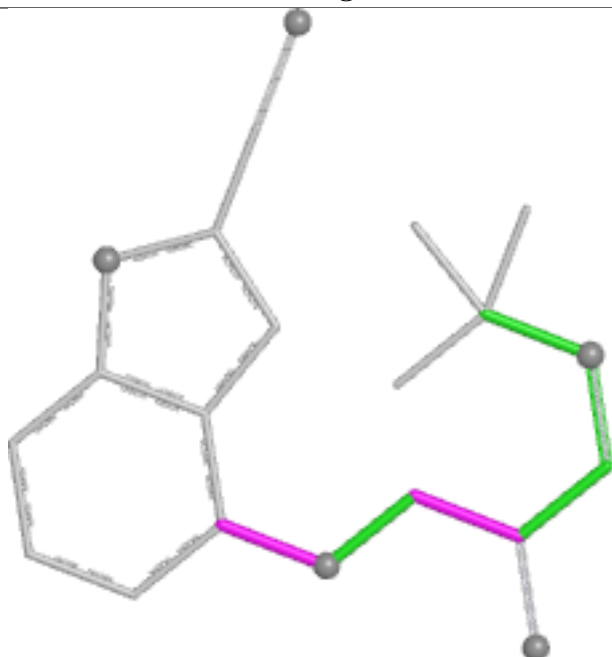
Ligand P32 A 400



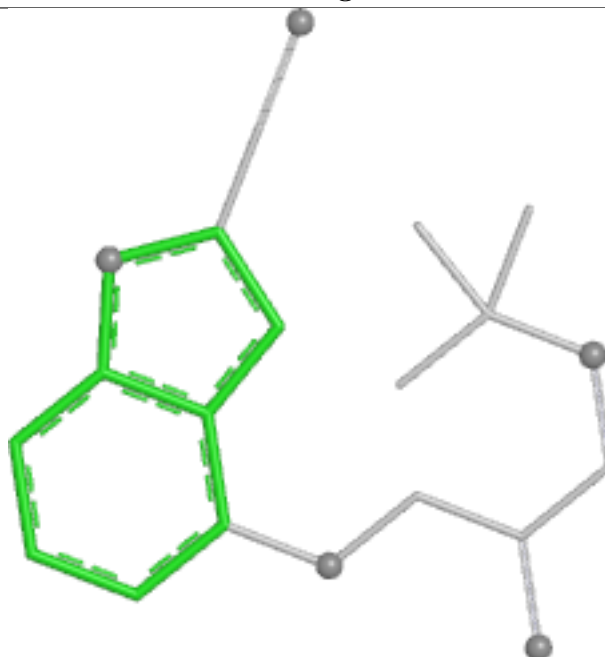
Bond lengths



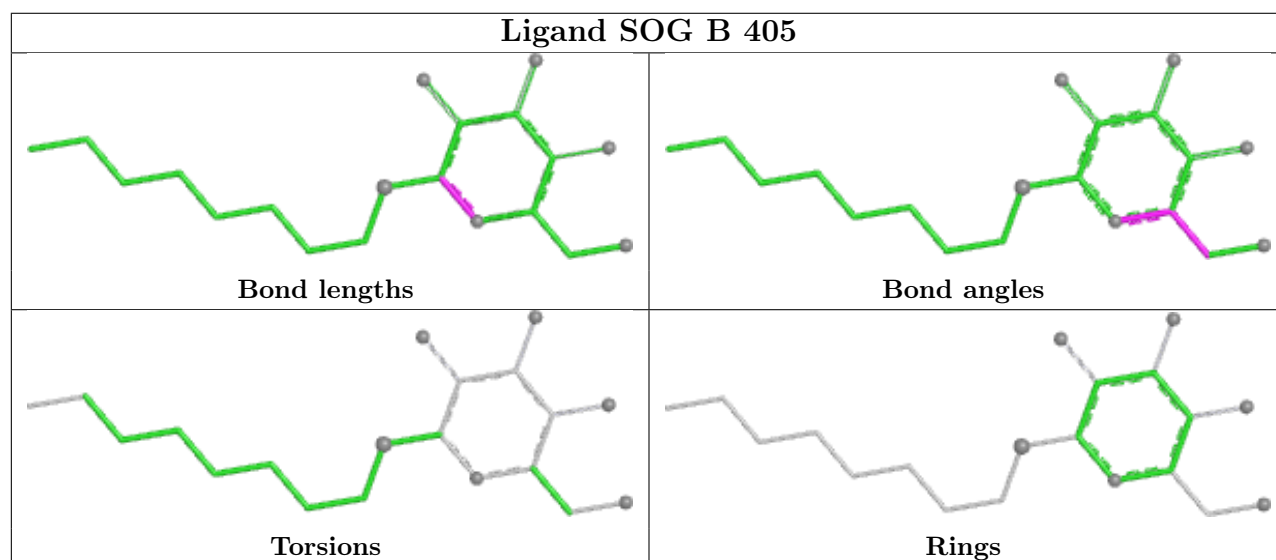
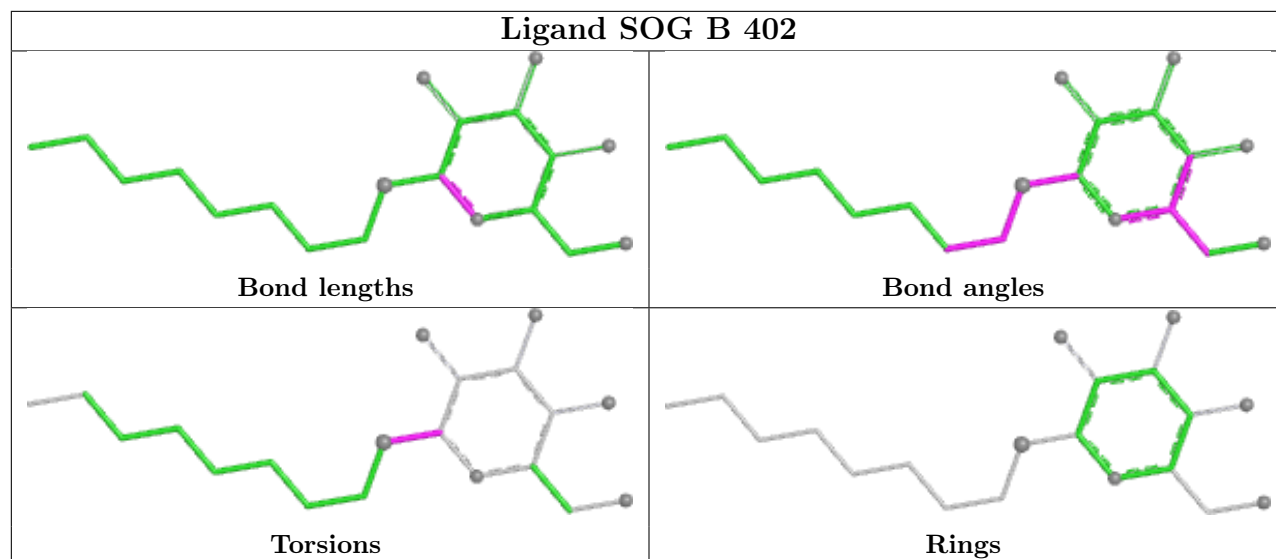
Bond angles



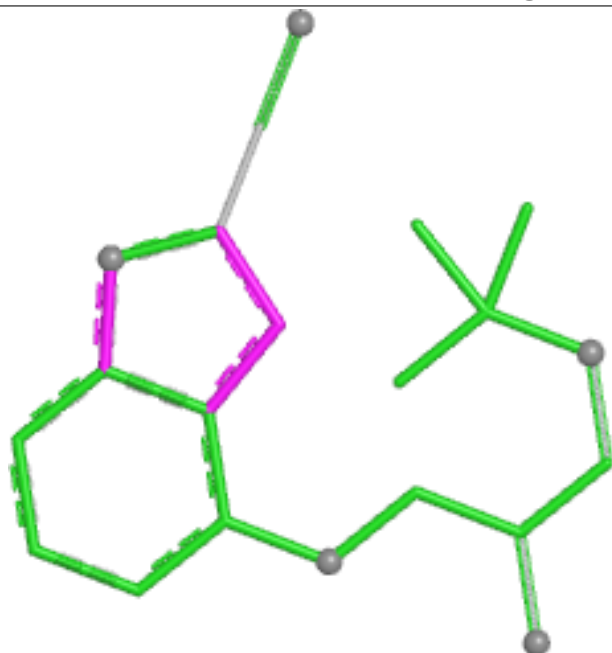
Torsions



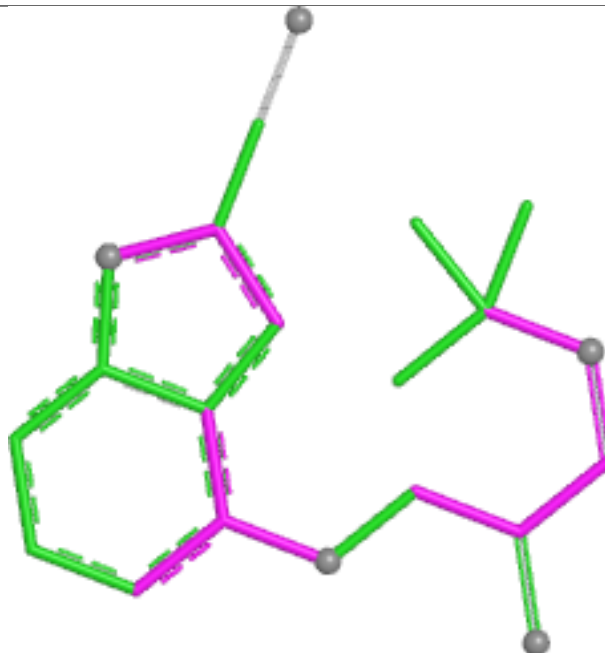
Rings



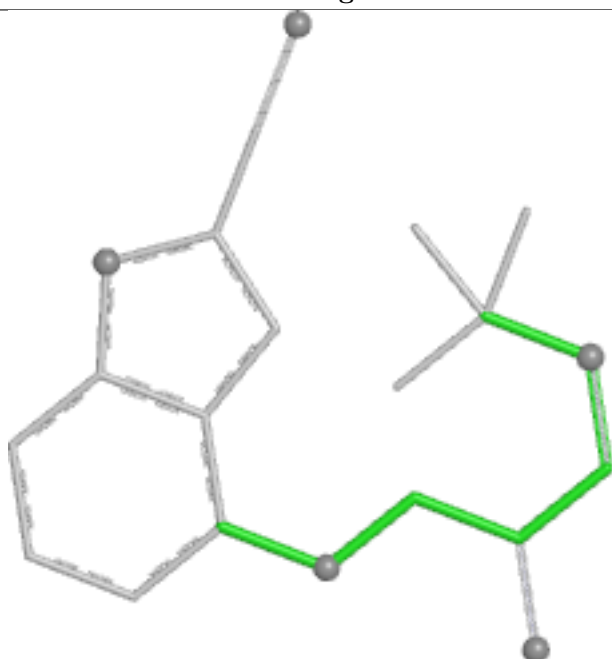
Ligand P32 B 400



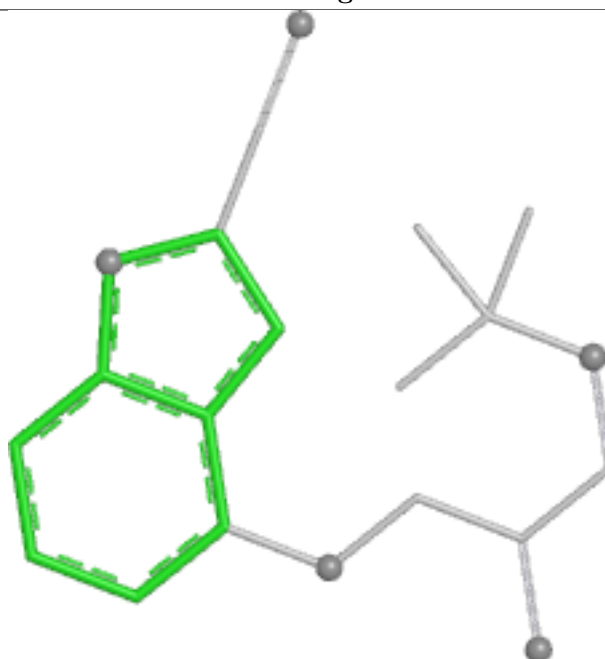
Bond lengths



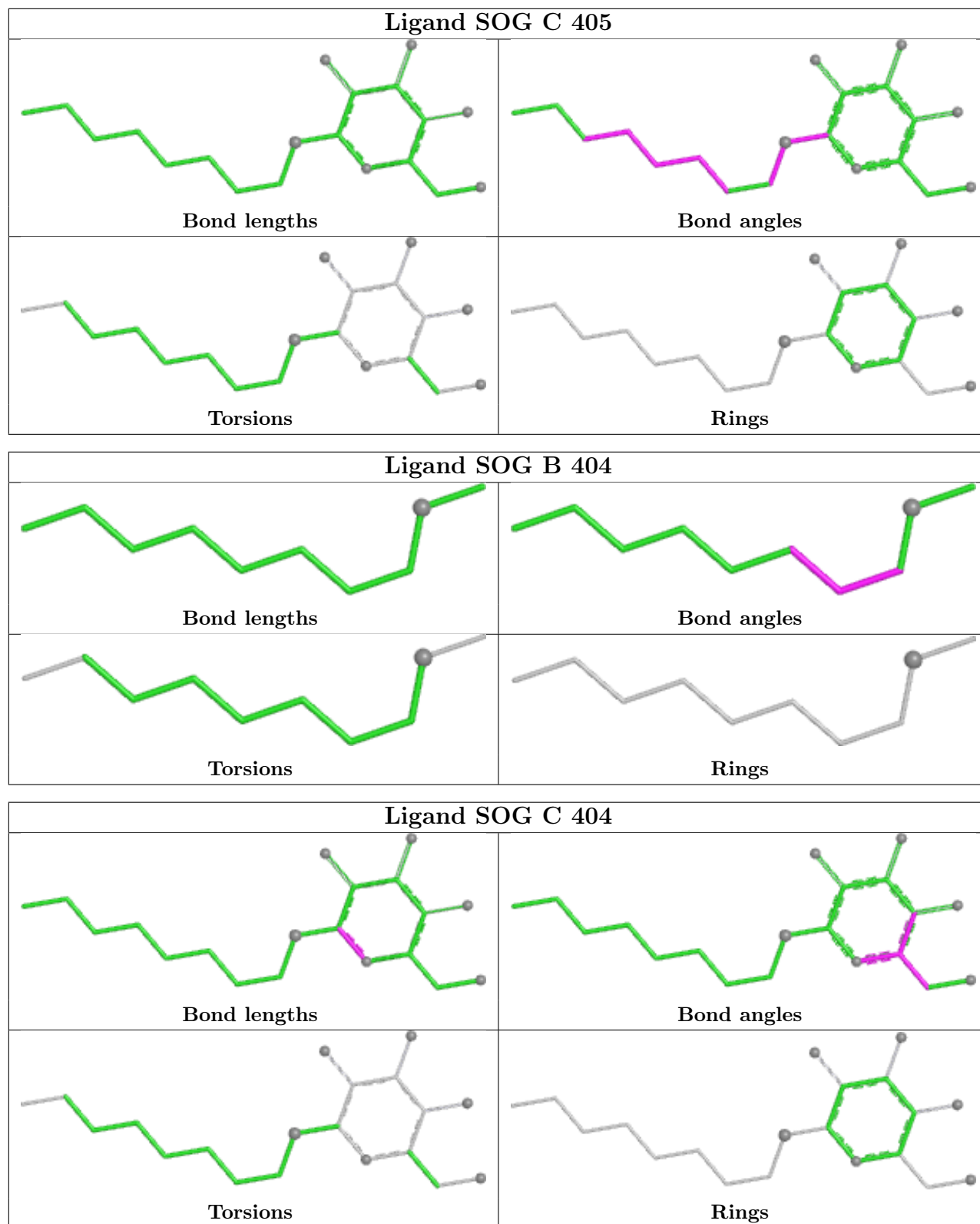
Bond angles



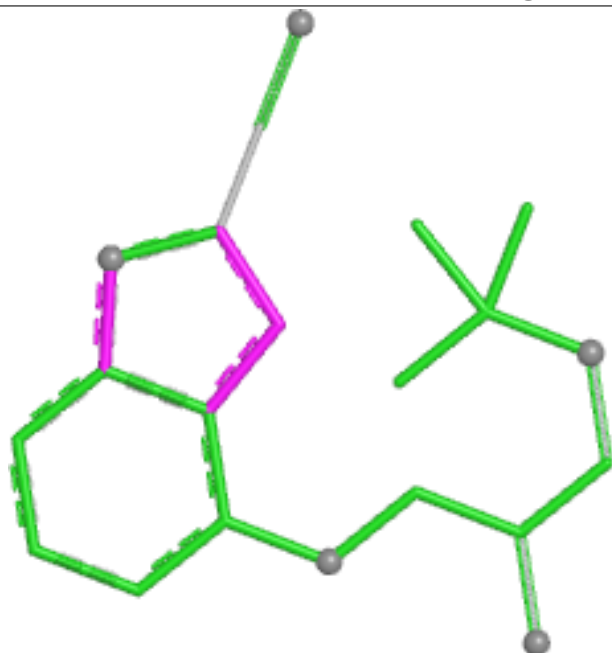
Torsions



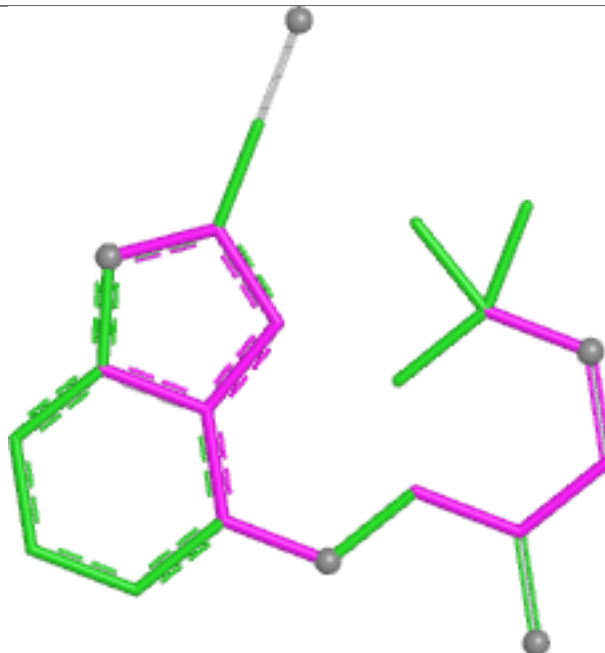
Rings



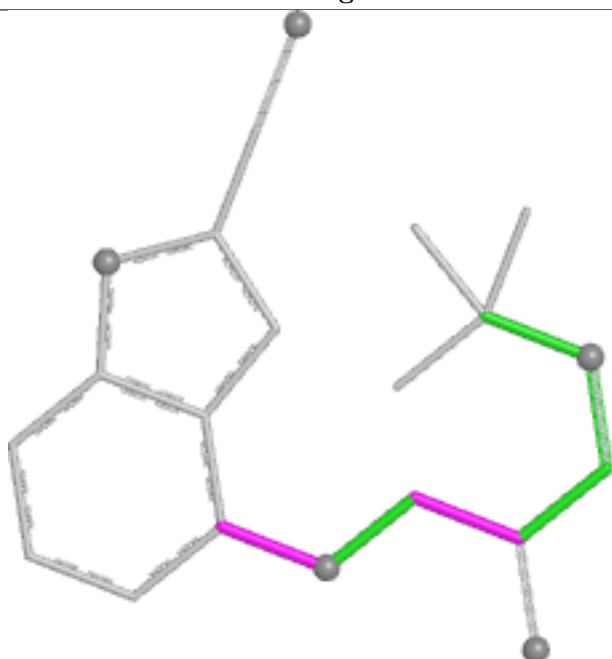
Ligand P32 D 400



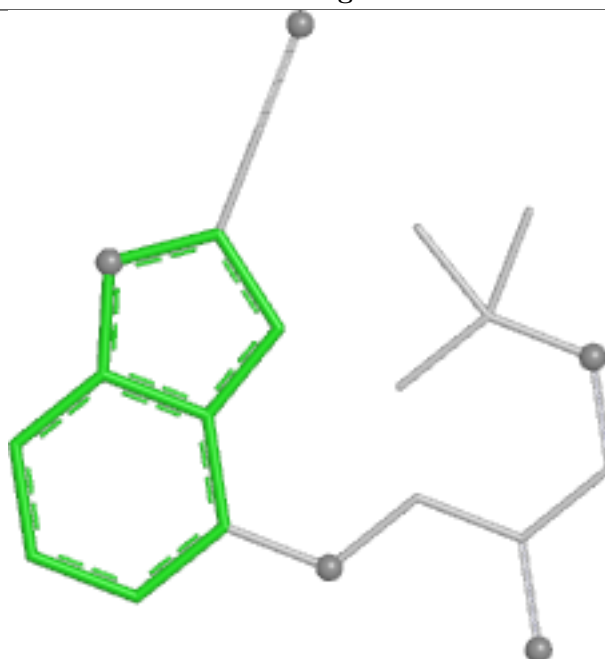
Bond lengths



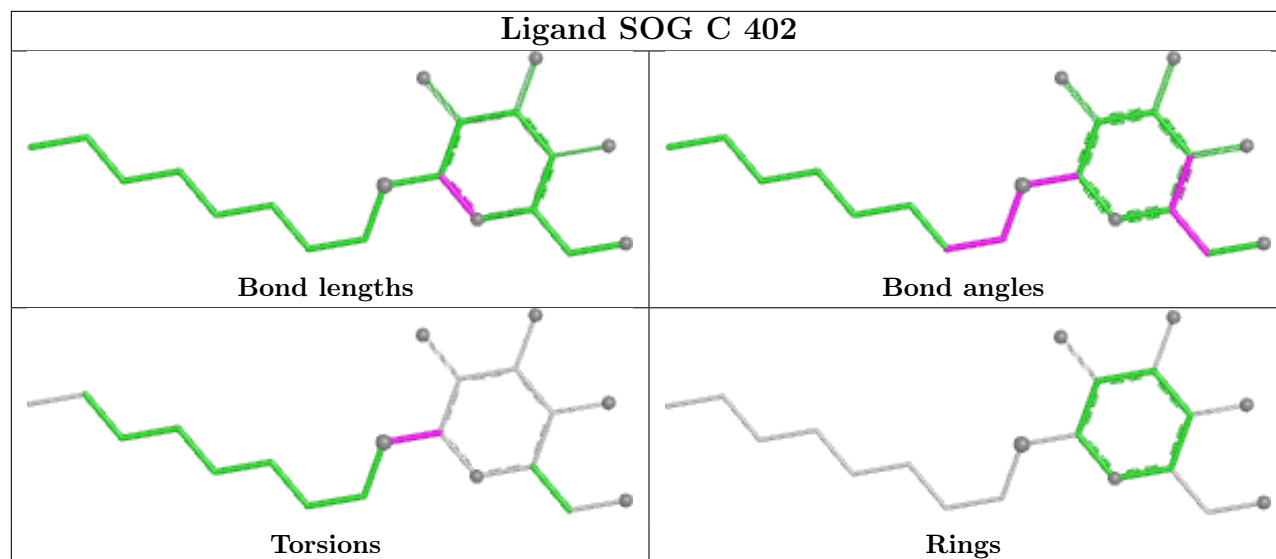
Bond angles



Torsions



Rings



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	274/313 (87%)	-0.02	6 (2%) 62 59	19, 36, 82, 119	0
1	B	276/313 (88%)	0.19	20 (7%) 21 18	21, 41, 104, 138	0
1	C	263/313 (84%)	0.13	13 (4%) 35 31	21, 40, 87, 120	0
1	D	272/313 (86%)	0.01	12 (4%) 39 35	20, 36, 81, 119	0
All	All	1085/1252 (86%)	0.08	51 (4%) 36 33	19, 38, 88, 138	0

All (51) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	353	PHE	5.1
1	B	70	GLN	4.2
1	B	349	PHE	4.1
1	B	352	ALA	4.0
1	D	237	GLN	3.9
1	B	69	THR	3.7
1	A	40	TRP	3.6
1	B	357	LEU	3.5
1	C	344	CYS	3.4
1	C	40	TRP	3.4
1	D	357	LEU	3.4
1	D	238	ILE	3.3
1	A	358	ALA	3.3
1	B	359	PHE	3.2
1	C	284	ARG	3.1
1	B	348	ASP	3.0
1	A	238	ILE	3.0
1	B	344	CYS	3.0
1	D	236	GLU	3.0
1	A	70	GLN	3.0
1	D	356	LEU	2.8

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Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	C	69	THR	2.8
1	C	316	ASN	2.7
1	C	147	PHE	2.7
1	B	72	LEU	2.6
1	C	346	SER	2.6
1	C	47	LEU	2.6
1	B	354	LYS	2.6
1	C	68	SER	2.5
1	C	238	ILE	2.5
1	D	285	GLU	2.5
1	B	147	PHE	2.5
1	B	71	ARG	2.5
1	D	313	ASN	2.4
1	B	47	LEU	2.4
1	B	347	PRO	2.4
1	C	70	GLN	2.4
1	D	284	ARG	2.4
1	B	322	ASP	2.3
1	C	319	LEU	2.3
1	D	69	THR	2.3
1	D	103	VAL	2.2
1	B	346	SER	2.2
1	D	286	HIS	2.2
1	A	235	LYS	2.1
1	D	315	PHE	2.1
1	C	44	MET	2.1
1	B	236	GLU	2.1
1	B	317	ARG	2.1
1	A	103	VAL	2.1
1	B	39	GLN	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 ⓘ

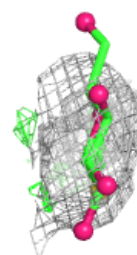
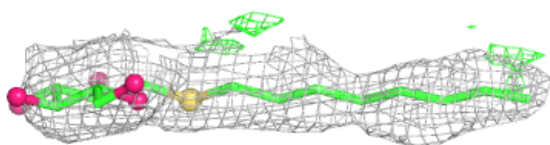
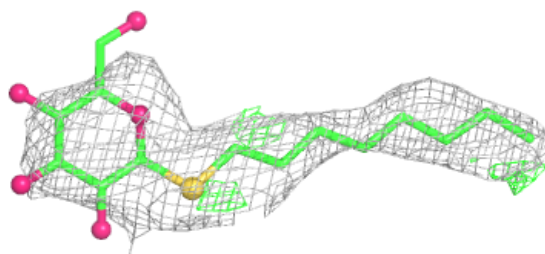
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists 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	SOG	B	405	20/20	0.70	0.19	55,114,128,131	0
4	SOG	D	405	20/20	0.73	0.18	41,115,131,133	0
4	SOG	C	405	20/20	0.77	0.17	55,107,126,128	0
4	SOG	A	405	20/20	0.83	0.17	39,89,110,116	0
5	D10	C	406	10/10	0.83	0.18	36,49,60,60	0
4	SOG	C	404	20/20	0.85	0.17	29,95,123,123	0
4	SOG	B	404	10/20	0.85	0.15	30,42,48,81	0
5	D10	A	406	10/10	0.86	0.20	25,41,47,48	0
4	SOG	C	402	20/20	0.86	0.14	18,91,111,113	0
4	SOG	A	402	20/20	0.87	0.16	33,54,74,94	0
4	SOG	B	402	20/20	0.88	0.14	19,90,122,125	0
4	SOG	D	402	20/20	0.89	0.15	26,54,88,90	0
4	SOG	D	403	20/20	0.90	0.12	18,36,54,72	0
3	NA	B	401	1/1	0.93	0.08	40,40,40,40	0
4	SOG	A	403	20/20	0.93	0.10	28,39,49,50	0
2	P32	A	400	21/21	0.94	0.09	17,32,41,44	0
2	P32	C	400	21/21	0.96	0.08	22,33,43,51	0
2	P32	B	400	21/21	0.96	0.07	19,33,39,50	0
3	NA	D	401	1/1	0.97	0.05	39,39,39,39	0
2	P32	D	400	21/21	0.97	0.07	17,33,40,42	0
3	NA	C	401	1/1	0.97	0.04	36,36,36,36	0
3	NA	A	401	1/1	0.98	0.05	33,33,33,33	0

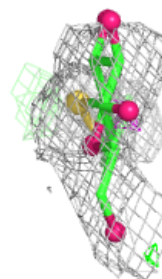
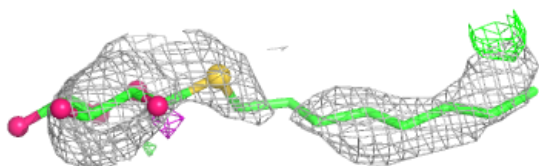
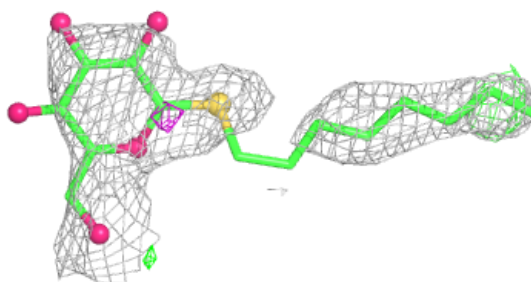
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around SOG B 405:

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

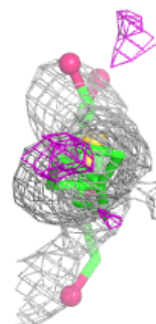
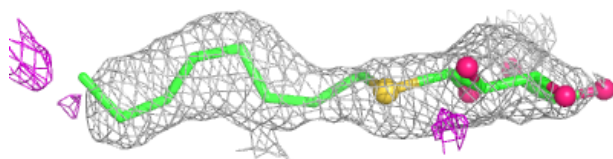
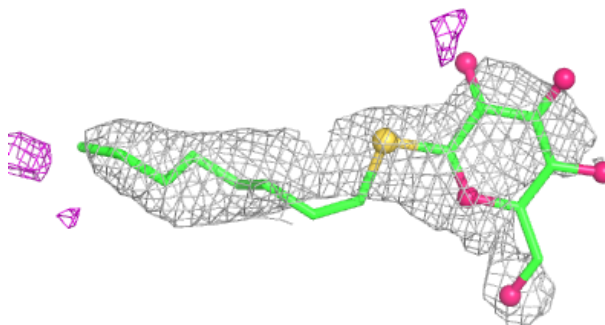
**Electron density around SOG C 405:**

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

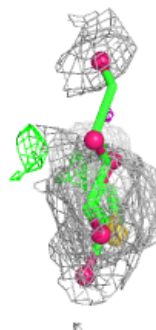
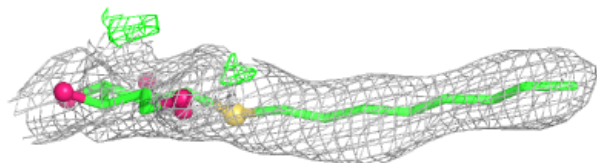
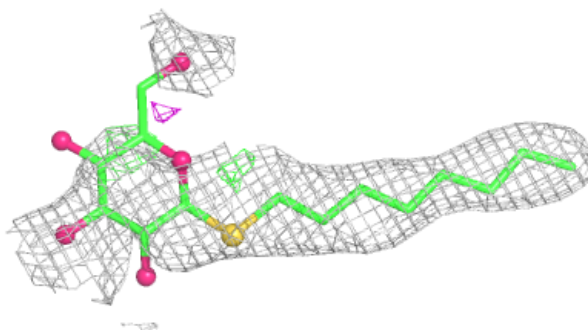


Electron density around SOG A 405:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

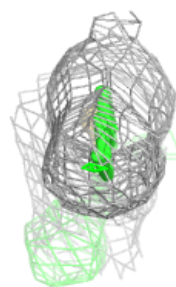
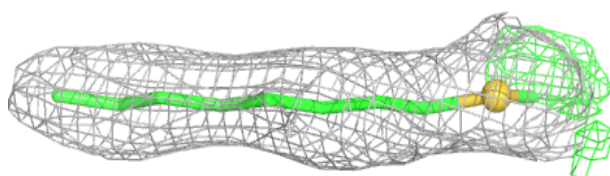
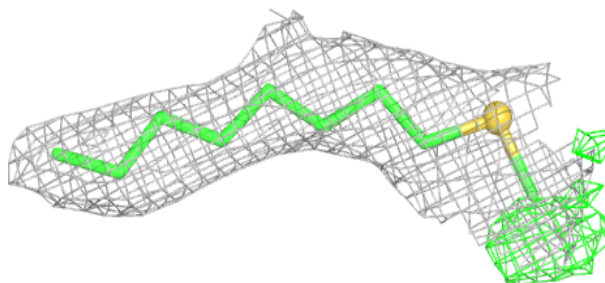
**Electron density around SOG C 404:**

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

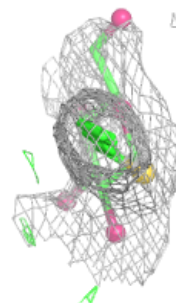
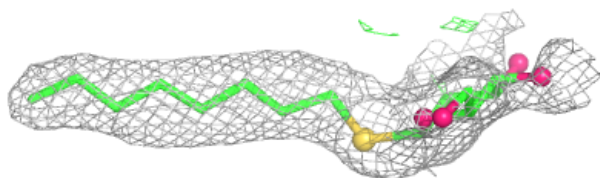
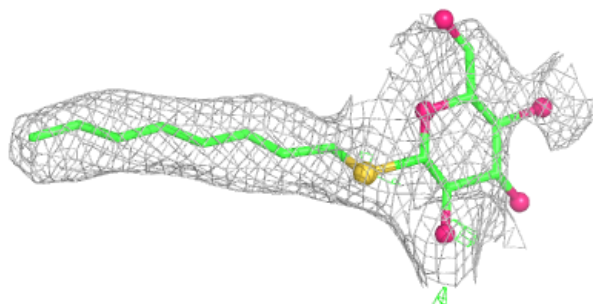


Electron density around SOG B 404:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

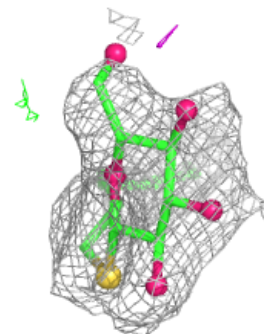
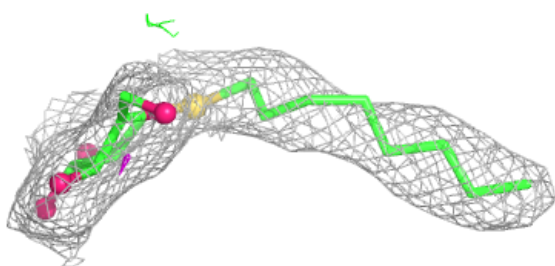
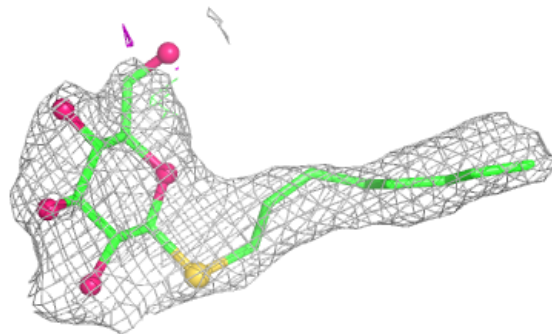
**Electron density around SOG C 402:**

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and green (positive)

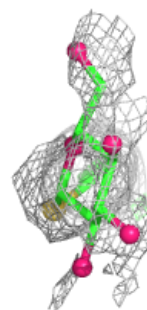
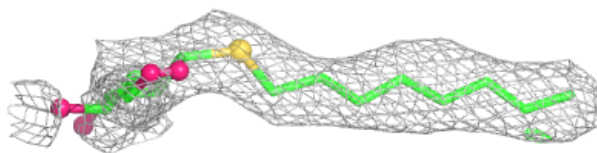
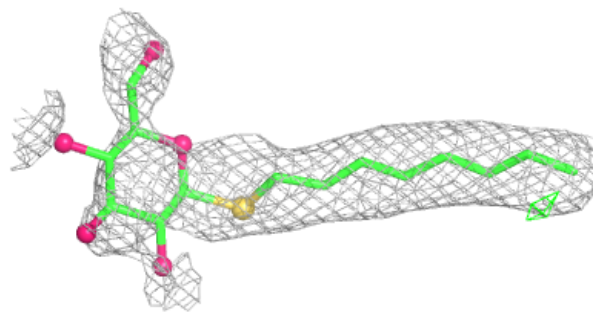


Electron density around SOG A 402:

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

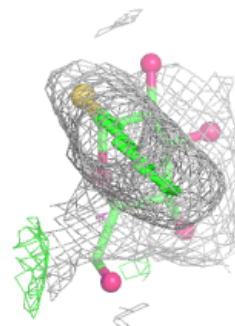
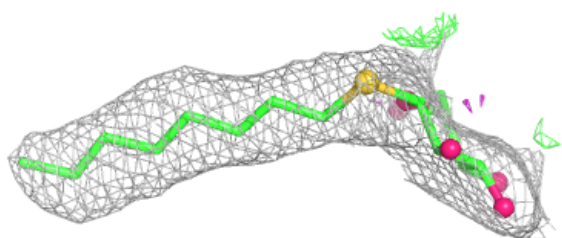
**Electron density around SOG B 402:**

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and green (positive)

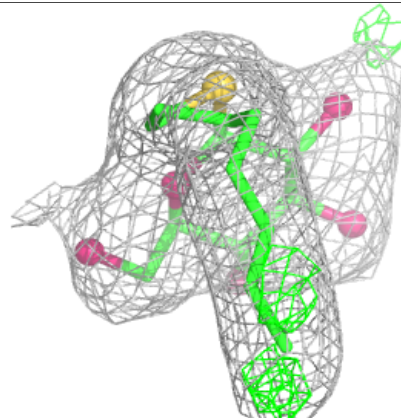
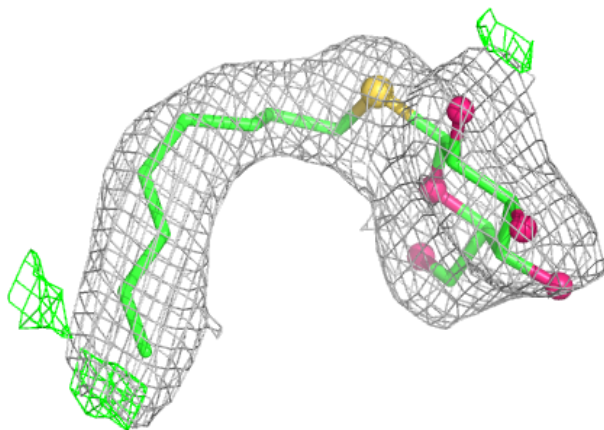


Electron density around SOG D 402:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

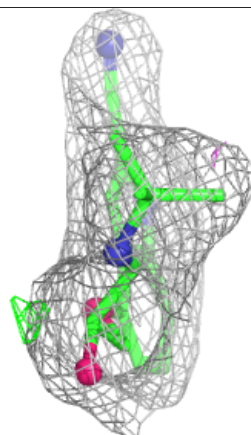
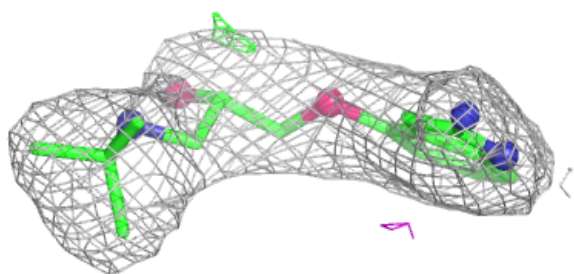
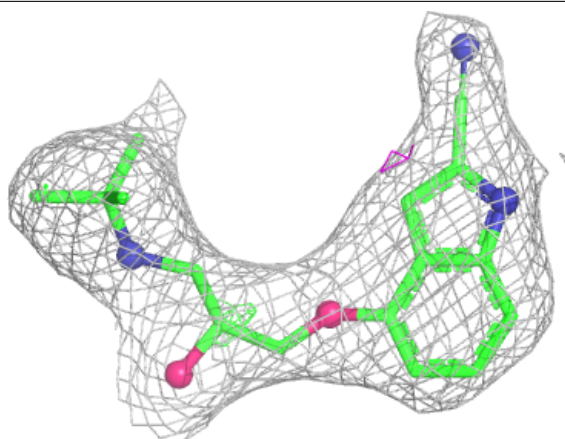
**Electron density around SOG D 403:**

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



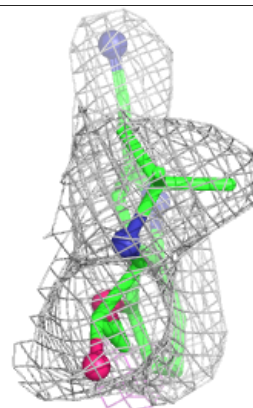
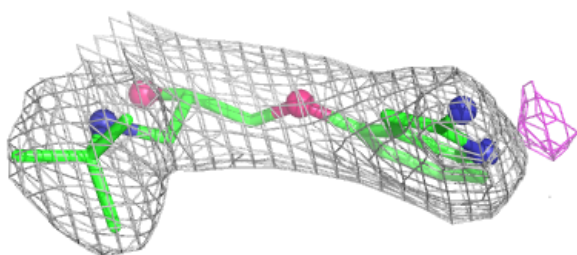
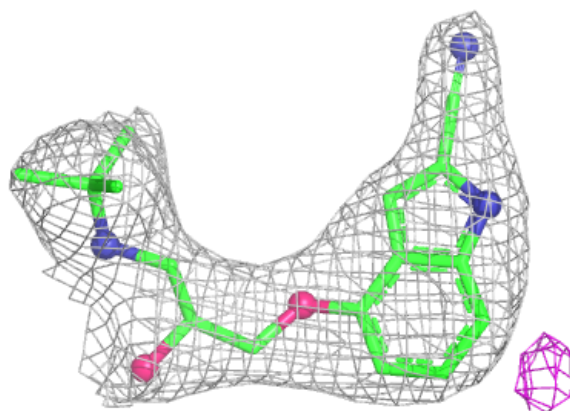
Electron density around P32 A 400:

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



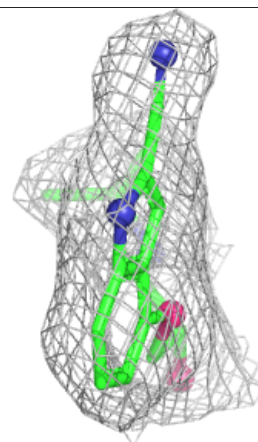
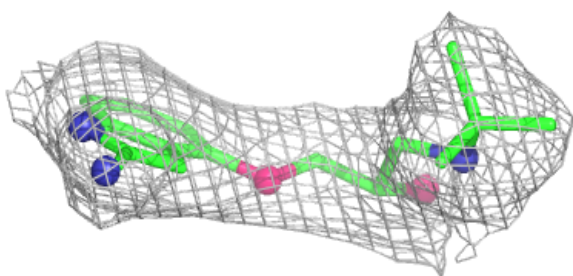
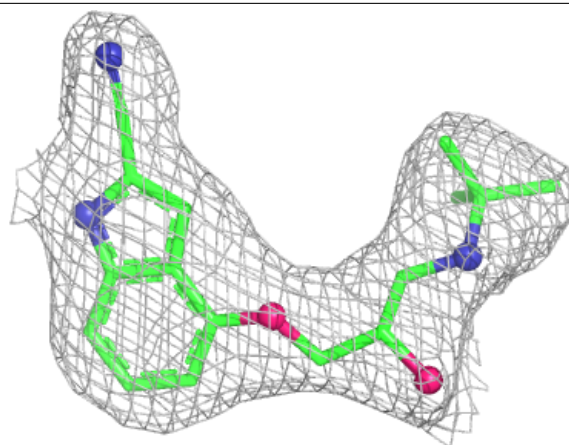
Electron density around P32 C 400:

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



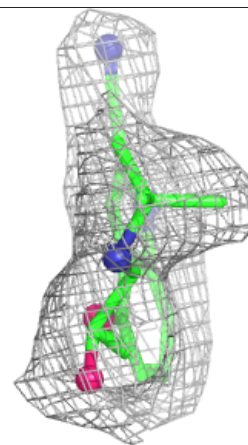
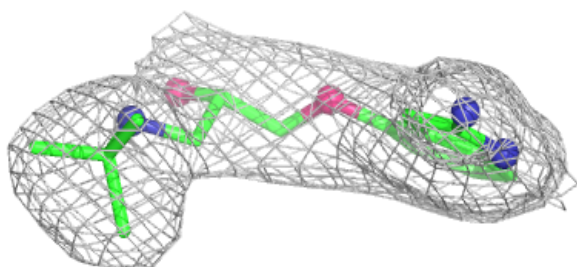
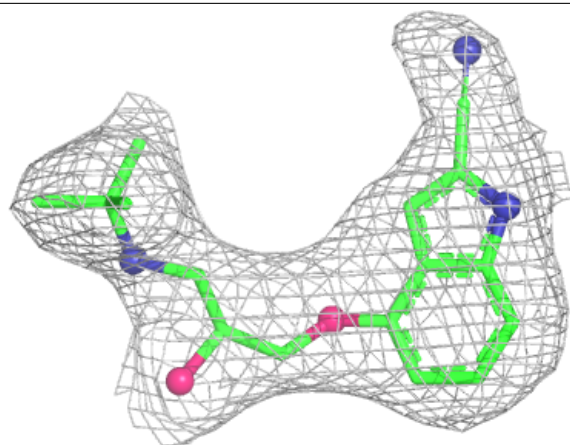
Electron density around P32 B 400:

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



Electron density around P32 D 400:

$2mF_o-DF_c$ (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.