



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2026 – 12:14 AM UTC

PDB ID : 8FYV / pdb_00008fyv
Title : Salmonella enterica serovar Typhimurium chemoreceptor Tsr (taxis to serine and repellents) ligand-binding domain in complex with l-serine
Authors : Baylink, A.; Gentry-Lear, Z.; Glenn, S.
Deposited on : 2023-01-26
Resolution : 2.20 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

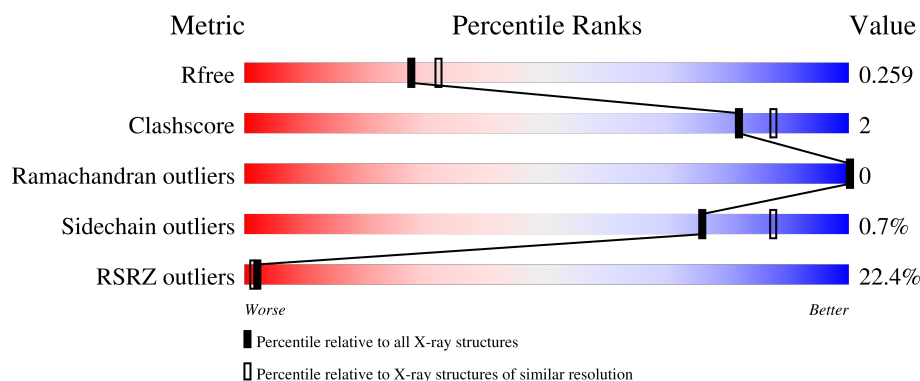
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 Å.

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



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

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

Mol	Chain	Length	Quality of chain
1	A	156	22% 85% 6% 10%
1	B	156	22% 85% 5% 10%
1	C	156	21% 83% 6% 10%
1	D	156	21% 85% . . 10%
1	E	156	16% 83% 6% 11%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 11197 atoms, of which 5300 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.

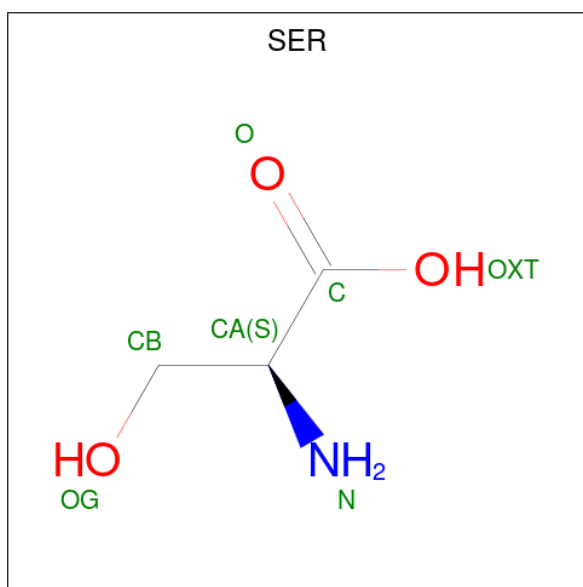
- Molecule 1 is a protein called Methyl-accepting chemotaxis protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	141	Total	C	H	N	O	S	0	2	0
			2213	714	1080	193	221	5			
1	B	140	Total	C	H	N	O	S	0	2	0
			2183	707	1061	192	218	5			
1	C	140	Total	C	H	N	O	S	0	4	0
			2201	708	1074	193	220	6			
1	D	140	Total	C	H	N	O	S	0	0	0
			2127	687	1035	188	212	5			
1	E	139	Total	C	H	N	O	S	0	0	0
			2126	688	1030	189	214	5			

- Molecule 2 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Cl	0	0
			2	2		
2	B	2	Total	Cl	0	0
			2	2		
2	C	2	Total	Cl	0	0
			2	2		
2	D	2	Total	Cl	0	0
			2	2		
2	E	1	Total	Cl	0	0
			1	1		

- Molecule 3 is SERINE (CCD ID: SER) (formula: C₃H₇NO₃) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	H	N	O	0	0
			11	3	4	1	3		
3	B	1	Total	C	H	N	O	0	0
			11	3	4	1	3		
3	C	1	Total	C	H	N	O	0	0
			11	3	4	1	3		
3	D	1	Total	C	H	N	O	0	0
			11	3	4	1	3		
3	E	1	Total	C	H	N	O	0	0
			11	3	4	1	3		

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	4	Total	Na	0	0
			4	4		
4	C	1	Total	Na	0	0
			1	1		
4	D	1	Total	Na	0	0
			1	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	52	Total	O	0	1
			53	53		

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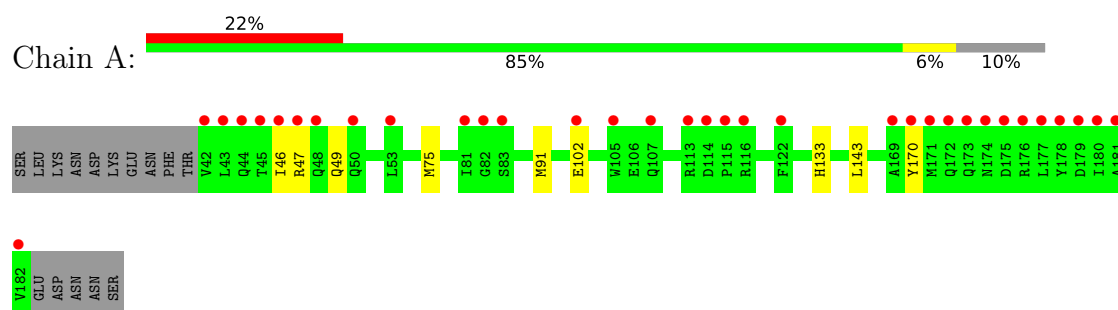
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	61	Total 61	O 61	0	0
5	C	63	Total 63	O 63	0	0
5	D	57	Total 57	O 57	0	0
5	E	42	Total 43	O 43	0	1

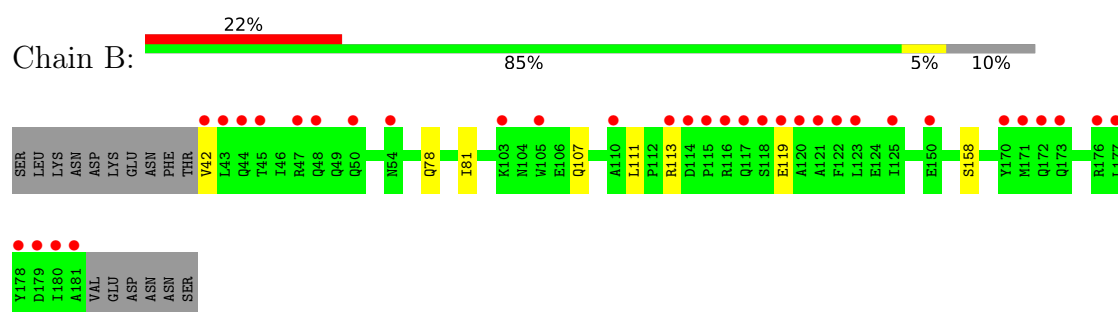
3 Residue-property plots [i](#)

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.

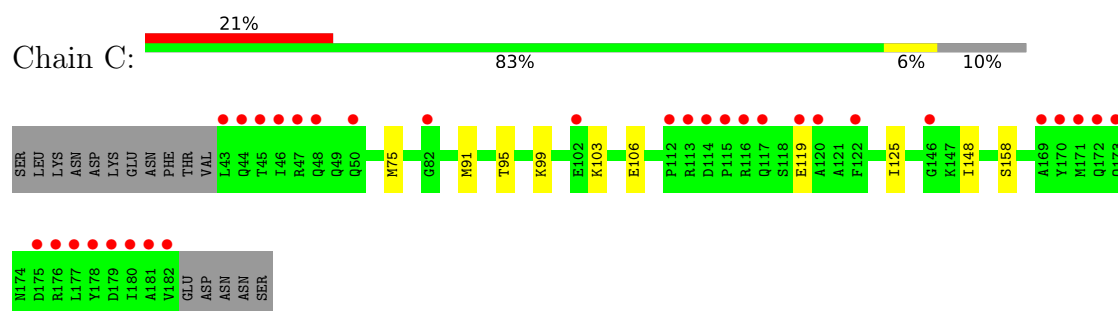
- Molecule 1: Methyl-accepting chemotaxis protein



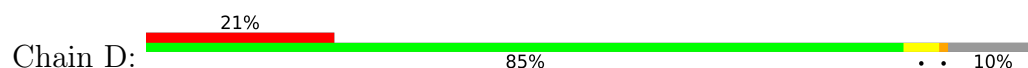
- Molecule 1: Methyl-accepting chemotaxis protein

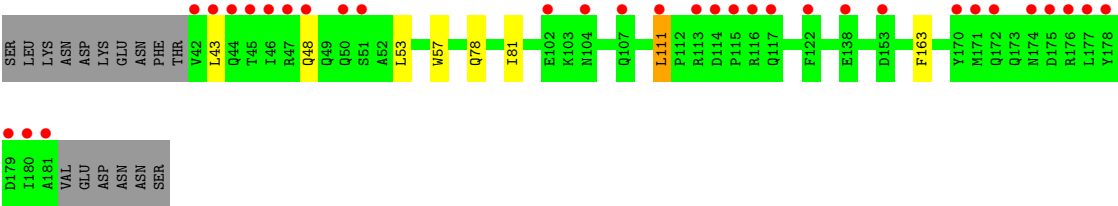


- Molecule 1: Methyl-accepting chemotaxis protein

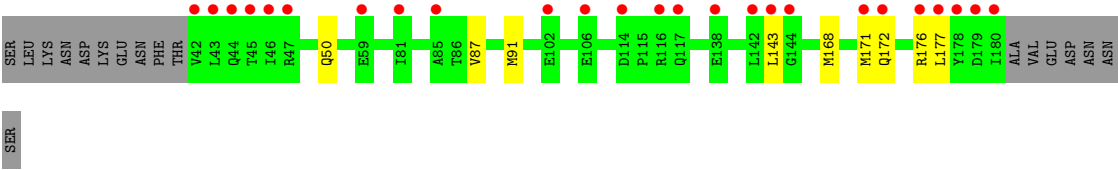
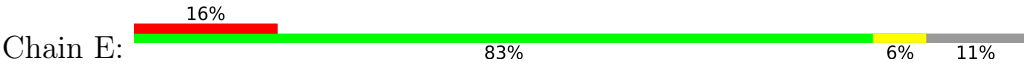


- Molecule 1: Methyl-accepting chemotaxis protein





● Molecule 1: Methyl-accepting chemotaxis protein



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	125.81Å 75.23Å 129.59Å 90.00° 116.47° 90.00°	Depositor
Resolution (Å)	58.00 – 2.20 58.00 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.8 (58.00-2.20) 100.0 (58.00-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.41 (at 2.20Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.241 , 0.258 0.242 , 0.259	Depositor DCC
R_{free} test set	2002 reflections (3.63%)	wwPDB-VP
Wilson B-factor (Å ²)	26.1	Xtriage
Anisotropy	0.844	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 51.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.003 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	11197	wwPDB-VP
Average B, all atoms (Å ²)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.27% 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: NA, CL

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.09	0/1162	0.23	0/1575
1	B	0.08	0/1151	0.21	0/1560
1	C	0.08	0/1164	0.22	0/1576
1	D	0.09	0/1111	0.24	0/1507
1	E	0.08	0/1116	0.22	0/1514
All	All	0.08	0/5704	0.23	0/7732

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	1133	1080	1068	6	0
1	B	1122	1061	1055	7	0
1	C	1127	1074	1053	9	0
1	D	1092	1035	1035	5	0
1	E	1096	1030	1030	7	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	2	0	0	0	0
2	E	1	0	0	0	0
3	A	7	4	4	0	0
3	B	7	4	4	0	0
3	C	7	4	4	0	0
3	D	7	4	4	0	0
3	E	7	4	4	0	0
4	B	4	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
5	A	53	0	0	0	0
5	B	61	0	0	2	0
5	C	63	0	0	1	0
5	D	57	0	0	0	0
5	E	43	0	0	0	0
All	All	5897	5300	5261	27	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 2.

All (27) 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:A:75:MET:HE1	1:B:81:ILE:HD11	1.71	0.72
1:C:99:LYS:NZ	5:C:301:HOH:O	2.32	0.63
1:B:107:GLN:O	1:B:111:LEU:HD13	2.00	0.61
1:C:106:GLU:OE1	1:E:176:ARG:NH2	2.36	0.58
1:C:75:MET:HE3	1:C:148:ILE:HG12	1.87	0.57
1:B:78:GLN:NE2	5:B:305:HOH:O	2.39	0.55
1:A:75:MET:CE	1:B:81:ILE:HD11	2.36	0.54
1:A:91:MET:SD	1:A:143:LEU:HD23	2.48	0.54
1:C:103:LYS:HG3	1:E:176:ARG:HD2	1.90	0.54
1:C:91:MET:O	1:C:95:THR:HG23	2.09	0.53
1:B:42:VAL:N	5:B:308:HOH:O	2.42	0.52
1:E:91:MET:SD	1:E:143:LEU:HD23	2.49	0.52
1:A:46:ILE:HD12	1:A:47:ARG:N	2.26	0.51
1:E:50:GLN:NE2	1:E:171:MET:SD	2.85	0.49
1:B:158:SER:HB2	1:C:158:SER:HB2	1.95	0.49
1:D:48:GLN:HG2	1:D:111:LEU:HD11	1.95	0.48
1:C:75:MET:HE1	1:D:81:ILE:HD11	1.95	0.48
1:A:102:GLU:HG3	1:A:133:HIS:CE1	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:119:GLU:HA	1:C:119:GLU:OE1	2.17	0.45
1:D:78:GLN:HA	1:D:78:GLN:OE1	2.18	0.44
1:E:168:MET:CE	1:E:171:MET:HE2	2.49	0.43
1:B:113:ARG:NH1	1:B:119:GLU:OE2	2.52	0.43
1:D:57:TRP:CE3	1:D:163:PHE:CE2	3.07	0.43
1:C:103:LYS:HE2	1:E:172:GLN:HB3	2.00	0.42
1:D:53:LEU:HD12	1:D:53:LEU:O	2.19	0.42
1:E:87:VAL:O	1:E:91:MET:HG2	2.19	0.42
1:A:49:GLN:HG2	1:A:170:TYR:OH	2.21	0.41

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	141/156 (90%)	140 (99%)	1 (1%)	0	100	100
1	B	140/156 (90%)	139 (99%)	1 (1%)	0	100	100
1	C	142/156 (91%)	140 (99%)	2 (1%)	0	100	100
1	D	138/156 (88%)	137 (99%)	1 (1%)	0	100	100
1	E	137/156 (88%)	136 (99%)	1 (1%)	0	100	100
All	All	698/780 (90%)	692 (99%)	6 (1%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	116/134 (87%)	116 (100%)	0	100	100
1	B	114/134 (85%)	114 (100%)	0	100	100
1	C	116/134 (87%)	115 (99%)	1 (1%)	70	84
1	D	108/134 (81%)	106 (98%)	2 (2%)	50	66
1	E	109/134 (81%)	108 (99%)	1 (1%)	70	84
All	All	563/670 (84%)	559 (99%)	4 (1%)	76	87

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	125	ILE
1	D	43	LEU
1	D	111	LEU
1	E	177	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	50	GLN
1	A	78	GLN
1	A	96	ASN
1	A	104	ASN
1	A	133	HIS
1	A	173	GLN
1	B	48	GLN
1	B	49	GLN
1	B	54	ASN
1	B	166	GLN
1	B	172	GLN
1	B	173	GLN
1	C	50	GLN
1	C	92	GLN
1	C	96	ASN
1	C	117	GLN
1	D	49	GLN
1	D	50	GLN
1	D	96	ASN

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Mol	Chain	Res	Type
1	D	117	GLN
1	D	173	GLN
1	E	78	GLN
1	E	173	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 20 ligands modelled in this entry, 15 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	$\# Z > 2$
3	SER	A	203	-	4,6,6	1.14	1 (25%)	2,7,7	1.89	1 (50%)
3	SER	D	204	-	4,6,6	1.13	1 (25%)	2,7,7	1.81	1 (50%)
3	SER	E	202	-	4,6,6	1.15	1 (25%)	2,7,7	1.87	1 (50%)
3	SER	B	207	-	4,6,6	1.11	1 (25%)	2,7,7	1.82	1 (50%)
3	SER	C	204	-	4,6,6	1.15	1 (25%)	2,7,7	1.85	1 (50%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns.

'-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	SER	A	203	-	-	0/6/6/6	-
3	SER	D	204	-	-	0/6/6/6	-
3	SER	E	202	-	-	0/6/6/6	-
3	SER	B	207	-	-	0/6/6/6	-
3	SER	C	204	-	-	1/6/6/6	-

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	202	SER	OXT-C	-2.19	1.23	1.30
3	C	204	SER	OXT-C	-2.17	1.23	1.30
3	A	203	SER	OXT-C	-2.14	1.23	1.30
3	D	204	SER	OXT-C	-2.13	1.23	1.30
3	B	207	SER	OXT-C	-2.09	1.24	1.30

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	203	SER	OXT-C-O	-2.64	118.09	124.08
3	E	202	SER	OXT-C-O	-2.63	118.12	124.08
3	C	204	SER	OXT-C-O	-2.59	118.21	124.08
3	B	207	SER	OXT-C-O	-2.54	118.31	124.08
3	D	204	SER	OXT-C-O	-2.54	118.32	124.08

There are no chirality outliers.

All (1) torsion outliers are listed below:

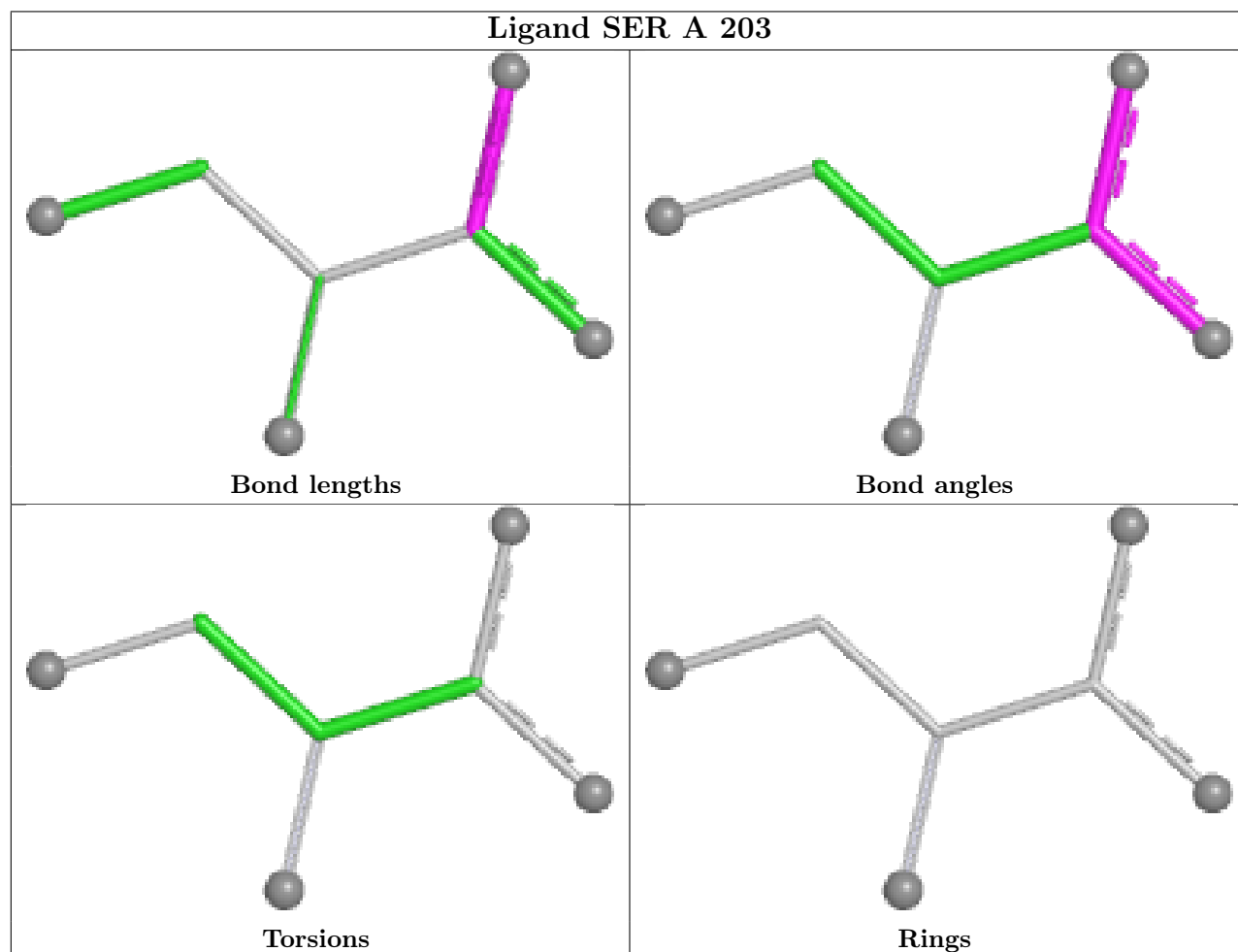
Mol	Chain	Res	Type	Atoms
3	C	204	SER	O-C-CA-N

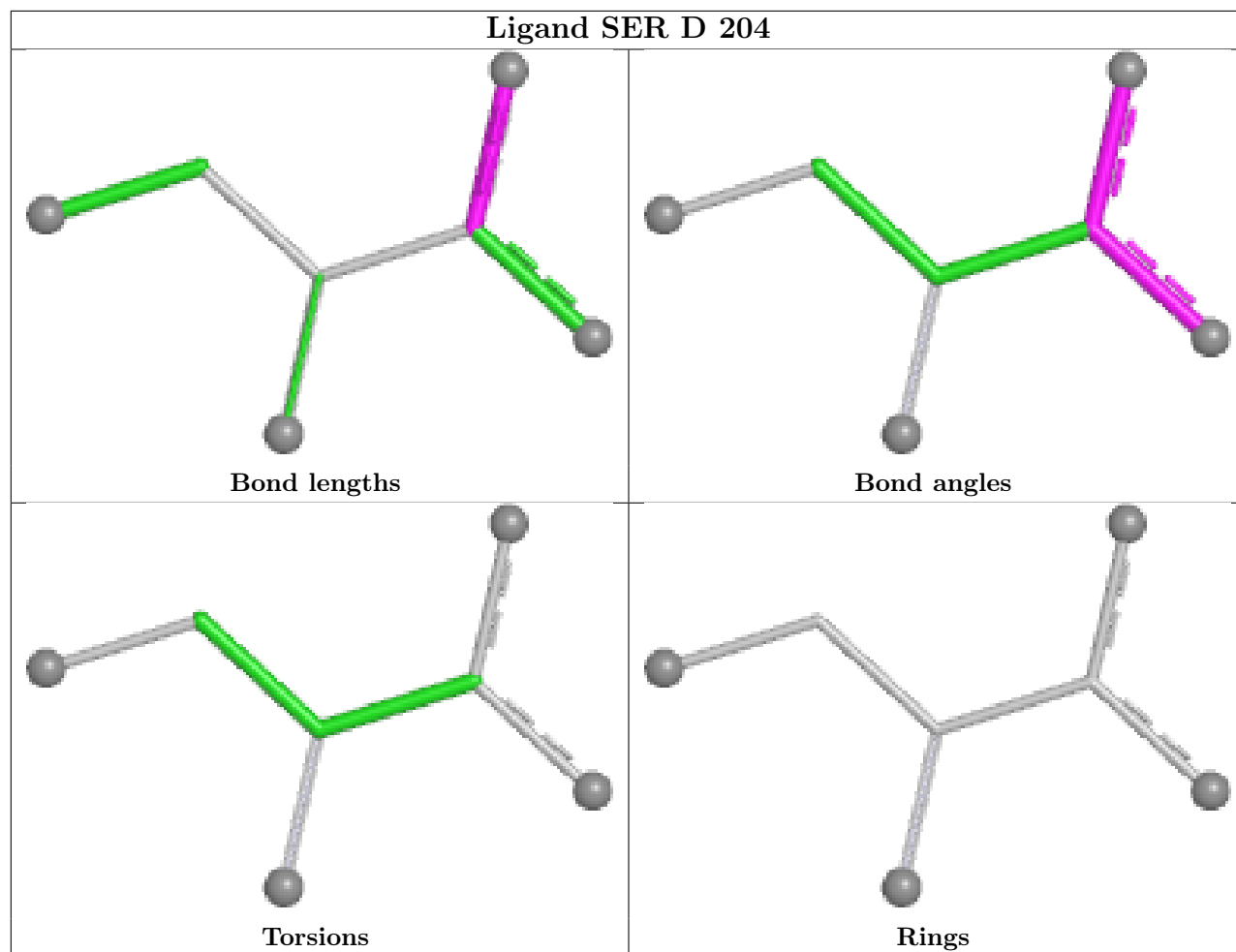
There are no ring outliers.

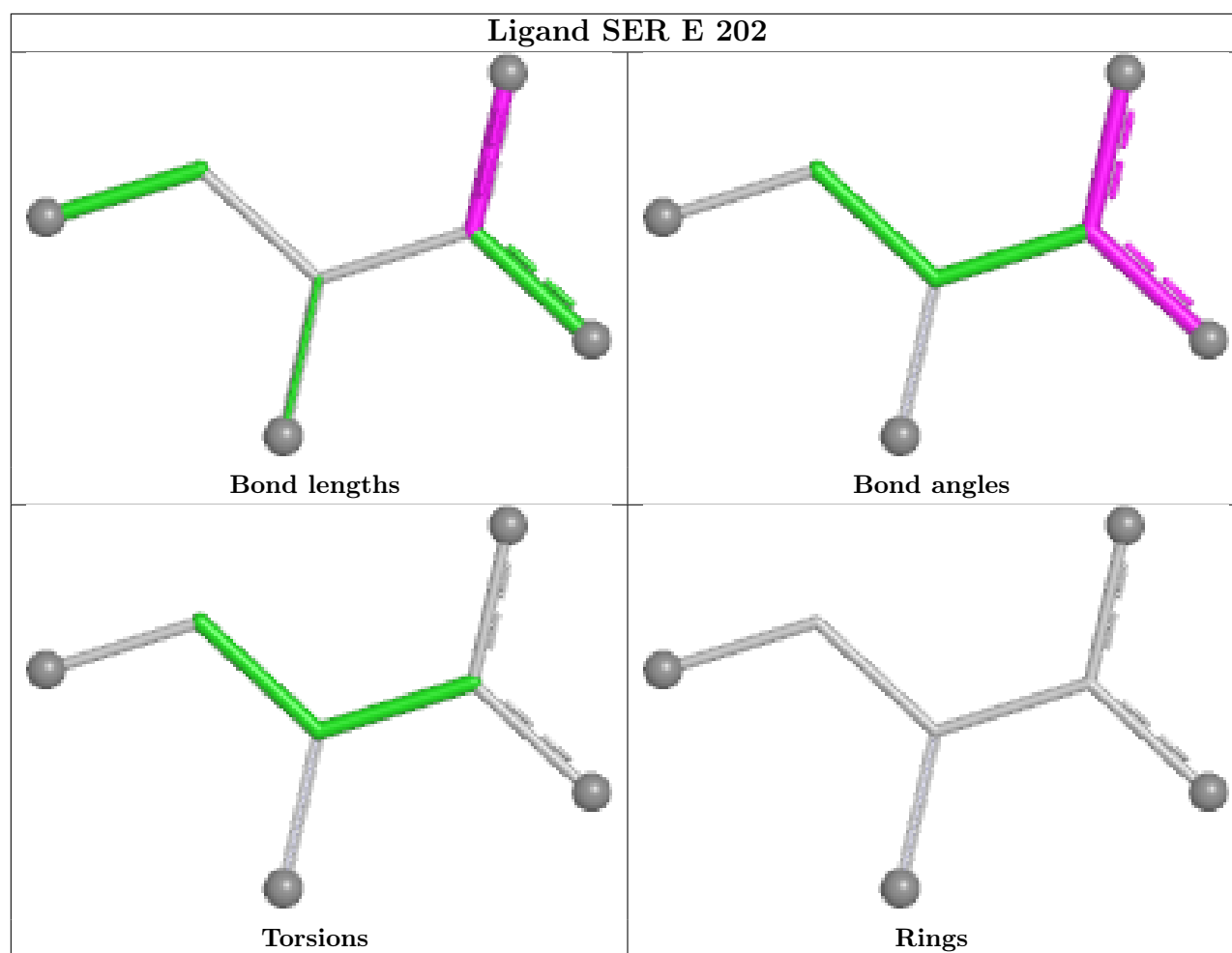
No monomer is involved in short contacts.

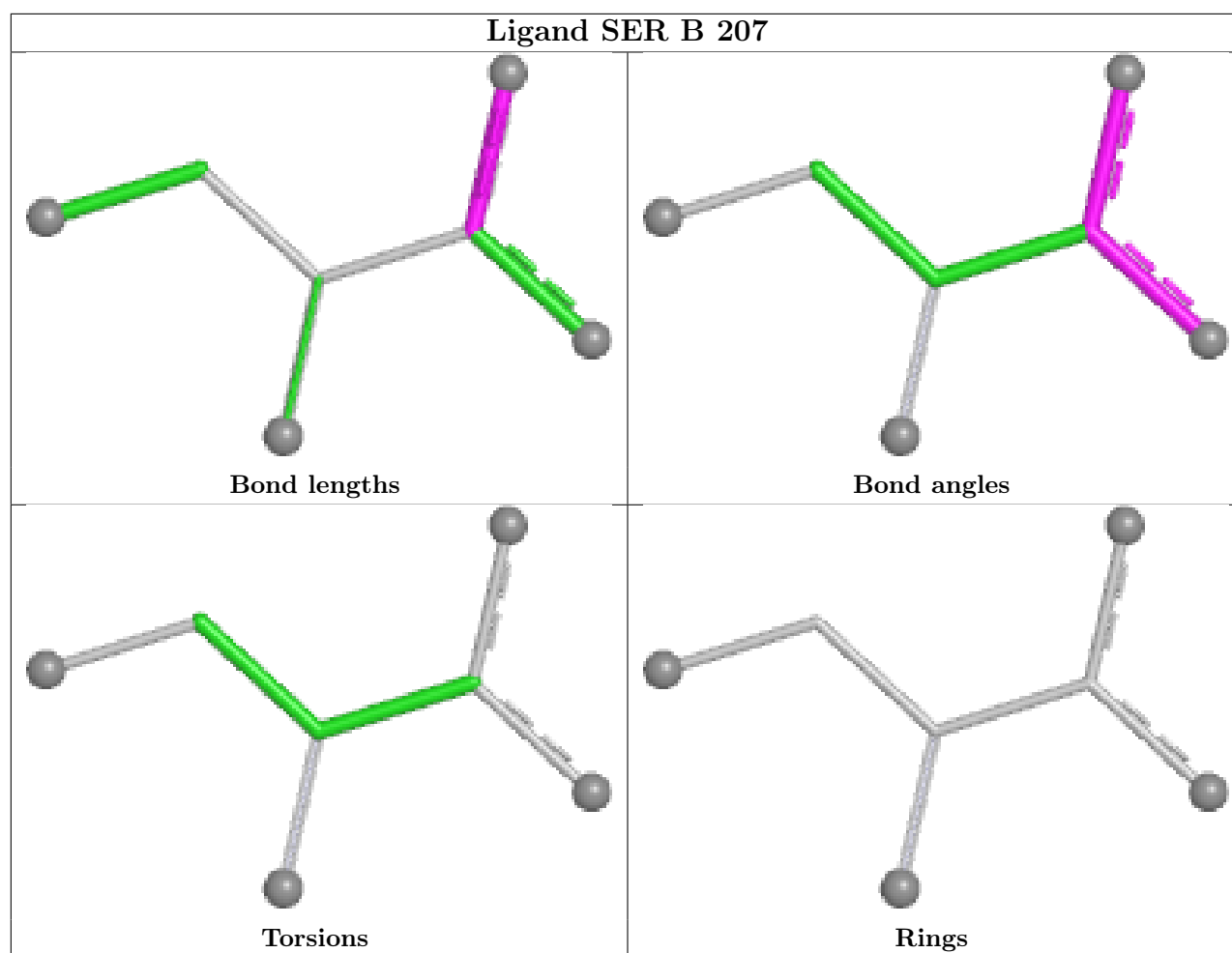
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the

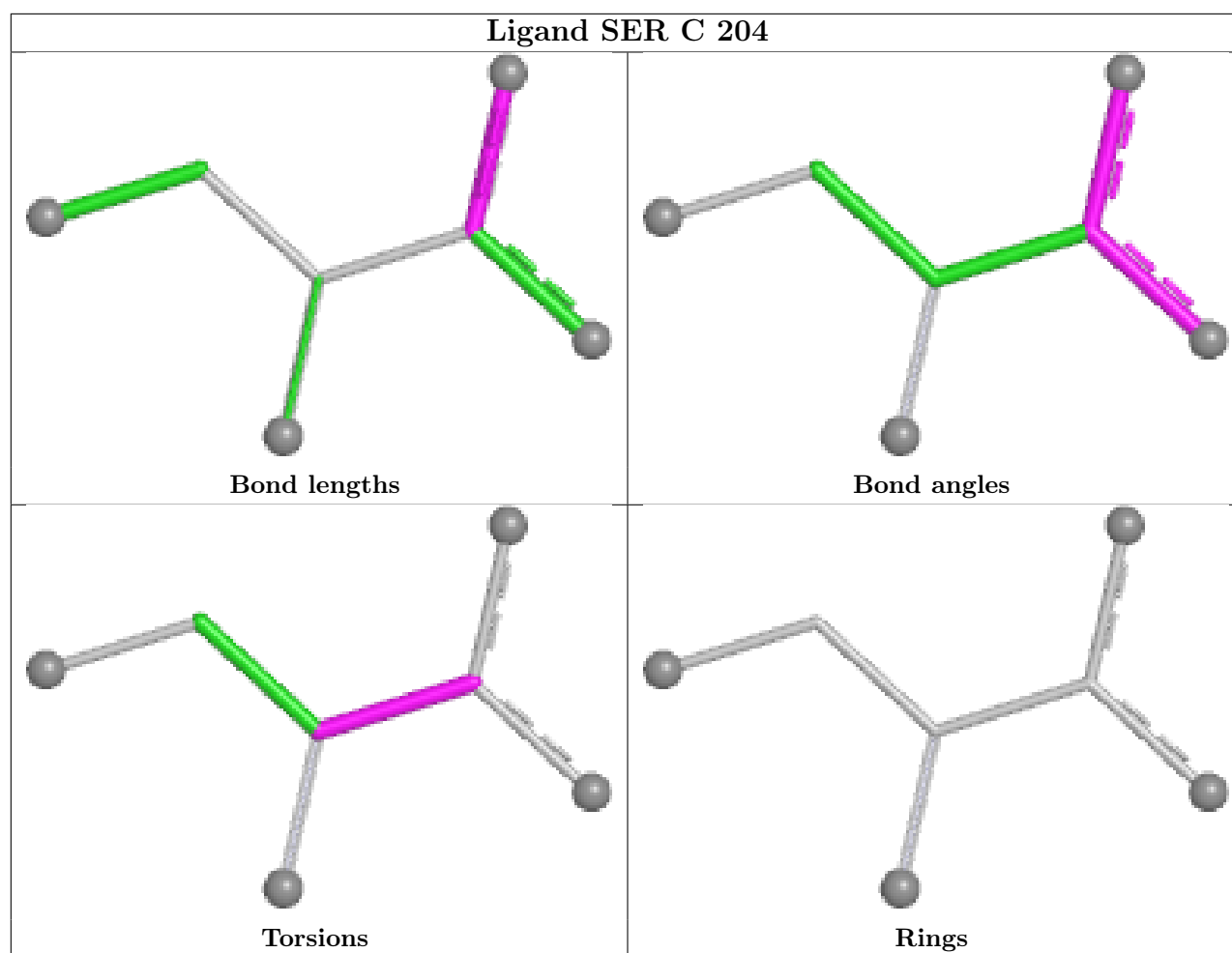
average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.











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(Å²)	Q<0.9
1	A	141/156 (90%)	1.18	34 (24%)	2	1	30, 46, 107, 144	1 (0%)
1	B	140/156 (89%)	1.19	34 (24%)	2	1	29, 46, 105, 122	1 (0%)
1	C	140/156 (89%)	1.24	32 (22%)	2	1	32, 48, 112, 130	2 (1%)
1	D	140/156 (89%)	1.24	32 (22%)	2	1	32, 47, 105, 118	0
1	E	139/156 (89%)	1.33	25 (17%)	3	2	37, 60, 101, 122	0
All	All	700/780 (89%)	1.24	157 (22%)	2	1	29, 50, 107, 144	4 (0%)

All (157) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	122[A]	PHE	12.1
1	A	122[A]	PHE	8.8
1	E	42	VAL	8.0
1	C	182	VAL	6.8
1	E	43	LEU	6.7
1	E	180	ILE	6.6
1	D	42	VAL	6.6
1	C	43	LEU	6.5
1	A	178	TYR	6.2
1	D	178	TYR	6.1
1	E	44	GLN	6.0
1	A	182	VAL	6.0
1	A	42	VAL	5.9
1	B	42	VAL	5.8
1	A	43	LEU	5.8
1	E	178	TYR	5.7
1	C	178	TYR	5.3
1	C	176	ARG	5.0
1	D	181	ALA	5.0
1	C	180	ILE	4.9

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Mol	Chain	Res	Type	RSRZ
1	E	116	ARG	4.8
1	B	43	LEU	4.6
1	D	179	ASP	4.5
1	D	47	ARG	4.5
1	C	122[A]	PHE	4.5
1	D	177	LEU	4.4
1	A	180	ILE	4.4
1	B	121	ALA	4.3
1	B	178	TYR	4.3
1	D	176	ARG	4.2
1	B	176	ARG	4.1
1	B	177	LEU	4.1
1	D	122	PHE	4.1
1	D	46	ILE	4.1
1	A	181	ALA	4.1
1	D	43	LEU	4.1
1	C	113	ARG	4.1
1	A	177	LEU	3.9
1	B	172	GLN	3.9
1	B	181	ALA	3.9
1	A	46	ILE	3.8
1	D	116	ARG	3.8
1	E	179	ASP	3.8
1	C	44	GLN	3.8
1	A	171	MET	3.8
1	C	47	ARG	3.7
1	B	180	ILE	3.7
1	D	170	TYR	3.7
1	D	113	ARG	3.6
1	C	46	ILE	3.5
1	D	180	ILE	3.5
1	A	115	PRO	3.5
1	A	172	GLN	3.5
1	E	47	ARG	3.5
1	C	177	LEU	3.4
1	A	174	ASN	3.4
1	E	46	ILE	3.4
1	C	115	PRO	3.3
1	D	115	PRO	3.2
1	B	115	PRO	3.2
1	A	47	ARG	3.2
1	B	47	ARG	3.2

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Mol	Chain	Res	Type	RSRZ
1	D	175	ASP	3.2
1	A	173	GLN	3.2
1	A	114	ASP	3.1
1	B	44	GLN	3.1
1	B	113	ARG	3.1
1	C	169	ALA	3.1
1	B	119	GLU	3.0
1	A	44	GLN	3.0
1	E	144	GLY	2.9
1	C	116	ARG	2.9
1	E	81	ILE	2.9
1	E	143	LEU	2.9
1	C	171[A]	MET	2.9
1	A	169	ALA	2.8
1	D	107	GLN	2.8
1	C	181	ALA	2.8
1	E	45	THR	2.7
1	C	114	ASP	2.7
1	D	44	GLN	2.7
1	E	138	GLU	2.7
1	A	83	SER	2.7
1	C	112	PRO	2.7
1	B	170	TYR	2.7
1	B	48	GLN	2.7
1	B	179	ASP	2.7
1	D	102	GLU	2.7
1	A	170	TYR	2.7
1	D	171	MET	2.7
1	E	177	LEU	2.7
1	A	107	GLN	2.6
1	C	170	TYR	2.6
1	E	176	ARG	2.6
1	C	173	GLN	2.6
1	E	114	ASP	2.6
1	B	116	ARG	2.6
1	A	45	THR	2.6
1	C	82	GLY	2.6
1	B	171	MET	2.6
1	E	85	ALA	2.5
1	B	54	ASN	2.5
1	C	45	THR	2.5
1	B	125	ILE	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	82	GLY	2.5
1	B	45	THR	2.5
1	D	45	THR	2.5
1	B	150	GLU	2.5
1	B	103	LYS	2.5
1	D	50	GLN	2.5
1	E	102	GLU	2.4
1	A	50	GLN	2.4
1	D	111	LEU	2.4
1	E	142	LEU	2.4
1	A	48	GLN	2.4
1	E	171	MET	2.4
1	A	116	ARG	2.4
1	A	102	GLU	2.4
1	C	146	GLY	2.4
1	B	173	GLN	2.4
1	C	48	GLN	2.3
1	C	172[A]	GLN	2.3
1	E	117	GLN	2.3
1	A	175	ASP	2.3
1	D	172	GLN	2.3
1	E	59	GLU	2.3
1	E	172	GLN	2.3
1	A	113	ARG	2.2
1	B	120	ALA	2.2
1	B	50	GLN	2.2
1	B	117	GLN	2.2
1	C	175	ASP	2.2
1	A	176	ARG	2.2
1	D	48	GLN	2.2
1	E	106	GLU	2.2
1	D	104	ASN	2.2
1	D	174	ASN	2.2
1	C	102	GLU	2.2
1	C	50	GLN	2.2
1	B	110	ALA	2.2
1	A	81	ILE	2.1
1	B	123[A]	LEU	2.1
1	D	153	ASP	2.1
1	C	117	GLN	2.1
1	C	120	ALA	2.1
1	D	51	SER	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	179	ASP	2.1
1	B	114	ASP	2.1
1	C	119	GLU	2.1
1	D	114	ASP	2.1
1	D	117	GLN	2.1
1	B	105	TRP	2.1
1	A	105	TRP	2.1
1	D	138	GLU	2.0
1	A	53	LEU	2.0
1	C	179	ASP	2.0
1	B	118	SER	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	NA	B	204	1/1	0.82	0.18	49,49,49,49	0
4	NA	B	202	1/1	0.88	0.24	58,58,58,58	0
2	CL	E	201	1/1	0.89	0.30	53,53,53,53	0
3	SER	B	207	7/7	0.90	0.13	31,35,43,44	0
2	CL	B	206	1/1	0.91	0.23	47,47,47,47	0
4	NA	B	201	1/1	0.92	0.16	41,41,41,41	0
2	CL	B	205	1/1	0.92	0.22	48,48,48,48	0
3	SER	E	202	7/7	0.92	0.11	42,50,64,64	0
4	NA	C	201	1/1	0.92	0.10	39,39,39,39	0
3	SER	D	204	7/7	0.93	0.10	35,39,49,49	0
4	NA	D	201	1/1	0.93	0.13	43,43,43,43	0
3	SER	A	203	7/7	0.95	0.08	28,29,40,41	0

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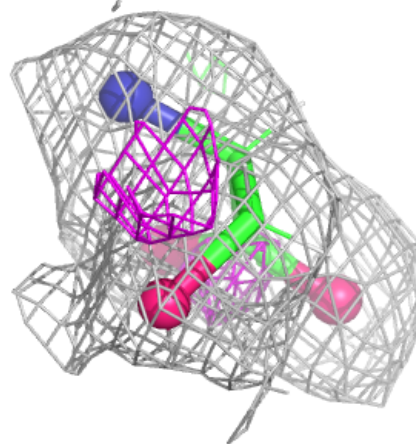
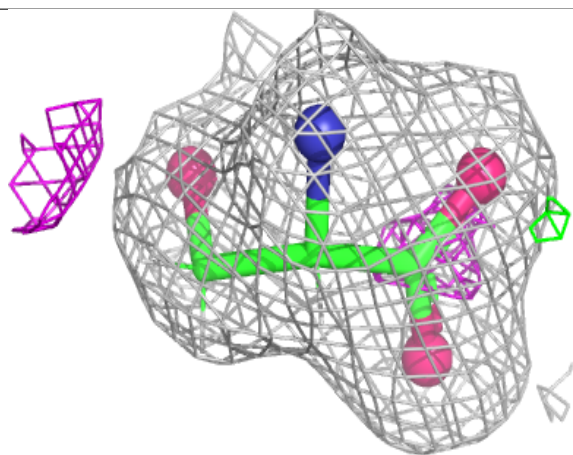
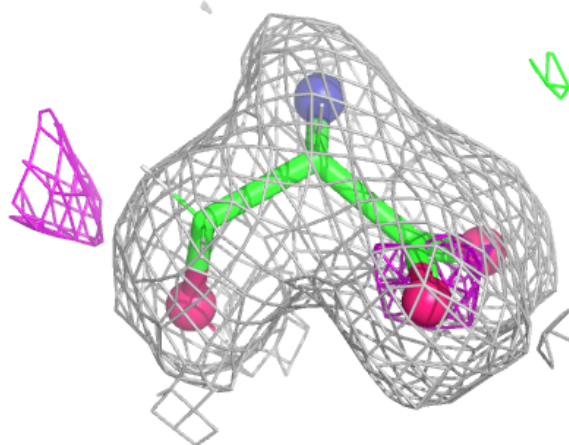
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	SER	C	204	7/7	0.95	0.08	30,33,40,41	0
4	NA	B	203	1/1	0.95	0.10	44,44,44,44	0
2	CL	A	201	1/1	0.96	0.13	45,45,45,45	0
2	CL	D	202	1/1	0.96	0.13	47,47,47,47	0
2	CL	C	202	1/1	0.97	0.17	39,39,39,39	0
2	CL	C	203	1/1	0.97	0.21	41,41,41,41	0
2	CL	A	202	1/1	0.97	0.14	43,43,43,43	0
2	CL	D	203	1/1	0.97	0.20	45,45,45,45	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.

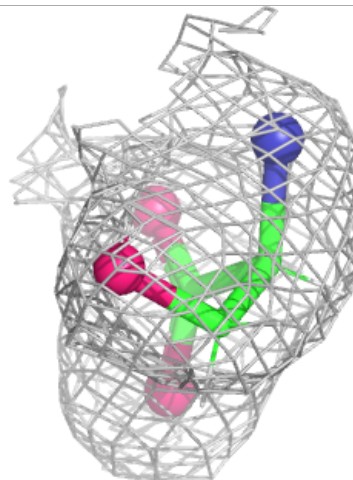
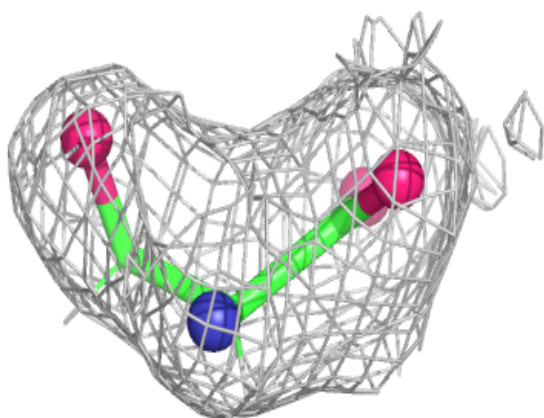
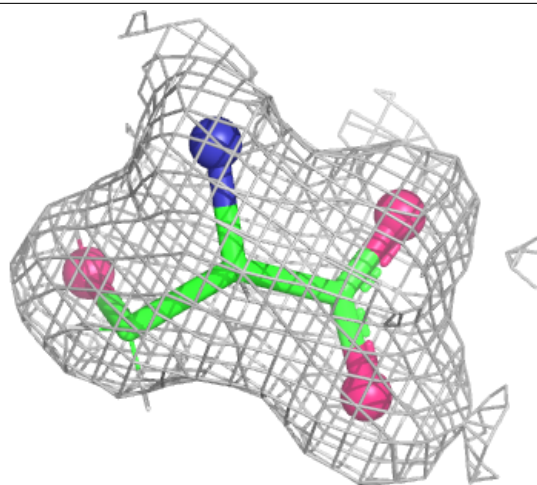
Electron density around SER B 207:

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



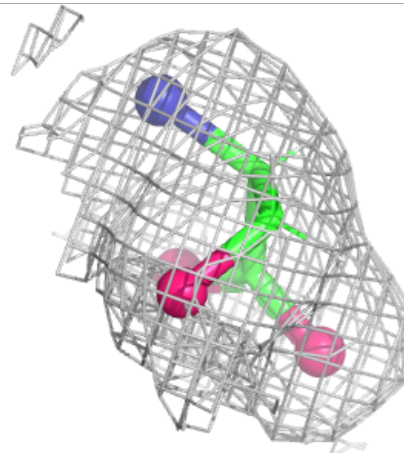
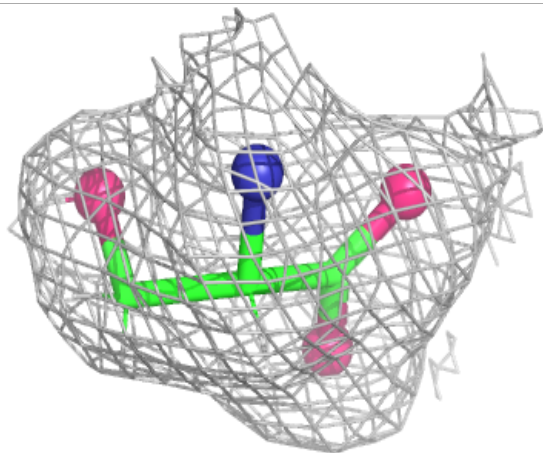
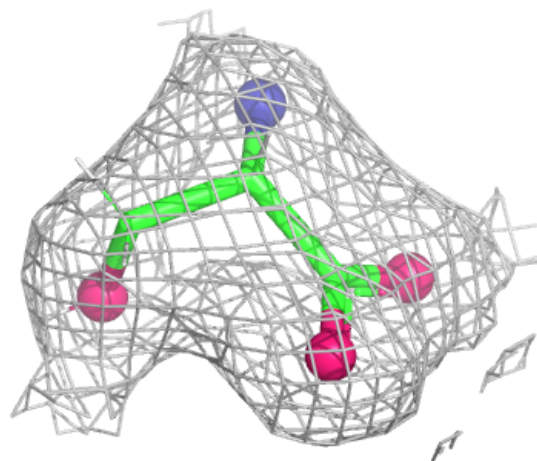
Electron density around SER E 202:

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



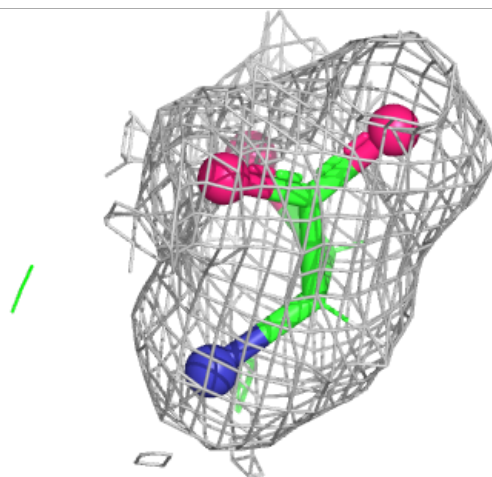
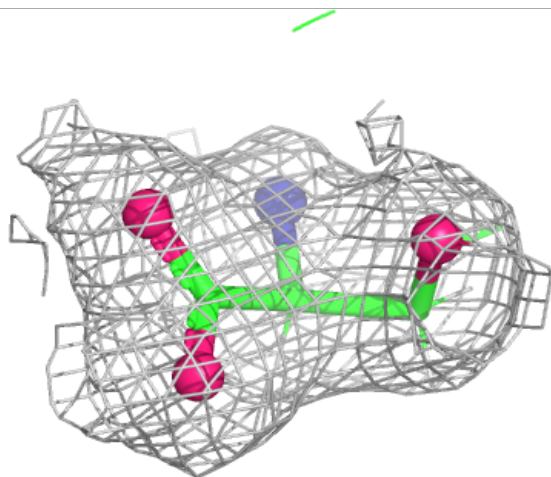
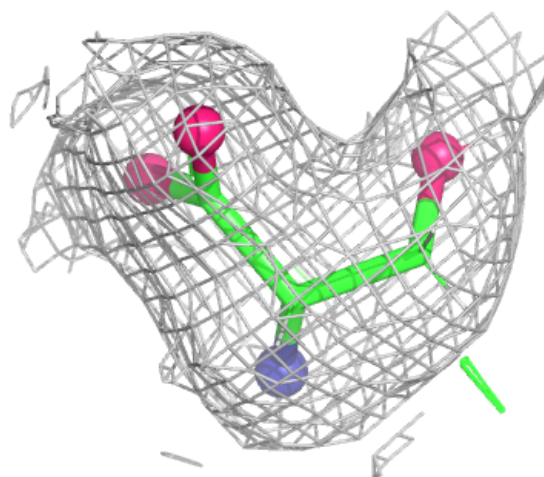
Electron density around SER D 204:

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



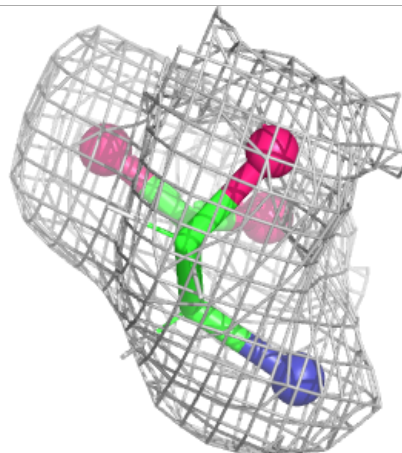
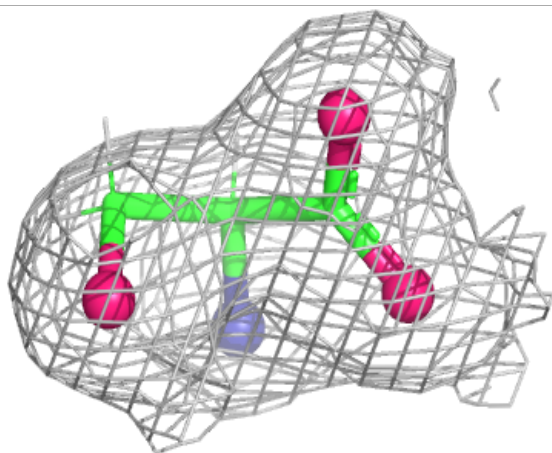
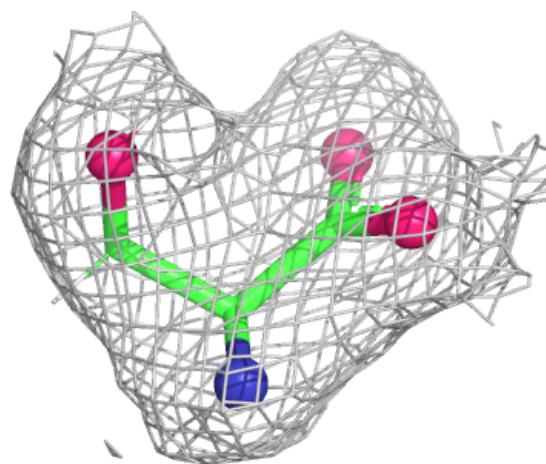
Electron density around SER A 203:

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



Electron density around SER C 204:

$2mF_o-DF_c$ (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.