



wwPDB X-ray Structure Validation Summary Report ⓘ

May 26, 2020 – 09:21 am BST

PDB ID : 2VU1
Title : Biosynthetic thiolase from *Z. ramigera*. Complex of with O-pantheteine- 11-pivalate.
Authors : Kursula, P.; Schmitz, W.; Wierenga, R.K.
Deposited on : 2008-05-19
Resolution : 1.51 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

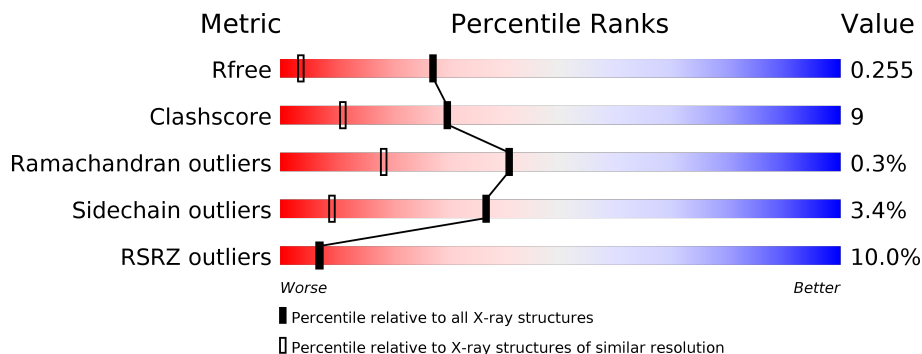
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.51 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}	130704	4009 (1.54-1.50)
Clashscore	141614	4249 (1.54-1.50)
Ramachandran outliers	138981	4148 (1.54-1.50)
Sidechain outliers	138945	4146 (1.54-1.50)
RSRZ outliers	127900	3943 (1.54-1.50)

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 on 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	392	<div style="display: flex; align-items: center;"> <div style="width: 2%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 84%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 15%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 2%; height: 10px; background-color: orange; margin-right: 5px;"></div> <div style="width: 2%; height: 10px; background-color: grey;"></div> </div> <p style="font-size: small; margin-top: 5px;">2% 84% 15% .</p>
1	B	392	<div style="display: flex; align-items: center;"> <div style="width: 2%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 87%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 13%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 2%; height: 10px; background-color: orange; margin-right: 5px;"></div> <div style="width: 2%; height: 10px; background-color: grey;"></div> </div> <p style="font-size: small; margin-top: 5px;">2% 87% 13% .</p>
1	C	392	<div style="display: flex; align-items: center;"> <div style="width: 8%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 81%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 18%; height: 10px; background-color: yellow; margin-right: 5px;"></div> </div> <p style="font-size: small; margin-top: 5px;">8% 81% 18%</p>
1	D	392	<div style="display: flex; align-items: center;"> <div style="width: 28%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 72%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 26%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 2%; height: 10px; background-color: orange; margin-right: 5px;"></div> <div style="width: 2%; height: 10px; background-color: grey;"></div> </div> <p style="font-size: small; margin-top: 5px;">28% 72% 26% ..</p>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 12631 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.

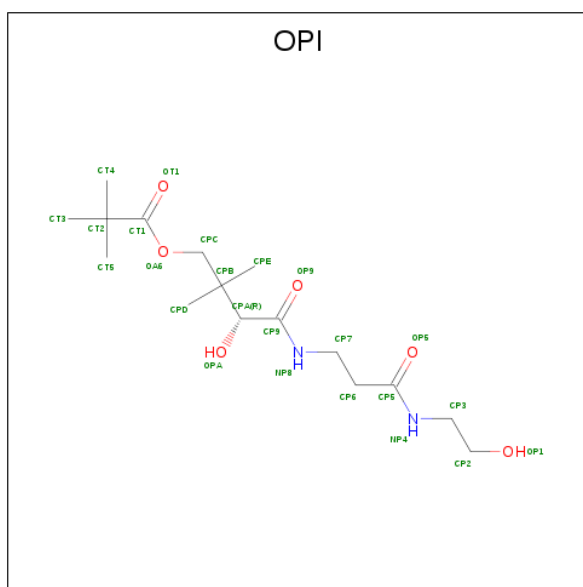
- Molecule 1 is a protein called ACETYL-COA ACETYLTRANSFERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	391	2909	1819	520	547	23	0	16	0
1	B	392	2903	1811	521	549	22	0	13	0
1	C	391	2828	1755	511	541	21	0	0	0
1	D	390	2821	1751	510	539	21	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	11	ALA	-	insertion	UNP P07097
A	129	ARG	ALA	conflict	UNP P07097
B	11	ALA	-	insertion	UNP P07097
B	129	ARG	ALA	conflict	UNP P07097
C	11	ALA	-	insertion	UNP P07097
C	129	ARG	ALA	conflict	UNP P07097
D	11	ALA	-	insertion	UNP P07097
D	129	ARG	ALA	conflict	UNP P07097

- Molecule 2 is PANTOTHENYL-AMINOETHANOL-11-PIVALIC ACID (three-letter code: OPI) (formula: $C_{16}H_{30}N_2O_6$).

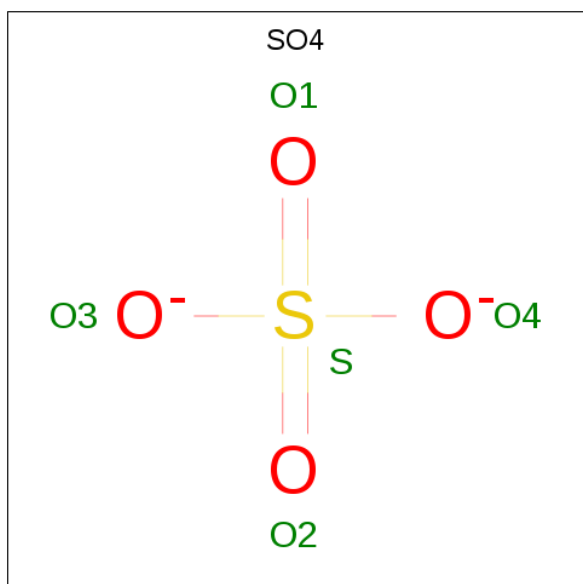


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	
			Total	C	N			O
2	A	1	24	16	2	6	0	0
2	B	1	24	16	2	6	0	0

- Molecule 3 is SODIUM ION (three-letter code: NA) (formula: Na).

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

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total O S 5 4 1	0	0
4	A	1	Total O S 5 4 1	0	0
4	B	1	Total O S 5 4 1	0	0
4	B	1	Total O S 5 4 1	0	0
4	D	1	Total O S 5 4 1	0	0

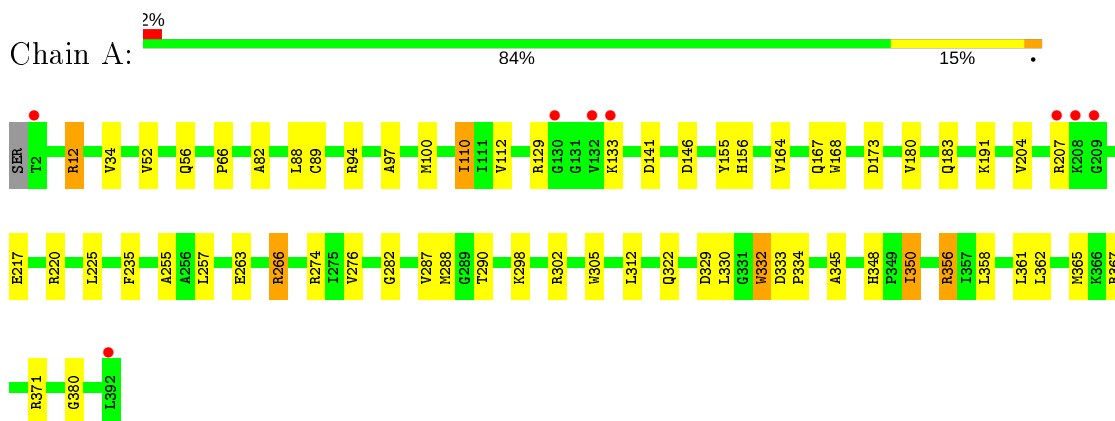
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	408	Total O 408 408	0	0
5	B	398	Total O 398 398	1	0
5	C	144	Total O 144 144	0	0
5	D	146	Total O 146 146	0	0

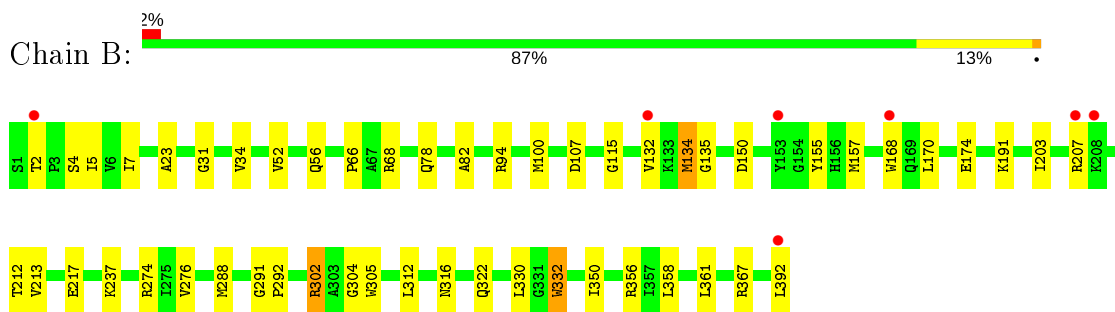
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA 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.

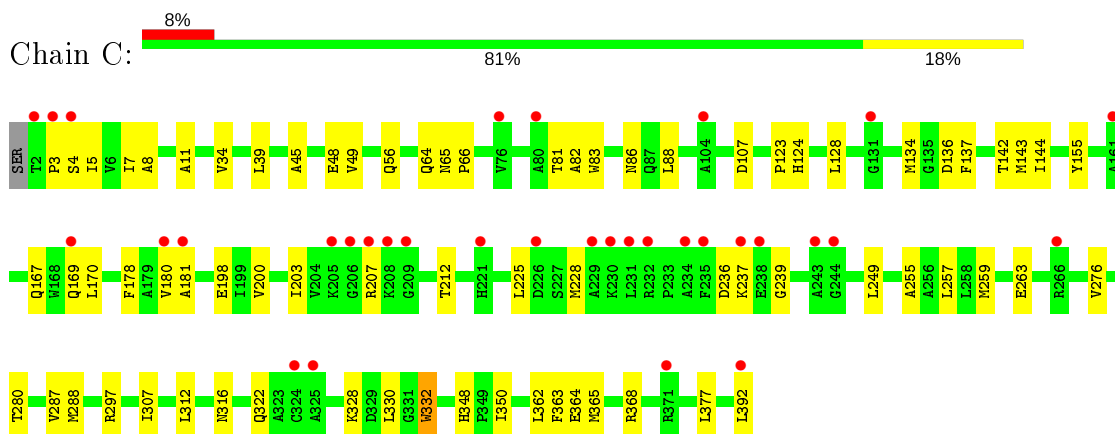
- Molecule 1: ACETYL-COA ACETYLTRANSFERASE



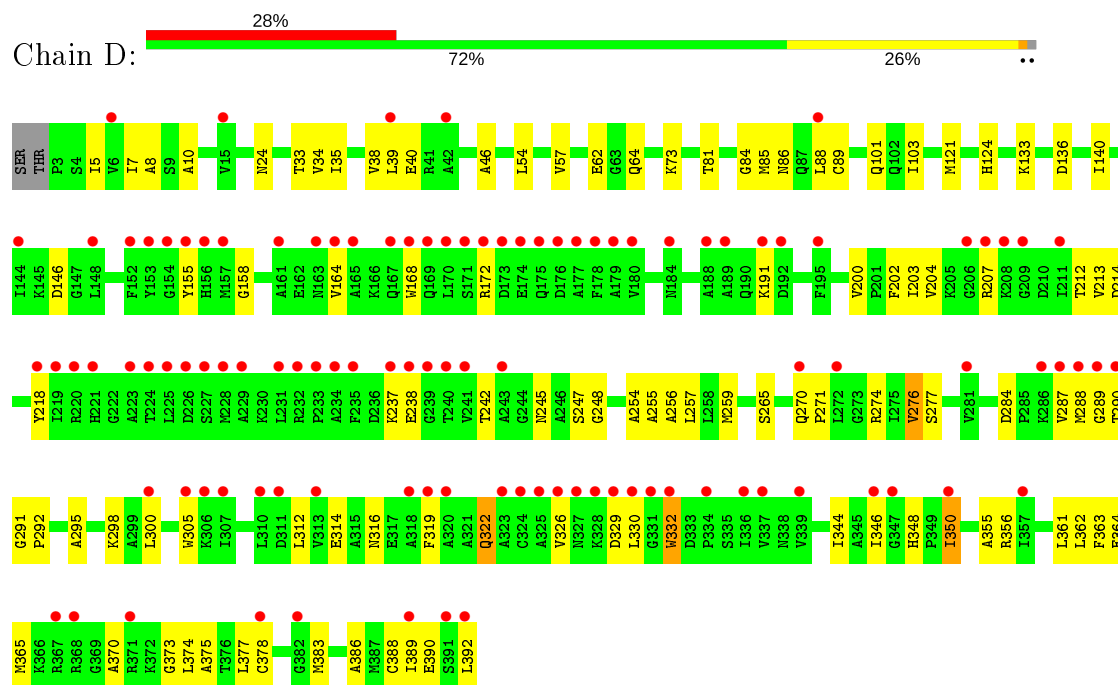
- Molecule 1: ACETYL-COA ACETYLTRANSFERASE



- Molecule 1: ACETYL-COA ACETYLTRANSFERASE



● Molecule 1: ACETYL-COA ACETYLTRANSFERASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	84.30Å 78.74Å 148.34Å 90.00° 92.93° 90.00°	Depositor
Resolution (Å)	20.00 – 1.51 19.98 – 1.51	Depositor EDS
% Data completeness (in resolution range)	92.5 (20.00-1.51) 79.9 (19.98-1.51)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.13 (at 1.51Å)	Xtriage
Refinement program	REFMAC 5.3.0028	Depositor
R, R_{free}	0.215 , 0.247 0.225 , 0.255	Depositor DCC
R_{free} test set	2810 reflections (1.00%)	wwPDB-VP
Wilson B-factor (Å ²)	13.1	Xtriage
Anisotropy	0.300	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 35.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.37$, $\langle L^2 \rangle = 0.20$	Xtriage
Estimated twinning fraction	0.136 for h,-k,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	12631	wwPDB-VP
Average B, all atoms (Å ²)	12.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.11% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: NA, CSO, SO4, OPI

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	1.11	3/2986 (0.1%)	1.06	11/4031 (0.3%)
1	B	1.07	7/2976 (0.2%)	1.01	8/4015 (0.2%)
1	C	0.63	0/2862	0.72	1/3864 (0.0%)
1	D	0.57	0/2855	0.69	1/3853 (0.0%)
All	All	0.88	10/11679 (0.1%)	0.89	21/15763 (0.1%)

The worst 5 of 10 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	115	GLY	N-CA	-6.20	1.36	1.46
1	B	168	TRP	CB-CG	5.85	1.60	1.50
1	A	217	GLU	CB-CG	-5.76	1.41	1.52
1	B	23	ALA	CA-CB	-5.61	1.40	1.52
1	A	362	LEU	C-O	-5.55	1.12	1.23

The worst 5 of 21 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	94	ARG	NE-CZ-NH2	-9.38	115.61	120.30
1	A	146	ASP	CB-CG-OD1	8.58	126.02	118.30
1	B	367	ARG	NE-CZ-NH2	-8.39	116.11	120.30
1	A	367	ARG	NE-CZ-NH1	7.70	124.15	120.30
1	A	129	ARG	NE-CZ-NH1	-6.56	117.02	120.30

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	2909	0	2972	44	0
1	B	2903	0	2956	39	0
1	C	2828	0	2833	53	0
1	D	2821	0	2827	87	0
2	A	24	0	30	4	0
2	B	24	0	30	3	0
3	A	1	0	0	0	0
4	A	10	0	0	2	0
4	B	10	0	0	0	0
4	D	5	0	0	0	0
5	A	408	0	0	8	2
5	B	398	0	0	9	2
5	C	144	0	0	11	0
5	D	146	0	0	22	0
All	All	12631	0	11648	218	4

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 9.

The worst 5 of 218 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:274[A]:ARG:NH1	1:B:276[A]:VAL:HG12	1.45	1.28
1:D:164:VAL:HG21	5:D:2122:HOH:O	1.47	1.11
1:A:164:VAL:HG22	1:A:287[B]:VAL:HG23	1.38	1.06
1:D:88:LEU:CD2	5:D:2059:HOH:O	2.09	1.00
1:D:146:ASP:O	5:D:2059:HOH:O	1.82	0.95

All (4) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A:2078:HOH:O	5:A:2211:HOH:O[2_655]	1.77	0.43
5:B:2101:HOH:O	5:B:2375:HOH:O[2_545]	2.03	0.17

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A:2118:HOH:O	5:A:2392:HOH:O[2_655]	2.11	0.09
5:B:2064:HOH:O	5:B:2222:HOH:O[2_545]	2.14	0.06

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	403/392 (103%)	388 (96%)	14 (4%)	1 (0%)	47	23
1	B	401/392 (102%)	390 (97%)	10 (2%)	1 (0%)	47	23
1	C	388/392 (99%)	375 (97%)	11 (3%)	2 (0%)	29	9
1	D	387/392 (99%)	373 (96%)	13 (3%)	1 (0%)	41	18
All	All	1579/1568 (101%)	1526 (97%)	48 (3%)	5 (0%)	41	18

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	350	ILE
1	C	350	ILE
1	D	350	ILE
1	A	350	ILE
1	C	4	SER

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	292/278 (105%)	285 (98%)	7 (2%)	49	19
1	B	291/278 (105%)	284 (98%)	7 (2%)	49	19
1	C	277/278 (100%)	267 (96%)	10 (4%)	35	8
1	D	276/278 (99%)	262 (95%)	14 (5%)	24	3
All	All	1136/1112 (102%)	1098 (97%)	38 (3%)	37	10

5 of 38 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	228	MET
1	C	288	MET
1	D	288	MET
1	C	263	GLU
1	C	322	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 17 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	156	HIS
1	C	169	GLN
1	D	169	GLN
1	C	78	GLN
1	D	184	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](#)

5 non-standard protein/DNA/RNA residues are modelled in this entry.

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
1	CSO	A	89[A]	-	3,5,7	0.93	0	0,5,8	0.00	-
1	CSO	D	89	1	3,6,7	0.55	0	0,6,8	0.00	-
1	CSO	B	89	1	3,6,7	0.59	0	0,6,8	0.00	-
1	CSO	A	89[B]	-	3,6,7	1.28	1 (33%)	0,6,8	0.00	-
1	CSO	C	89	1	3,6,7	0.37	0	0,6,8	0.00	-

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
1	CSO	A	89[A]	-	-	0/1/4/7	-
1	CSO	D	89	1	-	0/1/5/7	-
1	CSO	B	89	1	-	1/1/5/7	-
1	CSO	A	89[B]	-	-	1/1/5/7	-
1	CSO	C	89	1	-	1/1/5/7	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	89[B]	CSO	CB-CA	2.08	1.58	1.53

There are no bond angle outliers.

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	C	89	CSO	N-CA-CB-SG
1	B	89	CSO	N-CA-CB-SG
1	A	89[B]	CSO	N-CA-CB-SG

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	89[A]	CSO	2	0
1	D	89	CSO	1	0
1	A	89[B]	CSO	1	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 1 is monoatomic - leaving 7 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	SO4	D	1393	-	4,4,4	0.17	0	6,6,6	0.27	0
4	SO4	A	1395	-	4,4,4	0.18	0	6,6,6	0.67	0
4	SO4	A	1396	-	4,4,4	0.12	0	6,6,6	0.36	0
2	OPI	A	1393	-	20,23,23	1.13	1 (5%)	30,32,32	1.89	6 (20%)
4	SO4	B	1395	-	4,4,4	0.18	0	6,6,6	0.66	0
2	OPI	B	1393	-	20,23,23	1.19	1 (5%)	30,32,32	1.57	4 (13%)
4	SO4	B	1394	-	4,4,4	0.13	0	6,6,6	0.28	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	OPI	A	1393	-	-	14/32/32/32	-
2	OPI	B	1393	-	-	12/32/32/32	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1393	OPI	OA6-CT1	5.00	1.44	1.33
2	A	1393	OPI	OA6-CT1	4.57	1.43	1.33

The worst 5 of 10 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1393	OPI	OA6-CT1-CT2	5.97	124.48	112.56
2	A	1393	OPI	OA6-CT1-CT2	5.75	124.03	112.56
2	A	1393	OPI	OA6-CT1-OT1	-3.97	116.37	124.54
2	A	1393	OPI	CPC-OA6-CT1	-3.94	111.24	116.81
2	B	1393	OPI	OA6-CT1-OT1	-3.49	117.36	124.54

There are no chirality outliers.

5 of 26 torsion outliers are listed below:

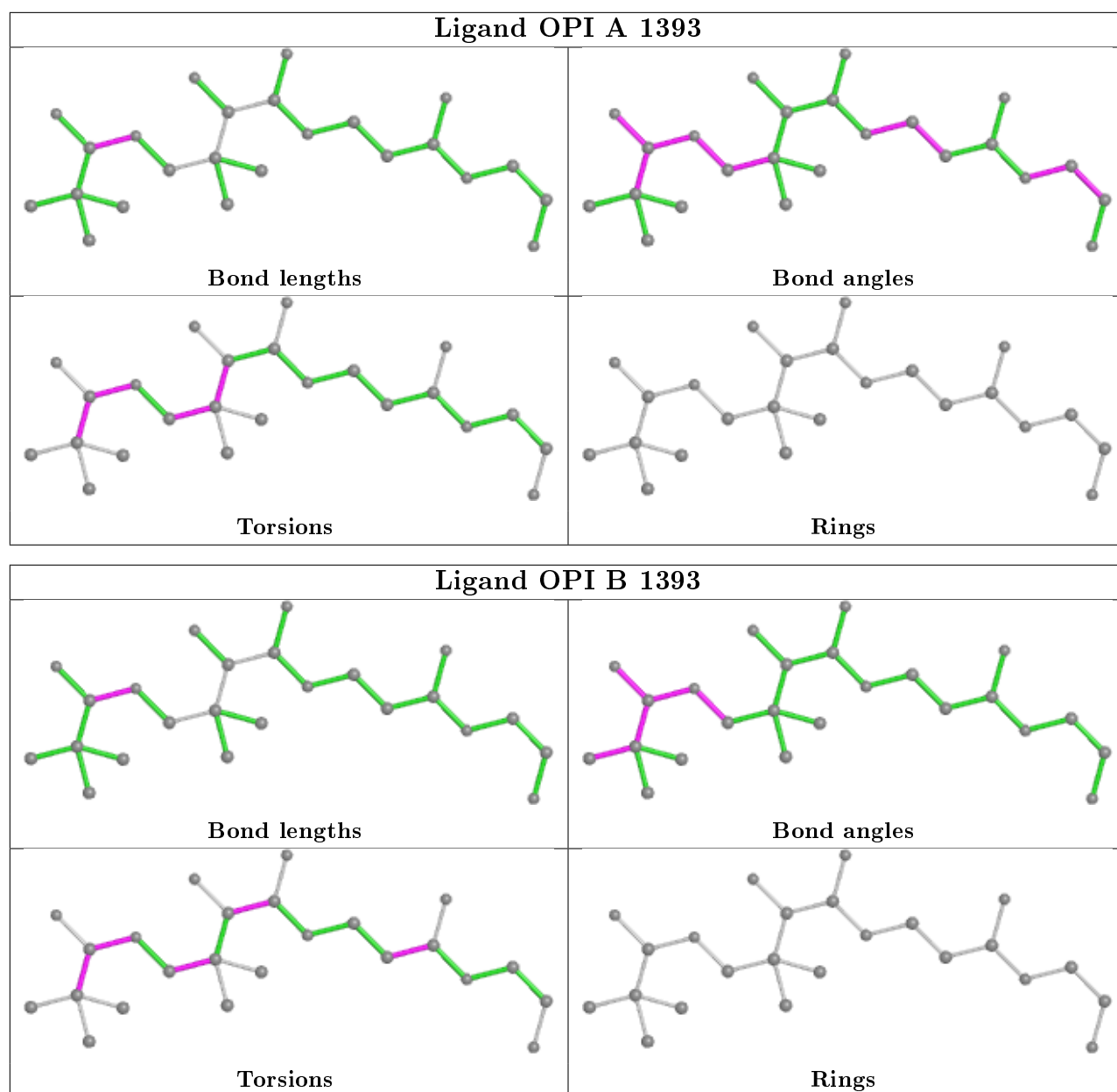
Mol	Chain	Res	Type	Atoms
2	A	1393	OPI	CP9-CPA-CPB-CPD
2	A	1393	OPI	CP9-CPA-CPB-CPE
2	A	1393	OPI	CP9-CPA-CPB-CPC
2	A	1393	OPI	OPA-CPA-CPB-CPD
2	A	1393	OPI	OPA-CPA-CPB-CPE

There are no ring outliers.

4 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1395	SO4	1	0
4	A	1396	SO4	1	0
2	A	1393	OPI	4	0
2	B	1393	OPI	3	0

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



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	390/392 (99%)	0.09	8 (2%) 63 68	4, 9, 20, 42	0
1	B	391/392 (99%)	-0.00	7 (1%) 68 73	3, 9, 20, 44	0
1	C	390/392 (99%)	0.78	33 (8%) 10 11	4, 11, 19, 40	0
1	D	389/392 (99%)	1.49	108 (27%) 0 0	3, 10, 19, 40	0
All	All	1560/1568 (99%)	0.59	156 (10%) 7 7	3, 10, 20, 44	0

The worst 5 of 156 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	229	ALA	8.3
1	D	232	ARG	7.8
1	D	170	LEU	7.5
1	D	331	GLY	6.3
1	D	319	PHE	6.2

6.2 Non-standard residues in protein, DNA, RNA chains [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
1	CSO	D	89	7/8	0.87	0.11	8,9,11,14	0
1	CSO	C	89	7/8	0.94	0.09	10,11,14,15	0
1	CSO	A	89[B]	7/8	0.96	0.09	2,7,9,10	4
1	CSO	A	89[A]	6/8	0.96	0.09	7,9,9,15	3
1	CSO	B	89	7/8	0.97	0.06	7,8,12,20	0

6.3 Carbohydrates

There are no carbohydrates in this entry.

6.4 Ligands

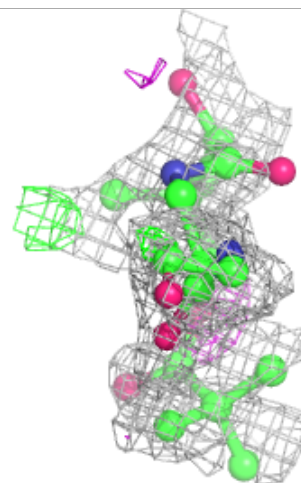
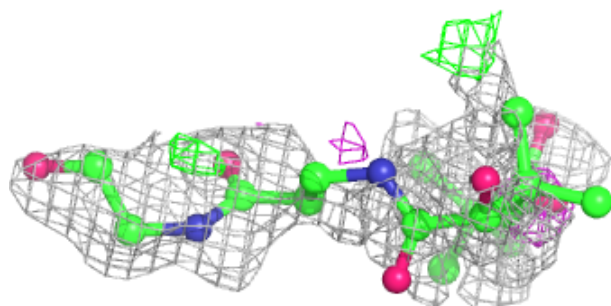
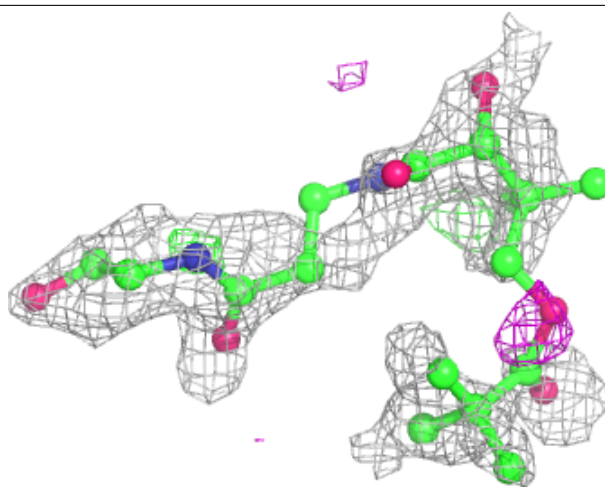
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

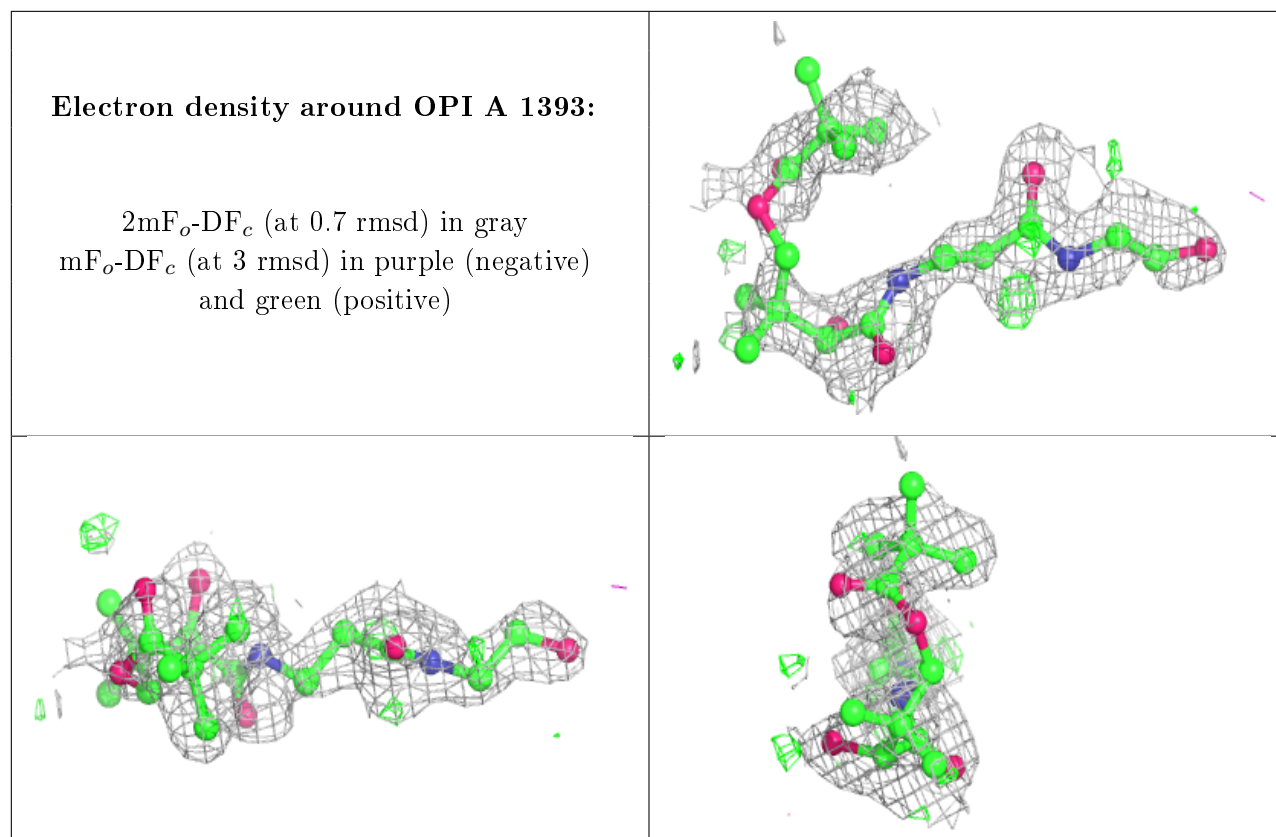
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	OPI	B	1393	24/24	0.48	0.33	55,63,65,67	0
2	OPI	A	1393	24/24	0.67	0.20	36,46,50,50	0
3	NA	A	1394	1/1	0.69	0.17	57,57,57,57	0
4	SO4	A	1396	5/5	0.91	0.22	45,47,47,49	0
4	SO4	D	1393	5/5	0.92	0.20	63,64,65,65	0
4	SO4	B	1394	5/5	0.92	0.18	57,57,58,58	0
4	SO4	A	1395	5/5	0.97	0.11	33,33,36,36	0
4	SO4	B	1395	5/5	0.98	0.10	25,25,28,30	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 OPI B 1393:

$2mF_o-DF_c$ (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.