



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 04:44 PM UTC

PDB ID : 1D6H / pdb_00001d6h
Title : CHALONE SYNTHASE (N336A MUTANT COMPLEXED WITH COA)
Authors : Jez, J.M.; Ferrer, J.L.; Bowman, M.E.; Dixon, R.A.; Noel, J.P.
Deposited on : 1999-10-13
Resolution : 2.15 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

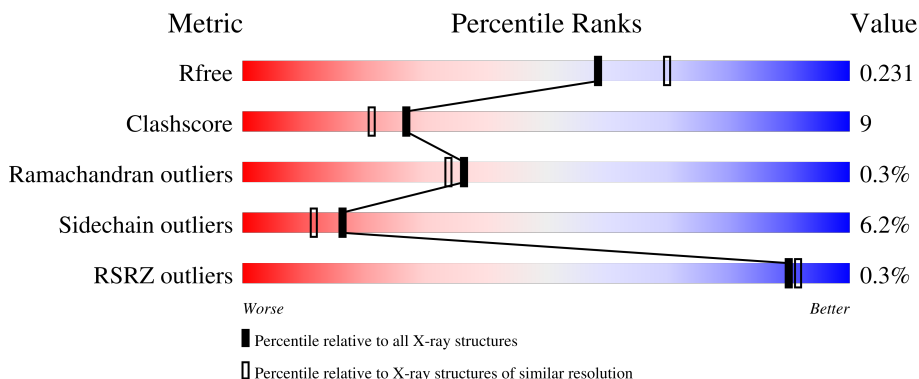
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.15 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	2057 (2.16-2.16)
Clashscore	190562	2159 (2.16-2.16)
Ramachandran outliers	187476	2134 (2.16-2.16)
Sidechain outliers	187428	2133 (2.16-2.16)
RSRZ outliers	180081	2059 (2.16-2.16)

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	387	 59% 31% 7%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 3224 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.

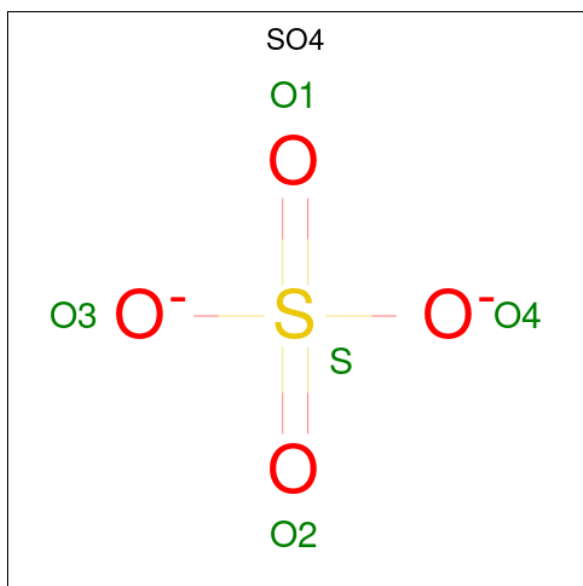
- Molecule 1 is a protein called CHALCONE SYNTHASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	387	2987	1899	503	565	20	0	1	0

There is a discrepancy between the modelled and reference sequences:

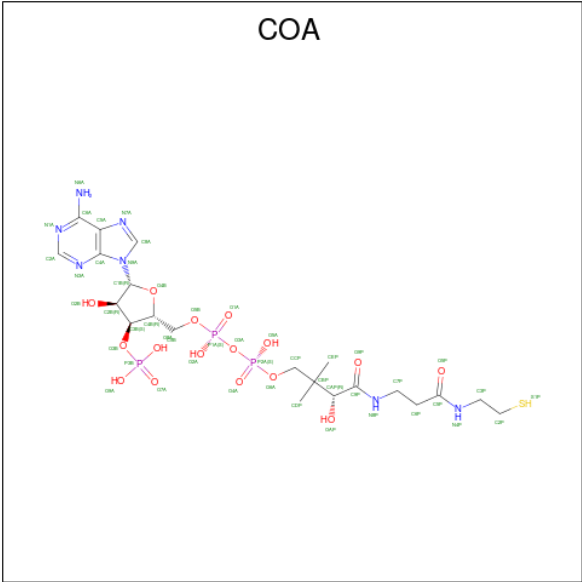
Chain	Residue	Modelled	Actual	Comment	Reference
A	336	ALA	ASN	engineered mutation	UNP P30074

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

- Molecule 3 is COENZYME A (CCD ID: COA) (formula: C₂₁H₃₆N₇O₁₆P₃S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
3	A	1	Total	C	N	O	P	S	0	0
			48	21	7	16	3	1		

- Molecule 4 is water.

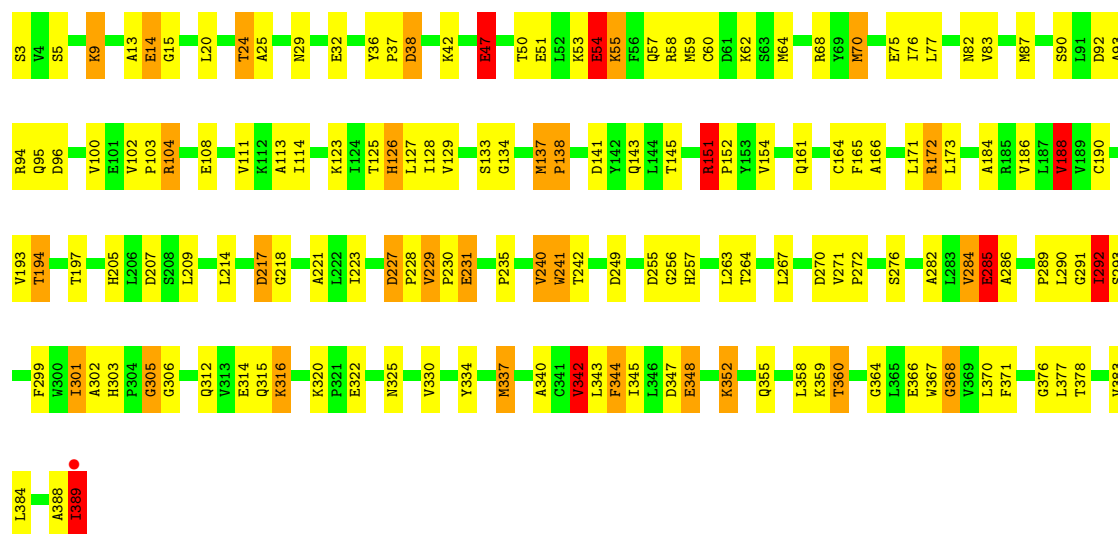
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	184	Total	O	0	0
			184	184		

3 Residue-property plots

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: CHALCONE SYNTHASE

Chain A: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	97.36Å 97.36Å 65.21Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	85.00 – 2.15 84.32 – 2.15	Depositor EDS
% Data completeness (in resolution range)	(Not available) (85.00-2.15) 91.8 (84.32-2.15)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.16 (at 2.15Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.184 , 0.257 (Not available) , 0.231	Depositor DCC
R_{free} test set	932 reflections (5.14%)	wwPDB-VP
Wilson B-factor (Å ²)	29.1	Xtriage
Anisotropy	0.091	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 56.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.043 for -h,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3224	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.34% 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: CSD, SO4, COA

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	1.06	3/3036 (0.1%)	2.16	141/4108 (3.4%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	11

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	126	HIS	CG-ND1	6.43	1.45	1.38
1	A	165	PHE	C-N	5.72	1.42	1.33
1	A	229	VAL	N-CA	5.26	1.50	1.46

All (141) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	376	GLY	O-C-N	-11.37	107.92	122.70
1	A	137	MET	CA-C-O	-10.42	107.62	120.23
1	A	166	ALA	N-CA-C	9.25	124.20	113.19
1	A	94	ARG	CD-NE-CZ	8.90	136.86	124.40
1	A	228	PRO	CA-C-N	-8.82	116.60	122.60
1	A	228	PRO	C-N-CA	-8.82	116.60	122.60
1	A	227	ASP	CA-C-O	8.69	127.17	120.48
1	A	138	PRO	N-CA-CB	8.22	111.65	102.60
1	A	389	ILE	CA-CB-CG1	-8.15	96.54	110.40
1	A	194	THR	CA-C-N	7.96	132.00	120.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	194	THR	C-N-CA	7.96	132.00	120.38
1	A	342	VAL	N-CA-CB	-7.94	95.06	110.78
1	A	322	GLU	CA-C-O	-7.86	109.98	119.49
1	A	59	MET	CA-C-O	-7.78	112.58	120.90
1	A	229	VAL	O-C-N	-7.71	116.07	121.55
1	A	249	ASP	CA-C-O	-7.34	112.54	121.28
1	A	358	LEU	O-C-N	-7.13	113.88	122.65
1	A	368	GLY	CA-C-N	-7.10	113.08	122.94
1	A	368	GLY	C-N-CA	-7.10	113.08	122.94
1	A	291	GLY	CA-C-N	-6.90	114.63	123.19
1	A	291	GLY	C-N-CA	-6.90	114.63	123.19
1	A	285	GLU	CB-CG-CD	6.86	124.26	112.60
1	A	70	MET	CA-C-O	6.81	128.65	120.28
1	A	57	GLN	OE1-CD-NE2	-6.80	115.80	122.60
1	A	284	VAL	N-CA-CB	6.71	120.65	110.58
1	A	128	ILE	CA-C-O	-6.70	113.36	120.53
1	A	384	LEU	CA-C-O	6.61	128.58	121.11
1	A	301	ILE	CA-C-O	-6.52	113.79	120.25
1	A	184	ALA	O-C-N	-6.49	114.89	122.87
1	A	214	LEU	N-CA-C	6.43	121.26	113.41
1	A	165	PHE	CA-C-O	-6.39	109.94	120.80
1	A	151	ARG	CA-C-N	6.39	126.31	119.28
1	A	151	ARG	C-N-CA	6.39	126.31	119.28
1	A	9	LYS	O-C-N	-6.36	114.90	122.15
1	A	344	PHE	CA-CB-CG	6.35	120.15	113.80
1	A	165	PHE	CA-C-N	-6.31	106.76	121.52
1	A	165	PHE	C-N-CA	-6.31	106.76	121.52
1	A	54	GLU	CA-CB-CG	6.28	126.66	114.10
1	A	47	GLU	CB-CG-CD	-6.25	101.97	112.60
1	A	360	THR	CA-C-O	-6.24	114.25	121.49
1	A	165	PHE	O-C-N	6.20	132.91	123.00
1	A	184	ALA	CA-C-N	6.19	131.55	122.39
1	A	184	ALA	C-N-CA	6.19	131.55	122.39
1	A	68	ARG	NE-CZ-NH1	-6.19	115.31	121.50
1	A	113	ALA	CA-C-O	6.17	127.09	120.55
1	A	145	THR	CA-C-N	6.12	128.77	120.44
1	A	145	THR	C-N-CA	6.12	128.77	120.44
1	A	241	TRP	CA-C-O	-6.10	114.18	120.89
1	A	154	VAL	O-C-N	-6.03	116.24	122.63
1	A	92	ASP	O-C-N	6.01	129.00	122.15
1	A	188	VAL	CA-C-O	-6.01	114.10	120.53
1	A	376	GLY	N-CA-C	-5.98	99.00	113.18

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	93	ALA	O-C-N	-5.92	115.97	122.07
1	A	217	ASP	CA-C-O	-5.91	114.45	121.66
1	A	25	ALA	O-C-N	-5.89	116.05	123.17
1	A	128	ILE	O-C-N	5.86	129.96	123.10
1	A	366	GLU	N-CA-C	5.86	118.79	111.24
1	A	347	ASP	CA-C-N	5.83	128.09	120.28
1	A	347	ASP	C-N-CA	5.83	128.09	120.28
1	A	171	LEU	O-C-N	5.79	128.75	122.15
1	A	173	LEU	CA-C-N	5.78	128.30	120.38
1	A	173	LEU	C-N-CA	5.78	128.30	120.38
1	A	240	VAL	CB-CA-C	-5.77	105.76	111.30
1	A	325	ASN	O-C-N	-5.74	116.16	122.07
1	A	172	ARG	O-C-N	-5.73	116.05	122.12
1	A	337	MET	N-CA-CB	-5.70	101.91	110.91
1	A	60	CYS	N-CA-CB	5.69	118.26	110.01
1	A	172	ARG	CA-C-O	5.65	126.54	120.55
1	A	161	GLN	N-CA-C	5.64	119.18	111.39
1	A	270	ASP	CA-C-N	5.64	125.84	120.43
1	A	270	ASP	C-N-CA	5.64	125.84	120.43
1	A	384	LEU	CA-C-N	-5.63	114.50	122.94
1	A	384	LEU	C-N-CA	-5.63	114.50	122.94
1	A	255	ASP	CA-CB-CG	5.59	118.19	112.60
1	A	24	THR	CA-C-O	-5.59	114.90	121.44
1	A	138	PRO	CA-N-CD	-5.58	103.69	111.50
1	A	38	ASP	N-CA-CB	5.57	118.15	110.07
1	A	96	ASP	CA-C-O	-5.54	113.60	119.97
1	A	306	GLY	N-CA-C	-5.54	101.04	112.34
1	A	229	VAL	CA-C-O	5.53	122.21	119.12
1	A	77	LEU	CA-C-O	-5.51	113.63	119.97
1	A	315	GLN	CA-CB-CG	5.51	125.13	114.10
1	A	93	ALA	CA-C-N	5.51	127.93	120.38
1	A	93	ALA	C-N-CA	5.51	127.93	120.38
1	A	151	ARG	CD-NE-CZ	-5.49	116.72	124.40
1	A	151	ARG	NE-CZ-NH2	-5.48	114.27	119.20
1	A	57	GLN	CG-CD-NE2	5.47	124.61	116.40
1	A	306	GLY	CA-C-O	-5.44	113.75	121.52
1	A	364	GLY	CA-C-O	5.41	125.13	119.07
1	A	342	VAL	O-C-N	-5.41	115.90	122.06
1	A	13	ALA	CA-C-O	-5.37	115.82	121.88
1	A	32	GLU	CA-C-O	-5.36	115.16	121.16
1	A	83	VAL	O-C-N	5.35	129.16	121.82
1	A	83	VAL	CA-C-O	-5.35	114.47	120.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	207	ASP	CA-CB-CG	5.35	117.95	112.60
1	A	305	GLY	CA-C-N	5.35	130.27	121.87
1	A	305	GLY	C-N-CA	5.35	130.27	121.87
1	A	366	GLU	CA-C-N	5.34	130.61	122.65
1	A	366	GLU	C-N-CA	5.34	130.61	122.65
1	A	389	ILE	CA-CB-CG2	5.31	119.53	110.50
1	A	137	MET	O-C-N	5.30	126.91	121.72
1	A	257	HIS	CA-CB-CG	-5.29	108.50	113.80
1	A	15	GLY	CA-C-N	5.27	125.03	119.76
1	A	15	GLY	C-N-CA	5.27	125.03	119.76
1	A	76	ILE	O-C-N	-5.26	116.75	121.91
1	A	47	GLU	CA-C-O	-5.26	114.16	120.10
1	A	55	LYS	CA-C-O	-5.24	115.29	120.90
1	A	114	ILE	CA-C-O	-5.24	115.62	121.17
1	A	186	VAL	CA-C-O	-5.21	114.99	120.57
1	A	330	VAL	CA-C-O	-5.21	115.27	121.05
1	A	388	ALA	N-CA-CB	5.19	117.60	109.97
1	A	276	SER	CA-C-N	5.17	127.93	120.38
1	A	276	SER	C-N-CA	5.17	127.93	120.38
1	A	55	LYS	O-C-N	5.16	127.45	122.09
1	A	205	HIS	CA-C-N	5.16	127.70	120.28
1	A	205	HIS	C-N-CA	5.16	127.70	120.28
1	A	104	ARG	O-C-N	-5.14	116.78	122.07
1	A	143	GLN	OE1-CD-NE2	5.11	127.71	122.60
1	A	343	LEU	CA-C-O	5.11	125.92	120.24
1	A	371	PHE	CA-C-N	5.11	131.42	121.41
1	A	371	PHE	C-N-CA	5.11	131.42	121.41
1	A	312	GLN	CA-C-O	-5.11	114.10	119.97
1	A	292	ILE	CA-C-O	5.10	125.89	120.48
1	A	223	ILE	CB-CA-C	-5.09	103.54	110.77
1	A	193	VAL	CA-C-N	5.08	131.58	122.38
1	A	193	VAL	C-N-CA	5.08	131.58	122.38
1	A	228	PRO	CA-C-O	-5.07	115.68	121.56
1	A	152	PRO	CB-CA-C	-5.06	104.31	112.21
1	A	292	ILE	CA-CB-CG2	5.05	119.09	110.50
1	A	282	ALA	CA-C-O	5.05	126.13	120.82
1	A	190	CYS	CA-C-O	-5.05	114.77	120.32
1	A	240	VAL	O-C-N	5.04	127.32	122.07
1	A	290	LEU	CA-C-O	5.03	125.31	119.32
1	A	292	ILE	O-C-N	-5.03	117.90	123.18
1	A	316	LYS	CA-C-O	5.02	126.27	120.90
1	A	209	LEU	O-C-N	-5.02	116.80	122.12

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	223	ILE	CA-C-O	-5.02	115.17	120.39
1	A	127	LEU	CA-C-O	5.01	126.31	120.20
1	A	340	ALA	CA-C-N	5.01	127.30	120.54
1	A	340	ALA	C-N-CA	5.01	127.30	120.54
1	A	292	ILE	CB-CA-C	5.00	118.29	110.83

There are no chirality outliers.

All (11) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	188	VAL	Mainchain
1	A	218	GLY	Mainchain
1	A	229	VAL	Mainchain
1	A	289	PRO	Mainchain
1	A	305	GLY	Mainchain,Peptide
1	A	345	ILE	Mainchain
1	A	348	GLU	Mainchain
1	A	360	THR	Mainchain
1	A	378	THR	Mainchain
1	A	64[B]	MET	Mainchain

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	2987	0	3029	56	0
2	A	5	0	0	0	0
3	A	48	0	31	4	0
4	A	184	0	0	6	2
All	All	3224	0	3060	56	2

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.

All (56) 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:342:VAL:HG13	1:A:370:LEU:HD11	1.55	0.89
1:A:51:GLU:HG2	4:A:530:HOH:O	1.78	0.83
1:A:104:ARG:NH2	1:A:108:GLU:OE1	2.16	0.74
1:A:58:ARG:HB3	1:A:62:LYS:HZ1	1.53	0.74
1:A:58:ARG:HB3	1:A:62:LYS:NZ	2.02	0.73
1:A:38:ASP:O	1:A:42:LYS:HG3	1.88	0.72
1:A:42:LYS:HG2	1:A:47:GLU:OE1	1.91	0.70
1:A:42:LYS:HG2	1:A:47:GLU:CD	2.23	0.64
1:A:62:LYS:HE3	3:A:390:COA:P1A	2.38	0.64
1:A:267:LEU:HD21	3:A:390:COA:S1P	2.40	0.62
1:A:271:VAL:HB	1:A:272:PRO:HD3	1.82	0.60
1:A:50:THR:O	1:A:54:GLU:HG2	2.03	0.58
1:A:55:LYS:NZ	3:A:390:COA:O7A	2.38	0.56
1:A:82:ASN:ND2	4:A:534:HOH:O	2.38	0.56
1:A:102:VAL:HB	1:A:103:PRO:HD3	1.87	0.55
1:A:352:LYS:HD3	4:A:504:HOH:O	2.05	0.55
1:A:230:PRO:O	1:A:231:GLU:HB2	2.07	0.54
1:A:292:ILE:HD12	1:A:293:SER:N	2.23	0.54
1:A:126:HIS:HE1	4:A:446:HOH:O	1.90	0.54
1:A:292:ILE:HD12	1:A:292:ILE:C	2.34	0.52
1:A:348:GLU:OE2	1:A:352:LYS:HD2	2.10	0.52
1:A:38:ASP:OD2	1:A:53:LYS:NZ	2.34	0.51
1:A:14:GLU:HG3	1:A:227:ASP:OD2	2.11	0.50
1:A:301:ILE:HG22	1:A:342:VAL:HG22	1.97	0.47
1:A:197:THR:CG2	1:A:263:LEU:HD23	2.44	0.47
1:A:256:GLY:HA2	1:A:264:THR:O	2.15	0.47
1:A:20:LEU:HD23	1:A:235:PRO:HA	1.97	0.47
1:A:241:TRP:HH2	1:A:285:GLU:HG3	1.80	0.46
1:A:24:THR:HB	1:A:344:PHE:CZ	2.50	0.45
1:A:137:MET:HA	1:A:138:PRO:C	2.42	0.45
1:A:5:SER:O	1:A:9:LYS:HG3	2.17	0.45
1:A:62:LYS:HE3	3:A:390:COA:O1A	2.17	0.45
1:A:100:VAL:O	1:A:103:PRO:HD2	2.17	0.45
1:A:129:VAL:HG21	1:A:141:ASP:HA	1.99	0.44
1:A:197:THR:HG22	1:A:263:LEU:HD23	1.99	0.44
1:A:271:VAL:CB	1:A:272:PRO:HD3	2.48	0.44
1:A:389:ILE:HD12	1:A:389:ILE:HA	1.54	0.44
1:A:194:THR:HG23	1:A:217:ASP:OD1	2.18	0.44
1:A:29:ASN:HB3	1:A:70:MET:O	2.17	0.43
1:A:299:PHE:CZ	1:A:368:GLY:HA3	2.53	0.43
1:A:334:TYR:CB	1:A:337:MET:HE2	2.49	0.43
1:A:188:VAL:O	1:A:221:ALA:HA	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:316:LYS:HE2	4:A:505:HOH:O	2.18	0.43
1:A:36:TYR:N	1:A:37:PRO:CD	2.81	0.43
1:A:240:VAL:HG21	1:A:367:TRP:HZ3	1.84	0.43
1:A:58:ARG:O	1:A:62:LYS:HD3	2.19	0.42
1:A:87:MET:HE2	1:A:87:MET:HB2	1.83	0.42
1:A:241:TRP:CH2	1:A:285:GLU:HG3	2.54	0.42
1:A:151:ARG:HH11	1:A:151:ARG:HD2	1.62	0.42
1:A:125:THR:OG1	1:A:126:HIS:HD2	2.03	0.41
1:A:172:ARG:HA	1:A:242:THR:HG21	2.02	0.41
1:A:342:VAL:HB	4:A:410:HOH:O	2.19	0.41
1:A:286:ALA:HB1	1:A:383:VAL:CG2	2.51	0.41
1:A:95:GLN:OE1	1:A:134:GLY:HA2	2.20	0.41
1:A:377:LEU:C	1:A:377:LEU:HD23	2.46	0.41
1:A:302:ALA:O	1:A:303:HIS:C	2.63	0.40

All (2) 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)
4:A:400:HOH:O	4:A:506:HOH:O[4_557]	2.10	0.10
4:A:470:HOH:O	4:A:543:HOH:O[4_557]	2.17	0.03

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	385/387 (100%)	374 (97%)	10 (3%)	1 (0%)	36 34

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	90	SER

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	322/321 (100%)	302 (94%)	20 (6%)	16	12

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	SER
1	A	14	GLU
1	A	47	GLU
1	A	54	GLU
1	A	75	GLU
1	A	111	VAL
1	A	123	LYS
1	A	133	SER
1	A	151	ARG
1	A	231	GLU
1	A	284	VAL
1	A	285	GLU
1	A	292	ILE
1	A	314	GLU
1	A	320	LYS
1	A	342	VAL
1	A	352	LYS
1	A	355	GLN
1	A	359	LYS
1	A	389	ILE

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	45	ASN
1	A	48	HIS
1	A	126	HIS
1	A	181	ASN
1	A	212	GLN

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Mol	Chain	Res	Type
1	A	312	GLN
1	A	355	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is 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	CSD	A	164	1	4,7,8	10.94	3 (75%)	1,8,10	8.79	1 (100%)

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	CSD	A	164	1	-	0/2/6/8	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	164	CSD	OD1-SG	-21.52	1.28	1.47
1	A	164	CSD	CB-SG	-2.93	1.62	1.79
1	A	164	CSD	CA-N	-2.59	1.40	1.48

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	164	CSD	OD1-SG-CB	-8.79	89.42	105.60

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 ligands 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	$\# Z > 2$
2	SO4	A	391	-	4,4,4	0.78	0	6,6,6	0.65	0
3	COA	A	390	-	47,50,50	1.45	10 (21%)	69,75,75	2.30	17 (24%)

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	COA	A	390	-	-	10/48/64/64	0/3/3/3

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	390	COA	OAP-CAP	-4.02	1.35	1.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	390	COA	C3P-N4P	3.20	1.53	1.46
3	A	390	COA	C7P-N8P	2.78	1.52	1.46
3	A	390	COA	P3B-O7A	2.62	1.58	1.50
3	A	390	COA	P3B-O8A	-2.57	1.45	1.54
3	A	390	COA	P3B-O9A	-2.47	1.45	1.54
3	A	390	COA	P2A-O6A	-2.43	1.49	1.59
3	A	390	COA	C2A-N1A	2.32	1.38	1.33
3	A	390	COA	CCP-CBP	2.28	1.56	1.52
3	A	390	COA	C2A-N3A	-2.03	1.30	1.33

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	390	COA	CDP-CBP-CAP	8.56	123.36	108.77
3	A	390	COA	C7P-N8P-C9P	-6.85	110.24	122.55
3	A	390	COA	C3P-N4P-C5P	-5.86	111.92	122.82
3	A	390	COA	O9A-P3B-O7A	-4.84	91.96	110.83
3	A	390	COA	C6P-C5P-N4P	4.62	124.76	116.34
3	A	390	COA	O5P-C5P-N4P	-4.16	114.87	123.03
3	A	390	COA	O9P-C9P-N8P	3.82	131.05	122.98
3	A	390	COA	O4B-C4B-C3B	-3.71	97.08	104.92
3	A	390	COA	O8A-P3B-O3B	3.04	117.71	105.85
3	A	390	COA	OAP-CAP-CBP	-2.90	103.47	110.18
3	A	390	COA	C2B-C3B-C4B	2.81	108.16	103.24
3	A	390	COA	O9A-P3B-O3B	2.48	115.50	105.85
3	A	390	COA	P1A-O5B-C5B	2.44	135.31	121.35
3	A	390	COA	O5A-P2A-O3A	2.29	113.46	107.27
3	A	390	COA	CEP-CBP-CDP	-2.23	104.75	109.20
3	A	390	COA	O6A-CCP-CBP	-2.14	107.11	110.55
3	A	390	COA	O5B-C5B-C4B	2.12	116.20	108.99

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	390	COA	C5B-O5B-P1A-O1A
3	A	390	COA	S1P-C2P-C3P-N4P
3	A	390	COA	C3B-O3B-P3B-O7A
3	A	390	COA	C5P-C6P-C7P-N8P
3	A	390	COA	C3B-C4B-C5B-O5B
3	A	390	COA	C5B-O5B-P1A-O3A
3	A	390	COA	P1A-O3A-P2A-O5A

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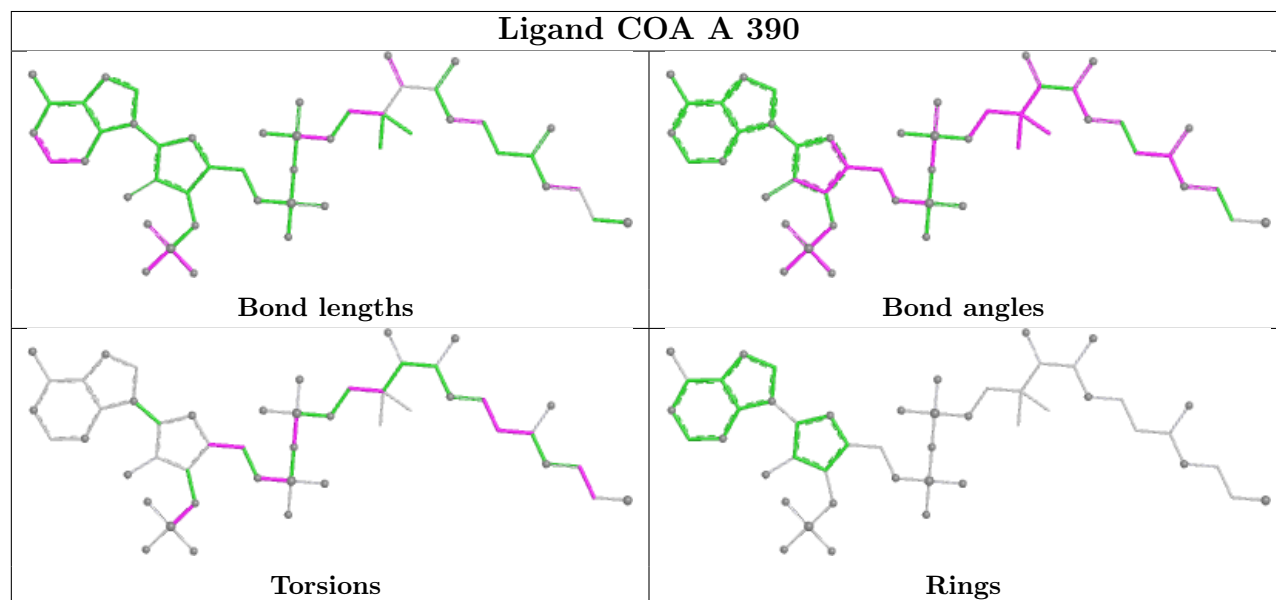
Mol	Chain	Res	Type	Atoms
3	A	390	COA	P1A-O3A-P2A-O4A
3	A	390	COA	CDP-CBP-CCP-O6A
3	A	390	COA	O5P-C5P-C6P-C7P

There are no ring outliers.

1 monomer is involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	390	COA	4	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 ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	386/387 (99%)	-0.30	1 (0%) 90 91	18, 32, 45, 59	1 (0%)

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	389	ILE	2.9

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	CSD	A	164	8/9	0.93	0.09	30,43,55,56	0

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.

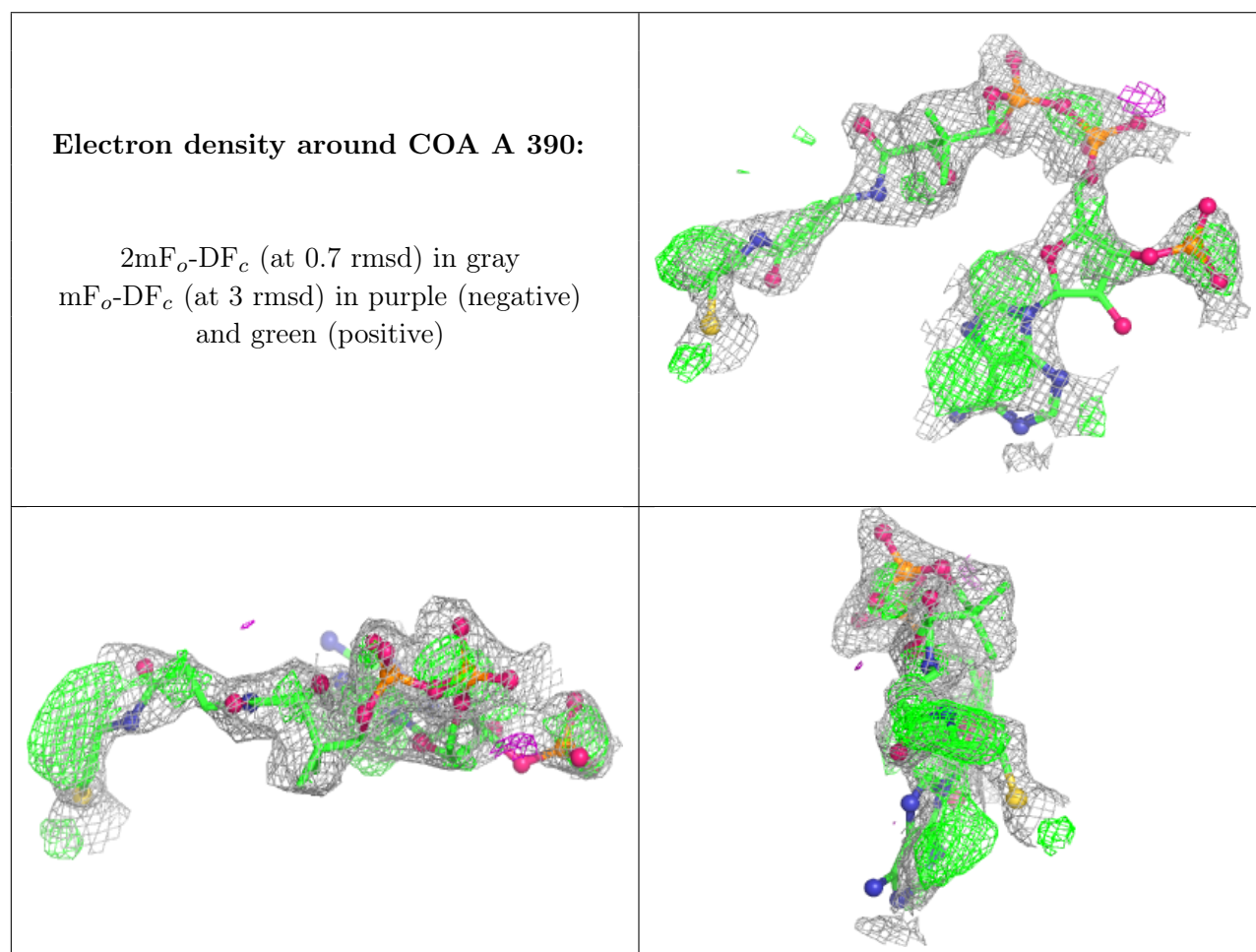
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	COA	A	390	48/48	0.76	0.23	24,40,64,65	48

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	SO4	A	391	5/5	0.91	0.10	52,53,53,55	5

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.



6.5 Other polymers ⓘ

There are no such residues in this entry.