



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:11 PM UTC

PDB ID : 1M1O / pdb_00001m1o
Title : Crystal structure of biosynthetic thiolase, C89A mutant, complexed with acetoacetyl-CoA
Authors : Kursula, P.; Ojala, J.; Lambeir, A.-M.; Wierenga, R.K.
Deposited on : 2002-06-20
Resolution : 1.95 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

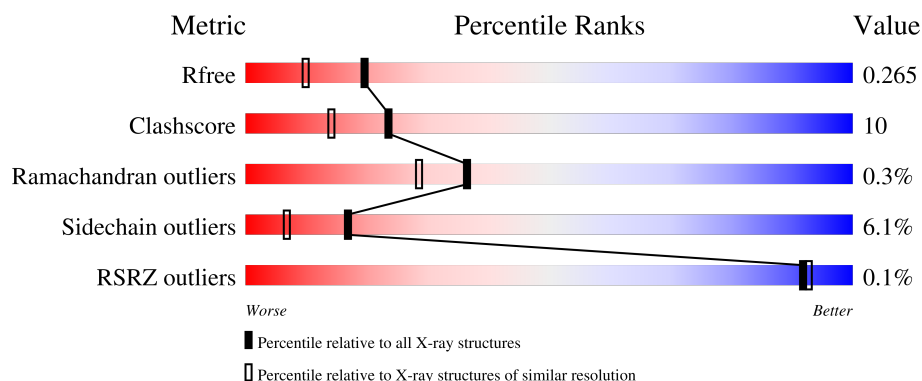
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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	3494 (1.96-1.96)
Clashscore	190562	3612 (1.96-1.96)
Ramachandran outliers	187476	3587 (1.96-1.96)
Sidechain outliers	187428	3587 (1.96-1.96)
RSRZ outliers	180081	3495 (1.96-1.96)

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	392	 84% 13% ..
1	B	392	 80% 18% ..
1	C	392	 76% 22% ..
1	D	392	 75% 22% ..

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 12414 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
1	A	390	Total	C	N	O	S	0	1	0
			2821	1752	510	539	20			
1	B	390	Total	C	N	O	S	0	1	0
			2821	1752	510	539	20			
1	C	390	Total	C	N	O	S	0	1	0
			2821	1752	510	539	20			
1	D	390	Total	C	N	O	S	0	1	0
			2821	1752	510	539	20			

There are 12 discrepancies between the modelled and reference sequences:

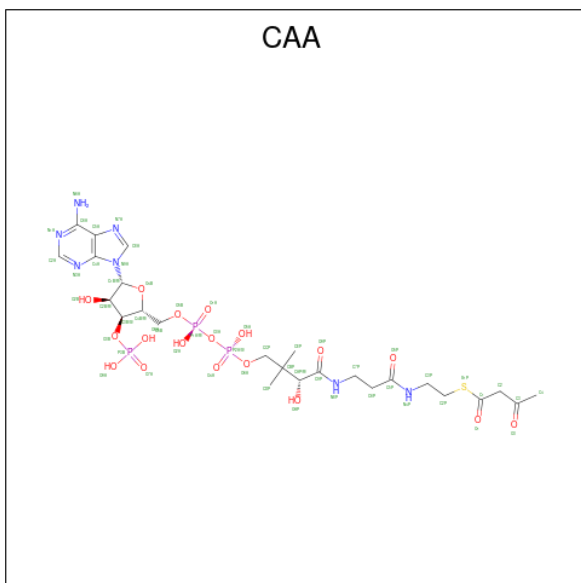
Chain	Residue	Modelled	Actual	Comment	Reference
A	10	ALA	-	insertion	UNP P07097
A	89	ALA	CYS	engineered mutation	UNP P07097
A	129	ARG	ALA	conflict	UNP P07097
B	10	ALA	-	insertion	UNP P07097
B	89	ALA	CYS	engineered mutation	UNP P07097
B	129	ARG	ALA	conflict	UNP P07097
C	10	ALA	-	insertion	UNP P07097
C	89	ALA	CYS	engineered mutation	UNP P07097
C	129	ARG	ALA	conflict	UNP P07097
D	10	ALA	-	insertion	UNP P07097
D	89	ALA	CYS	engineered mutation	UNP P07097
D	129	ARG	ALA	conflict	UNP P07097

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



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

- Molecule 3 is ACETOACETYL-COENZYME A (CCD ID: CAA) (formula: C₂₅H₄₀N₇O₁₈P₃S).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	S	0	0
			54	25	7	18	3	1		
3	B	1	Total	C	N	O	P	S	0	0
			54	25	7	18	3	1		

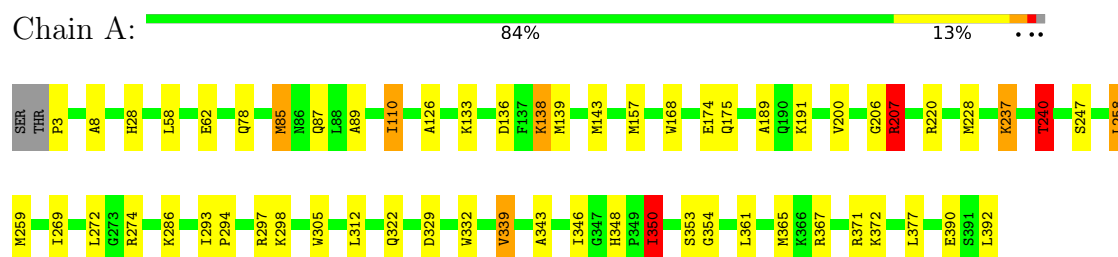
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	360	Total	O	0	0
			360	360		
4	B	372	Total	O	0	0
			372	372		
4	C	127	Total	O	0	0
			127	127		
4	D	143	Total	O	0	0
			143	143		

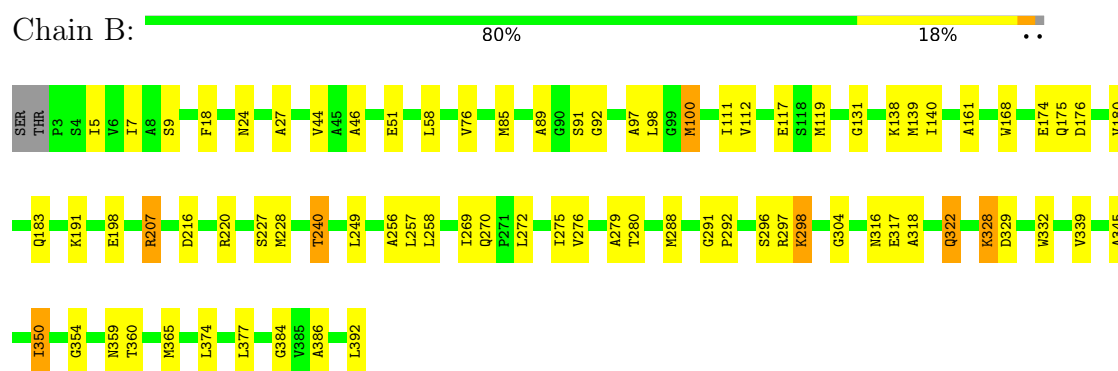
3 Residue-property plots

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.

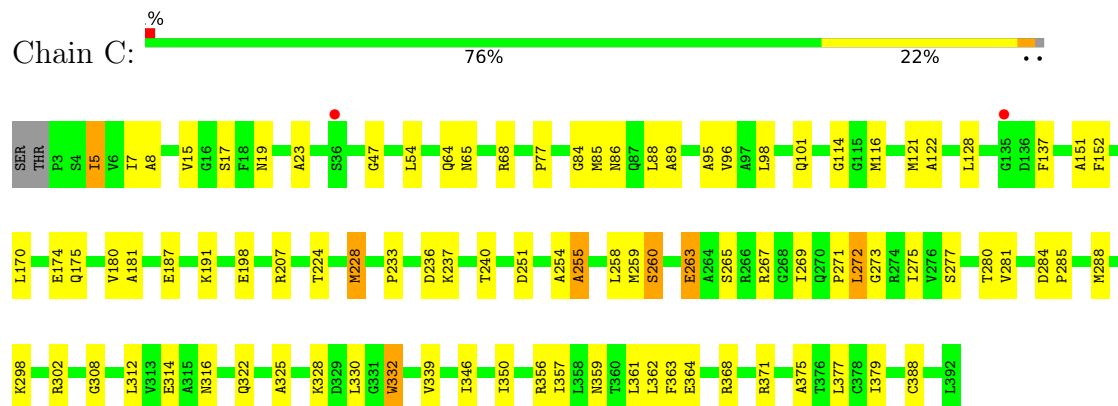
- Molecule 1: Acetyl-CoA acetyltransferase



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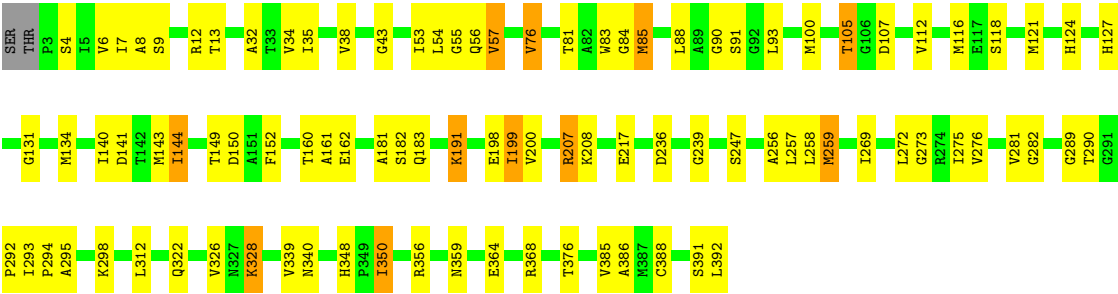
- Molecule 1: Acetyl-CoA acetyltransferase

Chain D:

75%

22%

..



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	84.22Å 79.22Å 148.29Å 90.00° 92.48° 90.00°	Depositor
Resolution (Å)	20.00 – 1.95 20.00 – 1.95	Depositor EDS
% Data completeness (in resolution range)	99.7 (20.00-1.95) 87.8 (20.00-1.95)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.32 (at 1.95Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.211 , 0.258 0.222 , 0.265	Depositor DCC
R_{free} test set	6373 reflections (4.50%)	wwPDB-VP
Wilson B-factor (Å ²)	20.6	Xtriage
Anisotropy	0.641	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 51.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.37$, $\langle L^2 \rangle = 0.19$	Xtriage
Estimated twinning fraction	0.166 for h,-k,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	12414	wwPDB-VP
Average B, all atoms (Å ²)	17.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.32% 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: CAA, SO4

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	1.31	9/2867 (0.3%)	1.19	6/3871 (0.2%)
1	B	1.26	12/2867 (0.4%)	1.21	7/3871 (0.2%)
1	C	0.86	5/2867 (0.2%)	0.97	2/3871 (0.1%)
1	D	0.82	6/2867 (0.2%)	0.97	1/3871 (0.0%)
All	All	1.09	32/11468 (0.3%)	1.09	16/15484 (0.1%)

All (32) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	126	ALA	CA-CB	-8.61	1.39	1.53
1	C	388	CYS	C-O	7.90	1.33	1.23
1	C	388	CYS	CA-C	7.88	1.62	1.52
1	C	260	SER	C-O	7.50	1.32	1.23
1	A	143	MET	SD-CE	-7.11	1.61	1.79
1	D	9	SER	CB-OG	7.06	1.56	1.42
1	D	388	CYS	C-O	6.87	1.32	1.24
1	C	308	GLY	C-O	6.25	1.31	1.23
1	B	386	ALA	CA-CB	-6.24	1.41	1.54
1	B	365	MET	SD-CE	-5.97	1.64	1.79
1	A	85	MET	CA-C	-5.92	1.45	1.52
1	D	259	MET	SD-CE	5.87	1.94	1.79
1	B	100	MET	SD-CE	5.84	1.94	1.79
1	B	97	ALA	C-O	-5.75	1.17	1.24
1	A	58	LEU	C-O	-5.72	1.18	1.24
1	D	105	THR	C-O	5.61	1.31	1.24
1	A	353	SER	C-O	-5.59	1.17	1.24
1	A	110	ILE	CA-C	-5.57	1.45	1.52
1	B	58	LEU	C-O	-5.57	1.17	1.23
1	B	46	ALA	CA-CB	5.54	1.62	1.53
1	B	216	ASP	CA-CB	-5.46	1.46	1.53
1	A	259	MET	SD-CE	5.40	1.93	1.79

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	27	ALA	CA-CB	-5.36	1.45	1.53
1	A	365	MET	SD-CE	-5.33	1.66	1.79
1	A	189	ALA	CA-CB	5.30	1.61	1.53
1	C	265	SER	C-O	5.29	1.30	1.24
1	B	76	VAL	N-CA	-5.21	1.40	1.47
1	B	112	VAL	CA-C	5.13	1.58	1.52
1	D	273	GLY	C-O	5.10	1.30	1.23
1	B	117	GLU	CA-C	-5.10	1.46	1.52
1	B	9	SER	C-O	5.08	1.29	1.23
1	D	107	ASP	CA-C	5.01	1.59	1.52

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	92	GLY	N-CA-C	-7.95	101.93	113.86
1	B	76	VAL	N-CA-C	-7.40	101.70	108.95
1	A	240	THR	CB-CA-C	6.50	117.93	109.28
1	D	76	VAL	N-CA-C	-6.29	102.78	108.95
1	B	91	SER	N-CA-C	6.28	117.79	111.07
1	C	350	ILE	N-CA-C	5.95	116.42	110.23
1	C	255	ALA	N-CA-C	5.78	116.18	108.38
1	A	343	ALA	N-CA-C	5.70	118.26	111.71
1	A	200	VAL	CA-C-O	5.50	122.77	119.19
1	A	339	VAL	N-CA-CB	-5.37	103.75	110.47
1	A	350	ILE	N-CA-C	5.23	120.23	109.34
1	A	240	THR	N-CA-CB	-5.22	102.93	111.18
1	B	304	GLY	N-CA-C	-5.11	107.76	115.32
1	B	207	ARG	N-CA-C	-5.11	105.89	112.68
1	B	98	LEU	CA-C-N	5.02	125.52	119.94
1	B	98	LEU	C-N-CA	5.02	125.52	119.94

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	2821	0	2828	41	0
1	B	2821	0	2828	57	0
1	C	2821	0	2828	64	0
1	D	2821	0	2828	77	0
2	A	10	0	0	0	0
2	B	10	0	0	0	0
3	A	54	0	36	6	0
3	B	54	0	36	4	0
4	A	360	0	0	19	0
4	B	372	0	0	22	0
4	C	127	0	0	23	0
4	D	143	0	0	39	0
All	All	12414	0	11384	232	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 10.

All (232) 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:296:SER:HB2	4:B:408:HOH:O	1.34	1.26
1:B:100:MET:HG3	4:B:466:HOH:O	1.49	1.12
1:B:258:LEU:HG	4:B:9882:HOH:O	1.55	1.06
3:A:1393:CAA:H4'3	4:A:467:HOH:O	1.56	1.03
1:A:89:ALA:HB2	3:A:1393:CAA:H2'2	1.44	0.98
1:B:269:ILE:HD11	4:B:473:HOH:O	1.67	0.94
1:B:89:ALA:HB2	3:B:2393:CAA:H2'2	1.50	0.93
1:A:258:LEU:HG	4:A:9763:HOH:O	1.66	0.93
1:D:207:ARG:HA	4:D:524:HOH:O	1.69	0.93
1:B:175:GLN:HE22	1:B:240:THR:CG2	1.85	0.90
1:B:175:GLN:HE22	1:B:240:THR:HG23	1.36	0.89
1:B:89:ALA:CB	3:B:2393:CAA:H2'2	2.04	0.88
1:A:258:LEU:CD2	4:A:9763:HOH:O	2.23	0.87
1:D:35:ILE:HD12	4:D:462:HOH:O	1.77	0.83
1:C:314:GLU:HB2	4:C:501:HOH:O	1.80	0.82
1:A:258:LEU:CG	4:A:9763:HOH:O	2.22	0.81
1:C:151:ALA:HB3	4:C:519:HOH:O	1.81	0.80
1:D:83:TRP:CZ2	4:D:489:HOH:O	2.33	0.80
1:A:286:LYS:HE3	4:A:9757:HOH:O	1.81	0.80
1:D:93:LEU:HA	4:D:454:HOH:O	1.81	0.80
1:B:227:SER:OG	4:B:9734:HOH:O	2.03	0.77
1:D:32:ALA:HA	4:D:462:HOH:O	1.84	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:157:MET:HG3	4:A:467:HOH:O	1.85	0.76
1:A:312:LEU:HD23	1:A:361:LEU:HD12	1.68	0.75
1:A:133:LYS:HA	4:A:401:HOH:O	1.85	0.75
1:B:175:GLN:NE2	1:B:240:THR:HG23	2.01	0.74
1:A:175:GLN:HE22	1:A:240:THR:CG2	2.01	0.73
3:A:1393:CAA:N6A	4:A:9827:HOH:O	2.16	0.73
1:D:83:TRP:HZ2	4:D:489:HOH:O	1.67	0.72
1:D:88:LEU:O	1:D:91:SER:OG	2.07	0.72
1:A:89:ALA:HB2	3:A:1393:CAA:C2	2.18	0.72
1:C:281:VAL:HA	4:C:458:HOH:O	1.90	0.72
1:D:91:SER:HB3	4:D:502:HOH:O	1.88	0.71
1:C:330:LEU:HD12	1:C:332:TRP:CH2	2.26	0.71
1:A:175:GLN:HE22	1:A:240:THR:HG23	1.55	0.70
1:C:15:VAL:HG22	4:C:480:HOH:O	1.91	0.70
1:C:5:ILE:HB	4:C:509:HOH:O	1.92	0.69
1:C:122:ALA:HA	4:C:431:HOH:O	1.93	0.69
1:A:371:ARG:HD2	4:A:454:HOH:O	1.92	0.68
1:D:385:VAL:HG22	4:D:510:HOH:O	1.94	0.68
1:C:272:LEU:O	1:C:362:LEU:HD22	1.95	0.66
1:C:330:LEU:HD12	1:C:332:TRP:CZ2	2.31	0.66
1:A:258:LEU:HD21	4:A:9763:HOH:O	1.90	0.66
1:D:12:ARG:NE	1:D:198:GLU:OE2	2.29	0.66
1:B:257:LEU:HD23	1:B:258:LEU:N	2.12	0.65
1:A:136:ASP:OD2	4:A:409:HOH:O	2.14	0.65
1:C:275:ILE:HB	4:C:509:HOH:O	1.97	0.65
1:B:288:MET:HE1	4:B:479:HOH:O	1.96	0.64
1:D:385:VAL:HG11	4:D:532:HOH:O	1.96	0.64
1:D:112:VAL:HG12	4:D:504:HOH:O	1.97	0.64
1:A:286:LYS:CE	4:A:9757:HOH:O	2.42	0.64
1:C:275:ILE:O	4:C:418:HOH:O	2.14	0.64
1:D:90:GLY:O	4:D:532:HOH:O	2.15	0.64
1:B:100:MET:HE2	4:B:466:HOH:O	1.98	0.63
1:D:312:LEU:HD13	1:D:368:ARG:HD2	1.79	0.63
4:C:504:HOH:O	1:D:140:ILE:HD13	1.99	0.63
1:D:183:GLN:HG2	4:D:461:HOH:O	1.98	0.63
1:C:375:ALA:HA	4:C:501:HOH:O	1.99	0.62
1:C:255:ALA:HB3	4:C:436:HOH:O	2.00	0.61
1:B:24:ASN:OD1	4:B:472:HOH:O	2.16	0.61
1:C:181:ALA:HB2	4:C:447:HOH:O	2.00	0.60
1:C:89:ALA:HB1	4:C:400:HOH:O	2.01	0.59
1:D:150:ASP:CG	4:D:521:HOH:O	2.45	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:161:ALA:HA	4:B:416:HOH:O	2.02	0.58
1:C:128:LEU:HD21	1:C:137:PHE:CE2	2.37	0.58
1:B:180:VAL:HG22	1:B:228:MET:HE2	1.84	0.58
1:B:228:MET:HE1	4:B:413:HOH:O	2.03	0.58
1:B:322:GLN:HB3	4:B:416:HOH:O	2.04	0.58
1:B:174:GLU:OE2	1:B:328:LYS:NZ	2.31	0.58
1:D:392:LEU:HD12	4:D:522:HOH:O	2.04	0.58
1:B:316:ASN:HB3	4:B:9738:HOH:O	2.05	0.57
1:D:124:HIS:HA	1:D:140:ILE:O	2.05	0.57
1:D:150:ASP:HB2	4:D:521:HOH:O	2.05	0.57
1:C:95:ALA:HB3	4:C:448:HOH:O	2.05	0.56
1:C:174:GLU:OE2	1:C:328:LYS:NZ	2.33	0.56
1:A:354:GLY:HA2	1:A:377:LEU:HD11	1.87	0.56
4:C:437:HOH:O	1:D:152:PHE:CE2	2.53	0.56
1:C:180:VAL:HG22	1:C:228:MET:HE2	1.87	0.56
1:C:263:GLU:O	1:C:267:ARG:HD2	2.06	0.56
1:C:8:ALA:HB1	1:C:269:ILE:HG21	1.88	0.55
1:C:64:GLN:O	1:C:65:ASN:C	2.50	0.55
1:D:326:VAL:HG23	4:D:456:HOH:O	2.06	0.55
1:A:87:GLN:C	4:A:444:HOH:O	2.50	0.54
1:C:356:ARG:HG3	4:C:480:HOH:O	2.06	0.54
1:C:330:LEU:CD1	1:C:332:TRP:CH2	2.89	0.54
1:B:168:TRP:CH2	1:B:329:ASP:HB2	2.42	0.54
1:C:312:LEU:HD23	1:C:361:LEU:HD12	1.89	0.54
1:C:7:ILE:N	1:C:273:GLY:O	2.27	0.54
1:A:175:GLN:NE2	1:A:240:THR:HG23	2.23	0.53
1:C:198:GLU:HB3	1:C:363:PHE:CD2	2.43	0.53
1:B:318:ALA:HB1	4:B:479:HOH:O	2.07	0.53
1:C:54:LEU:O	1:C:84:GLY:HA2	2.08	0.53
1:B:175:GLN:HE22	1:B:240:THR:HG21	1.70	0.52
1:C:68:ARG:HG3	1:D:152:PHE:HZ	1.74	0.52
1:A:168:TRP:CH2	1:A:329:ASP:HB2	2.44	0.52
1:B:297:ARG:NE	4:B:9829:HOH:O	2.41	0.52
1:D:340:ASN:ND2	1:D:364:GLU:OE1	2.36	0.52
1:B:374:LEU:C	1:B:374:LEU:HD23	2.36	0.51
1:D:289:GLY:O	1:D:292:PRO:HD2	2.10	0.51
1:D:56:GLN:OE1	1:D:116:MET:HE3	2.10	0.51
1:C:259:MET:HG3	1:C:260:SER:O	2.11	0.51
1:A:174:GLU:OE2	4:A:442:HOH:O	2.19	0.51
1:C:316:ASN:ND2	1:C:377:LEU:HD23	2.25	0.51
1:B:44:VAL:CG2	4:B:473:HOH:O	2.59	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:295:ALA:HA	4:D:503:HOH:O	2.10	0.50
1:A:157:MET:HA	1:A:157:MET:HE2	1.93	0.50
1:B:257:LEU:HD23	1:B:257:LEU:C	2.36	0.50
1:D:6:VAL:O	1:D:258:LEU:HD13	2.12	0.50
1:C:114:GLY:HA3	1:C:254:ALA:O	2.11	0.50
1:A:89:ALA:CB	3:A:1393:CAA:H2'2	2.29	0.50
1:A:297:ARG:NE	4:A:437:HOH:O	2.45	0.50
1:D:6:VAL:HG22	1:D:259:MET:O	2.11	0.50
1:A:8:ALA:HB1	1:A:269:ILE:HG21	1.94	0.49
1:D:272:LEU:C	4:D:530:HOH:O	2.54	0.49
1:C:96:VAL:HG23	4:C:448:HOH:O	2.12	0.49
1:D:236:ASP:HB3	1:D:239:GLY:HA3	1.95	0.49
1:B:119:MET:HE3	1:B:350:ILE:CD1	2.43	0.49
1:B:296:SER:CB	4:B:408:HOH:O	2.15	0.49
1:C:114:GLY:N	4:C:448:HOH:O	2.46	0.48
1:A:89:ALA:CB	3:A:1393:CAA:C2	2.90	0.48
1:B:7:ILE:HG23	1:B:256:ALA:HB1	1.95	0.48
1:D:275:ILE:HG21	4:D:535:HOH:O	2.13	0.48
1:D:257:LEU:HD23	1:D:258:LEU:N	2.28	0.48
1:D:76:VAL:HG23	4:D:531:HOH:O	2.13	0.48
1:D:6:VAL:O	1:D:258:LEU:CD1	2.62	0.48
1:D:217:GLU:HA	4:D:476:HOH:O	2.12	0.48
1:A:175:GLN:HE22	1:A:240:THR:HG21	1.76	0.48
1:C:95:ALA:HA	1:C:98:LEU:HD12	1.95	0.48
1:C:280:THR:HG23	1:D:81:THR:HG21	1.96	0.48
1:C:19:ASN:HB2	4:C:513:HOH:O	2.13	0.47
1:B:89:ALA:HB3	3:B:2393:CAA:H2'2	1.91	0.47
1:A:78:GLN:NE2	4:A:433:HOH:O	2.24	0.47
1:C:277:SER:HB2	4:D:528:HOH:O	2.13	0.47
1:B:318:ALA:CB	4:B:479:HOH:O	2.62	0.47
1:D:57:VAL:HG21	1:D:350:ILE:HG22	1.97	0.47
1:D:100:MET:C	1:D:100:MET:SD	2.98	0.47
1:C:54:LEU:HD13	1:C:116:MET:HE2	1.96	0.47
1:D:376:THR:HG21	4:D:496:HOH:O	2.14	0.47
1:B:270:GLN:NE2	1:B:392:LEU:OXT	2.48	0.46
1:C:152:PHE:CE2	4:C:519:HOH:O	2.56	0.46
1:D:55:GLY:CA	4:D:502:HOH:O	2.64	0.46
1:C:277:SER:HB2	1:C:302:ARG:HB3	1.98	0.46
1:C:269:ILE:O	1:C:271:PRO:HD3	2.15	0.46
1:D:181:ALA:HA	4:D:529:HOH:O	2.15	0.46
1:B:354:GLY:HA2	1:B:377:LEU:HD11	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:281:VAL:HG12	1:D:282:GLY:N	2.31	0.46
1:C:85:MET:HA	1:D:85:MET:HA	1.98	0.46
1:B:131:GLY:HA2	1:D:131:GLY:HA2	1.98	0.45
1:D:12:ARG:O	1:D:199:ILE:HA	2.17	0.45
1:D:257:LEU:HD23	1:D:257:LEU:C	2.41	0.45
1:B:257:LEU:C	1:B:257:LEU:CD2	2.90	0.45
1:D:7:ILE:HG23	1:D:256:ALA:HB1	1.98	0.45
1:D:160:THR:HG21	4:D:521:HOH:O	2.16	0.45
1:D:208:LYS:C	4:D:519:HOH:O	2.60	0.45
1:A:28:HIS:ND1	1:A:62:GLU:OE2	2.43	0.45
1:B:168:TRP:HH2	1:B:329:ASP:HB2	1.82	0.45
1:B:198:GLU:OE2	1:B:360:THR:OG1	2.32	0.45
1:B:280:THR:HA	1:B:384:GLY:O	2.17	0.45
1:D:118:SER:OG	1:D:121:MET:HB2	2.17	0.45
1:B:183:GLN:HA	1:B:345:ALA:HB2	1.99	0.45
1:B:279:ALA:CB	1:B:298:LYS:HB3	2.47	0.45
1:C:101:GLN:HG2	1:D:105:THR:HG21	1.98	0.45
1:D:247[B]:SER:OG	1:D:348:HIS:HB2	2.17	0.45
1:B:138:LYS:HB2	1:B:140:ILE:HD11	1.99	0.45
1:C:272:LEU:O	1:C:362:LEU:CD2	2.62	0.45
1:D:150:ASP:CB	4:D:521:HOH:O	2.64	0.45
1:D:386:ALA:N	4:D:510:HOH:O	2.50	0.45
1:C:121:MET:HA	1:D:127:HIS:CD2	2.52	0.44
1:D:144:ILE:O	1:D:149:THR:OG1	2.33	0.44
1:C:54:LEU:HD13	1:C:116:MET:SD	2.58	0.44
1:C:288:MET:HB2	1:C:379:ILE:O	2.17	0.44
1:C:5:ILE:N	1:C:5:ILE:HD13	2.33	0.44
1:B:44:VAL:HG23	4:B:473:HOH:O	2.17	0.44
1:D:83:TRP:HE3	1:D:84:GLY:O	2.00	0.44
1:D:281:VAL:HG23	4:D:503:HOH:O	2.17	0.44
1:C:88:LEU:HB3	4:C:443:HOH:O	2.18	0.44
1:A:206:GLY:O	1:A:207:ARG:C	2.61	0.44
1:A:3:PRO:N	4:A:9961:HOH:O	2.51	0.43
1:C:47:GLY:HA2	1:C:77:PRO:HG3	2.00	0.43
1:D:328:LYS:HG3	4:D:453:HOH:O	2.18	0.43
1:D:93:LEU:HB3	4:D:532:HOH:O	2.17	0.43
1:B:51:GLU:HB3	1:B:111:ILE:CD1	2.48	0.43
1:A:274:ARG:NH2	1:A:390:GLU:OE1	2.51	0.43
1:B:317:GLU:OE2	4:B:9724:HOH:O	2.21	0.43
1:A:85:MET:HA	1:B:85:MET:HA	2.00	0.43
1:D:161:ALA:HA	4:D:459:HOH:O	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:138:LYS:O	4:A:9793:HOH:O	2.22	0.43
1:D:141:ASP:OD1	1:D:143:MET:HB3	2.19	0.43
1:A:139:MET:HE1	1:B:139:MET:HE1	2.00	0.42
1:B:258:LEU:N	1:B:258:LEU:HD22	2.34	0.42
1:D:57:VAL:HG21	1:D:350:ILE:CG2	2.48	0.42
1:D:162:GLU:HG3	4:D:523:HOH:O	2.18	0.42
1:D:293:ILE:HB	1:D:294:PRO:CD	2.50	0.42
1:A:346:ILE:O	1:A:346:ILE:HG22	2.19	0.42
1:B:329:ASP:OD1	1:B:329:ASP:C	2.62	0.42
1:C:86:ASN:OD1	1:C:88:LEU:HD23	2.19	0.42
1:A:293:ILE:HB	1:A:294:PRO:CD	2.50	0.42
1:D:43:GLY:HA3	4:D:507:HOH:O	2.19	0.42
1:D:191:LYS:HA	4:D:431:HOH:O	2.19	0.42
1:B:18:PHE:HB2	1:B:249:LEU:O	2.19	0.42
1:B:131:GLY:HA2	1:D:131:GLY:CA	2.50	0.42
1:B:275:ILE:CG2	4:B:466:HOH:O	2.68	0.42
1:C:364:GLU:O	1:C:368:ARG:HG2	2.19	0.42
1:D:8:ALA:HB1	1:D:269:ILE:HG21	2.01	0.42
1:B:89:ALA:HB2	3:B:2393:CAA:C2	2.36	0.42
1:C:19:ASN:C	1:C:23:ALA:HB2	2.44	0.42
1:C:284:ASP:OD1	1:C:285:PRO:HD2	2.20	0.42
1:D:76:VAL:HB	4:D:396:HOH:O	2.20	0.41
1:C:356:ARG:NH2	1:C:357:ILE:HG22	2.35	0.41
1:C:7:ILE:O	1:C:272:LEU:N	2.54	0.41
1:C:346:ILE:HD13	1:C:356:ARG:NH1	2.35	0.41
1:C:54:LEU:HD13	1:C:116:MET:CE	2.50	0.41
1:C:89:ALA:CB	4:C:400:HOH:O	2.66	0.41
1:D:54:LEU:O	1:D:84:GLY:HA2	2.20	0.41
1:B:44:VAL:HG22	4:B:473:HOH:O	2.20	0.41
1:B:176:ASP:O	1:B:180:VAL:HG23	2.21	0.41
1:D:35:ILE:HG22	4:D:531:HOH:O	2.20	0.41
1:A:237:LYS:HD2	1:A:237:LYS:HA	1.87	0.41
1:A:247[B]:SER:OG	1:A:348:HIS:HB2	2.21	0.41
1:B:275:ILE:HG21	4:B:466:HOH:O	2.21	0.41
1:D:55:GLY:HA2	4:D:502:HOH:O	2.21	0.41
1:D:290:THR:O	1:D:294:PRO:HD2	2.21	0.41
1:A:305:TRP:CE2	1:A:372:LYS:HD3	2.55	0.41
1:C:170:LEU:HD11	1:C:325:ALA:HB2	2.02	0.41
1:C:175:GLN:HE22	1:C:240:THR:HG21	1.86	0.41
1:C:84:GLY:HA3	4:C:515:HOH:O	2.21	0.40
1:A:157:MET:CG	4:A:467:HOH:O	2.54	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:291:GLY:N	1:B:292:PRO:CD	2.83	0.40
1:A:350:ILE:HD13	1:A:350:ILE:HG21	1.91	0.40
1:D:34:VAL:O	1:D:38:VAL:HG13	2.20	0.40
1:C:233:PRO:HB2	1:C:236:ASP:O	2.21	0.40
1:D:88:LEU:O	1:D:91:SER:CB	2.70	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	389/392 (99%)	369 (95%)	18 (5%)	2 (0%)	24	16
1	B	389/392 (99%)	374 (96%)	14 (4%)	1 (0%)	36	28
1	C	389/392 (99%)	370 (95%)	19 (5%)	0	100	100
1	D	389/392 (99%)	366 (94%)	22 (6%)	1 (0%)	36	28
All	All	1556/1568 (99%)	1479 (95%)	73 (5%)	4 (0%)	36	28

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	207	ARG
1	A	350	ILE
1	D	350	ILE
1	B	350	ILE

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	277/278 (100%)	261 (94%)	16 (6%)	18	7
1	B	277/278 (100%)	264 (95%)	13 (5%)	23	12
1	C	277/278 (100%)	259 (94%)	18 (6%)	15	6
1	D	277/278 (100%)	257 (93%)	20 (7%)	13	4
All	All	1108/1112 (100%)	1041 (94%)	67 (6%)	17	7

All (67) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	110	ILE
1	A	138	LYS
1	A	191	LYS
1	A	207	ARG
1	A	220	ARG
1	A	228	MET
1	A	237	LYS
1	A	240	THR
1	A	258	LEU
1	A	272	LEU
1	A	298	LYS
1	A	322	GLN
1	A	332	TRP
1	A	339	VAL
1	A	367	ARG
1	A	392	LEU
1	B	5	ILE
1	B	191	LYS
1	B	207	ARG
1	B	220	ARG
1	B	240	THR
1	B	272	LEU
1	B	276	VAL
1	B	298	LYS
1	B	322	GLN
1	B	328	LYS
1	B	332	TRP
1	B	339	VAL
1	B	359	ASN
1	C	5	ILE

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Mol	Chain	Res	Type
1	C	17	SER
1	C	187	GLU
1	C	191	LYS
1	C	207	ARG
1	C	224	THR
1	C	228	MET
1	C	237	LYS
1	C	251	ASP
1	C	258	LEU
1	C	263	GLU
1	C	272	LEU
1	C	298	LYS
1	C	322	GLN
1	C	332	TRP
1	C	339	VAL
1	C	359	ASN
1	C	371	ARG
1	D	4	SER
1	D	13	THR
1	D	53	ILE
1	D	57	VAL
1	D	85	MET
1	D	134	MET
1	D	144	ILE
1	D	182	SER
1	D	191	LYS
1	D	199	ILE
1	D	200	VAL
1	D	207	ARG
1	D	276	VAL
1	D	298	LYS
1	D	322	GLN
1	D	328	LYS
1	D	339	VAL
1	D	356	ARG
1	D	359	ASN
1	D	391	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	78	GLN

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Mol	Chain	Res	Type
1	A	175	GLN
1	A	184	ASN
1	B	78	GLN
1	B	175	GLN
1	B	184	ASN
1	B	221	HIS
1	C	78	GLN
1	C	124	HIS
1	C	184	ASN
1	C	316	ASN
1	C	322	GLN
1	D	78	GLN
1	D	175	GLN
1	D	183	GLN
1	D	322	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](#)

6 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
3	CAA	B	2393	-	53,56,56	1.73	15 (28%)	77,83,83	2.38	16 (20%)
2	SO4	B	9721	-	4,4,4	0.28	0	6,6,6	0.44	0
2	SO4	A	9722	-	4,4,4	0.22	0	6,6,6	0.56	0
2	SO4	A	9720	-	4,4,4	0.25	0	6,6,6	0.56	0
3	CAA	A	1393	-	53,56,56	1.69	11 (20%)	77,83,83	2.19	16 (20%)
2	SO4	B	9719	-	4,4,4	0.23	0	6,6,6	0.25	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
3	CAA	A	1393	-	-	14/55/71/71	0/3/3/3
3	CAA	B	2393	-	-	12/55/71/71	0/3/3/3

All (26) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	2393	CAA	P3B-O9A	4.70	1.72	1.54
3	B	2393	CAA	P3B-O8A	4.52	1.71	1.54
3	A	1393	CAA	P3B-O8A	4.41	1.71	1.54
3	A	1393	CAA	P1A-O3A	4.26	1.64	1.59
3	A	1393	CAA	P3B-O9A	4.22	1.70	1.54
3	A	1393	CAA	P2A-O3A	4.00	1.63	1.59
3	B	2393	CAA	C5A-C4A	3.83	1.45	1.39
3	A	1393	CAA	P2A-O5A	3.57	1.71	1.55
3	B	2393	CAA	P2A-O5A	3.42	1.71	1.55
3	A	1393	CAA	C5A-C4A	3.39	1.45	1.39
3	B	2393	CAA	P1A-O2A	3.31	1.70	1.55
3	A	1393	CAA	P1A-O2A	3.25	1.70	1.55
3	A	1393	CAA	C5A-N7A	-2.90	1.33	1.39
3	B	2393	CAA	C1-S1P	2.75	1.82	1.76
3	B	2393	CAA	P1A-O3A	2.74	1.62	1.59
3	B	2393	CAA	C5A-C6A	2.73	1.48	1.41
3	A	1393	CAA	C4A-N9A	-2.72	1.32	1.37
3	B	2393	CAA	P2A-O3A	2.71	1.62	1.59
3	B	2393	CAA	C5A-N7A	-2.68	1.34	1.39
3	B	2393	CAA	O1-C1	-2.55	1.17	1.21
3	B	2393	CAA	C4A-N9A	-2.49	1.32	1.37
3	B	2393	CAA	C8A-N7A	2.48	1.36	1.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	2393	CAA	O3-C3	-2.41	1.13	1.21
3	A	1393	CAA	C1-S1P	2.39	1.81	1.76
3	A	1393	CAA	C5A-C6A	2.38	1.47	1.41
3	B	2393	CAA	P3B-O3B	2.21	1.63	1.59

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	2393	CAA	C2-C1-S1P	12.01	128.89	113.63
3	B	2393	CAA	O1-C1-C2	-10.28	106.27	123.40
3	A	1393	CAA	C2-C1-S1P	9.79	126.06	113.63
3	A	1393	CAA	O1-C1-C2	-7.73	110.53	123.40
3	A	1393	CAA	N3A-C4A-N9A	5.32	136.22	127.17
3	B	2393	CAA	C5A-C4A-N3A	-4.97	119.87	126.72
3	A	1393	CAA	C3-C2-C1	4.93	133.29	113.95
3	A	1393	CAA	C5A-C4A-N3A	-4.67	120.29	126.72
3	A	1393	CAA	C4A-N9A-C8A	4.48	110.44	105.74
3	B	2393	CAA	N3A-C4A-N9A	4.43	134.70	127.17
3	A	1393	CAA	N3A-C2A-N1A	-3.85	122.75	128.58
3	B	2393	CAA	N3A-C2A-N1A	-3.67	123.02	128.58
3	B	2393	CAA	C2A-N3A-C4A	3.42	120.18	111.83
3	A	1393	CAA	C2A-N3A-C4A	3.40	120.15	111.83
3	B	2393	CAA	C4A-N9A-C8A	3.27	109.17	105.74
3	B	2393	CAA	C4A-N9A-C1B	-3.26	119.02	126.63
3	A	1393	CAA	C2P-S1P-C1	-3.25	92.23	101.84
3	B	2393	CAA	O4B-C1B-N9A	3.19	114.21	108.09
3	A	1393	CAA	O3-C3-C2	-3.11	108.82	120.91
3	B	2393	CAA	C4-C3-C2	3.04	128.31	117.78
3	B	2393	CAA	C3-C2-C1	2.95	125.53	113.95
3	A	1393	CAA	C4A-N9A-C1B	-2.68	120.36	126.63
3	A	1393	CAA	N9A-C8A-N7A	-2.59	110.26	113.94
3	A	1393	CAA	C4-C3-C2	2.57	126.66	117.78
3	B	2393	CAA	C2A-N1A-C6A	2.48	122.81	118.73
3	B	2393	CAA	N9A-C8A-N7A	-2.42	110.50	113.94
3	B	2393	CAA	O3-C3-C2	-2.37	111.71	120.91
3	B	2393	CAA	C4B-O4B-C1B	-2.30	104.39	109.47
3	A	1393	CAA	C2A-N1A-C6A	2.24	122.41	118.73
3	A	1393	CAA	C5A-C4A-N9A	-2.13	103.49	105.81
3	B	2393	CAA	C1B-N9A-C8A	2.11	131.77	127.09
3	A	1393	CAA	N6A-C6A-N1A	2.09	123.02	118.38

There are no chirality outliers.

All (26) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1393	CAA	C5B-O5B-P1A-O1A
3	A	1393	CAA	CDP-CBP-CCP-O6A
3	A	1393	CAA	CEP-CBP-CCP-O6A
3	A	1393	CAA	CAP-CBP-CCP-O6A
3	A	1393	CAA	O1-C1-S1P-C2P
3	A	1393	CAA	C2-C1-S1P-C2P
3	A	1393	CAA	S1P-C1-C2-C3
3	B	2393	CAA	CCP-O6A-P2A-O3A
3	B	2393	CAA	CCP-O6A-P2A-O4A
3	B	2393	CAA	S1P-C2P-C3P-N4P
3	B	2393	CAA	C3P-C2P-S1P-C1
3	B	2393	CAA	O1-C1-S1P-C2P
3	B	2393	CAA	C2-C1-S1P-C2P
3	B	2393	CAA	C1-C2-C3-O3
3	B	2393	CAA	C1-C2-C3-C4
3	A	1393	CAA	S1P-C2P-C3P-N4P
3	A	1393	CAA	C1-C2-C3-O3
3	A	1393	CAA	C1-C2-C3-C4
3	B	2393	CAA	C2P-C3P-N4P-C5P
3	A	1393	CAA	C5B-O5B-P1A-O3A
3	A	1393	CAA	O1-C1-C2-C3
3	B	2393	CAA	O1-C1-C2-C3
3	A	1393	CAA	P1A-O3A-P2A-O5A
3	B	2393	CAA	C3B-O3B-P3B-O8A
3	A	1393	CAA	P1A-O3A-P2A-O4A
3	B	2393	CAA	C6P-C7P-N8P-C9P

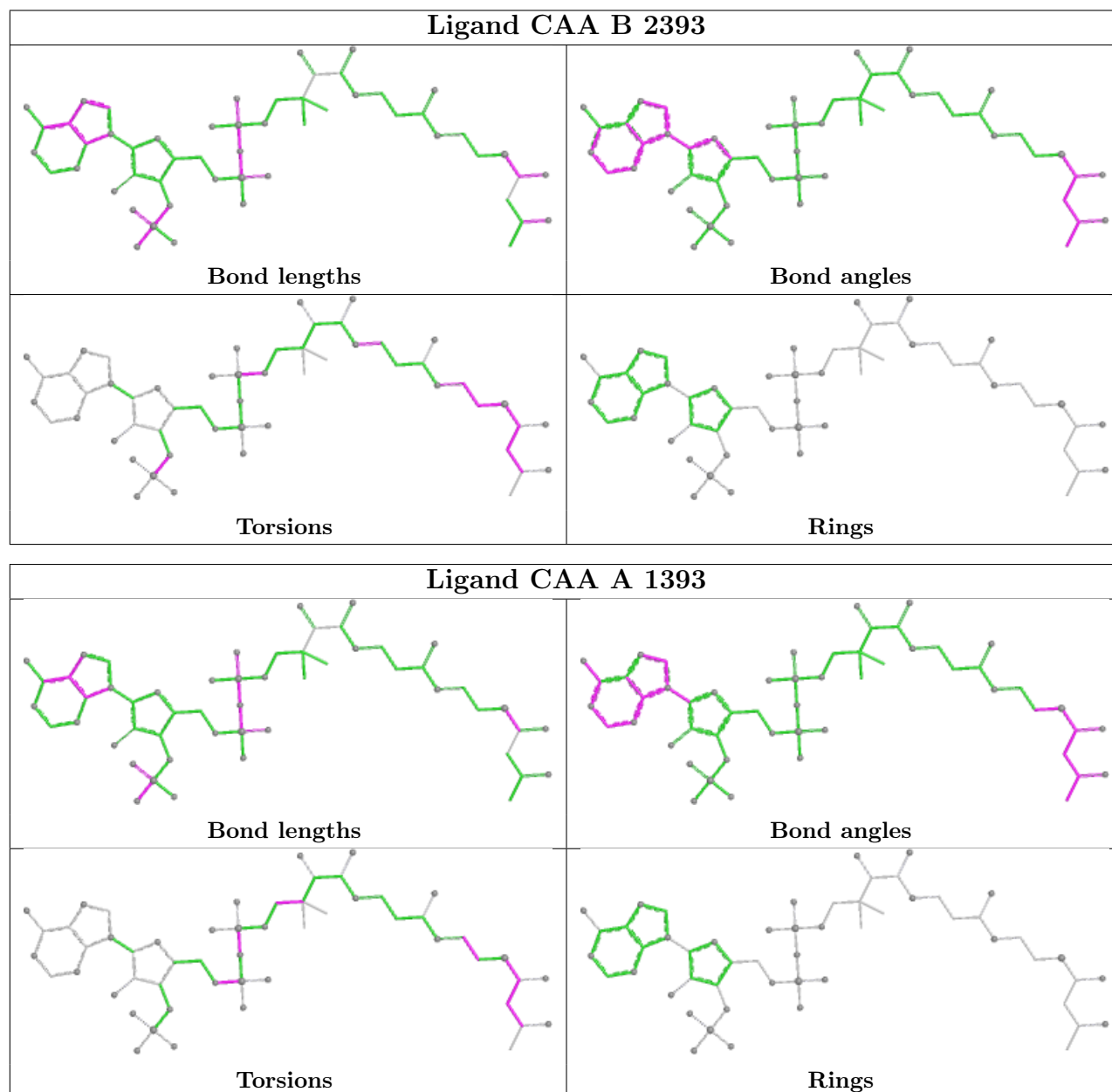
There are no ring outliers.

2 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	2393	CAA	4	0
3	A	1393	CAA	6	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%)	-1.18	0	100	100	6, 13, 27, 45	1 (0%)
1	B	390/392 (99%)	-1.18	0	100	100	6, 13, 26, 49	1 (0%)
1	C	390/392 (99%)	-0.81	2 (0%)	87	90	3, 14, 24, 38	1 (0%)
1	D	390/392 (99%)	-0.74	0	100	100	2, 14, 24, 41	1 (0%)
All	All	1560/1568 (99%)	-0.98	2 (0%)	92	93	2, 14, 25, 49	4 (0%)

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	36	SER	2.8
1	C	135	GLY	2.4

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](#)

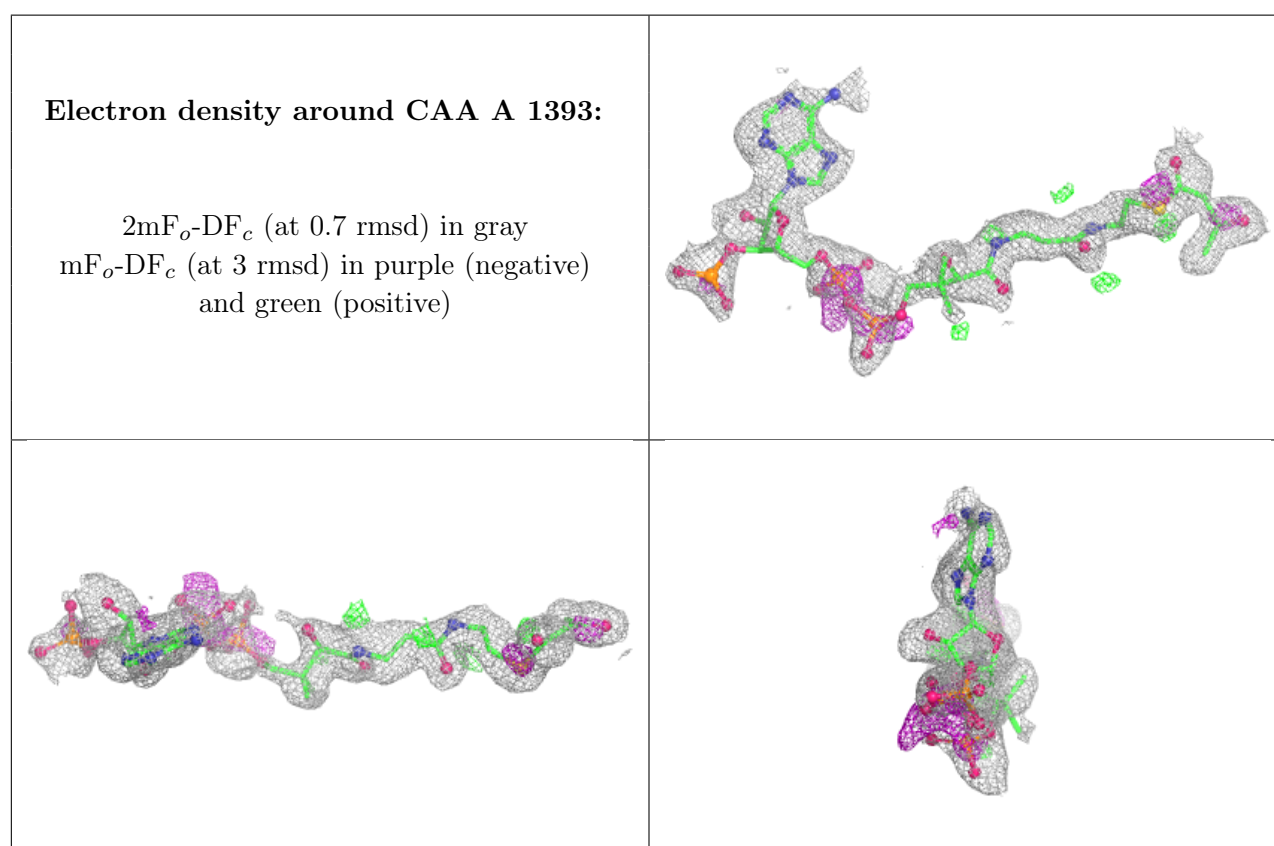
There are no oligosaccharides in this entry.

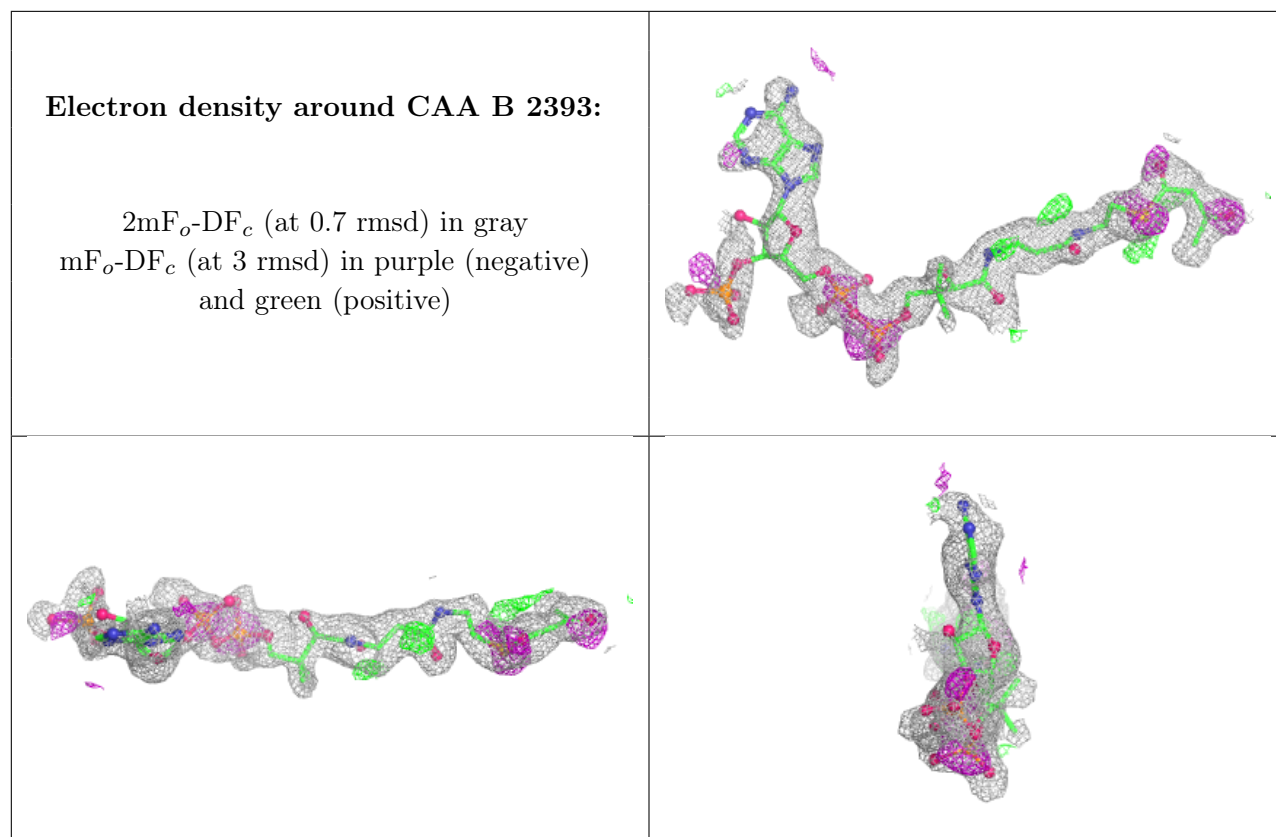
6.4 Ligands [i](#)

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($\text{\AA}^2$)	Q<0.9
3	CAA	A	1393	54/54	0.96	0.09	18,52,60,63	0
3	CAA	B	2393	54/54	0.96	0.09	11,52,57,60	0
2	SO4	A	9722	5/5	0.97	0.07	58,59,61,63	0
2	SO4	B	9719	5/5	0.97	0.07	66,67,68,68	0
2	SO4	A	9720	5/5	0.99	0.04	47,48,51,53	0
2	SO4	B	9721	5/5	0.99	0.05	47,47,50,51	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.





6.5 Other polymers [i](#)

There are no such residues in this entry.