



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 01:03 AM UTC

PDB ID : 4CKC / pdb_00004ckc
Title : Vaccinia virus capping enzyme complexed with SAH (monoclinic form)
Authors : Kyrieleis, O.J.P.; Chang, J.; de la Pena, M.; Shuman, S.; Cusack, S.
Deposited on : 2014-01-02
Resolution : 2.90 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

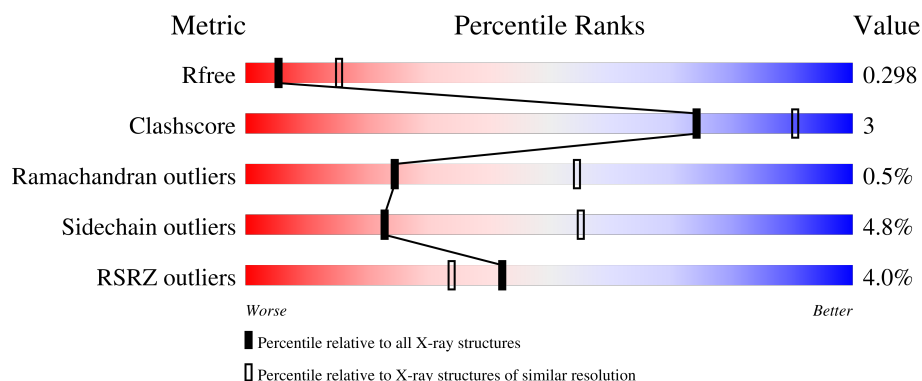
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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	844	6% 83% 13% ..
1	D	844	3% 85% 12% .
2	B	287	3% 86% 11% ..
2	E	287	% 84% 13% ..

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 18065 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.

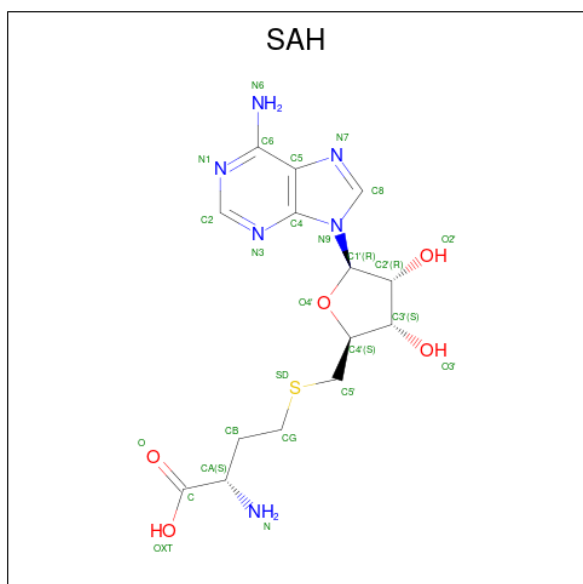
- Molecule 1 is a protein called MRNA-CAPPING ENZYME CATALYTIC SUBUNIT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	819	Total	C	N	O	S	0	1	0
			6640	4280	1091	1250	19			
1	D	821	Total	C	N	O	S	0	1	0
			6659	4293	1094	1253	19			

- Molecule 2 is a protein called MRNA-CAPPING ENZYME REGULATORY SUBUNIT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	284	Total	C	N	O	S	0	0	0
			2320	1492	385	430	13			
2	E	283	Total	C	N	O	S	0	0	0
			2312	1488	383	428	13			

- Molecule 3 is S-ADENOSYL-L-HOMOCYSTEINE (CCD ID: SAH) (formula: $C_{14}H_{20}N_6O_5S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			26	14	6	5	1		
3	D	1	Total	C	N	O	S	0	0
			26	14	6	5	1		

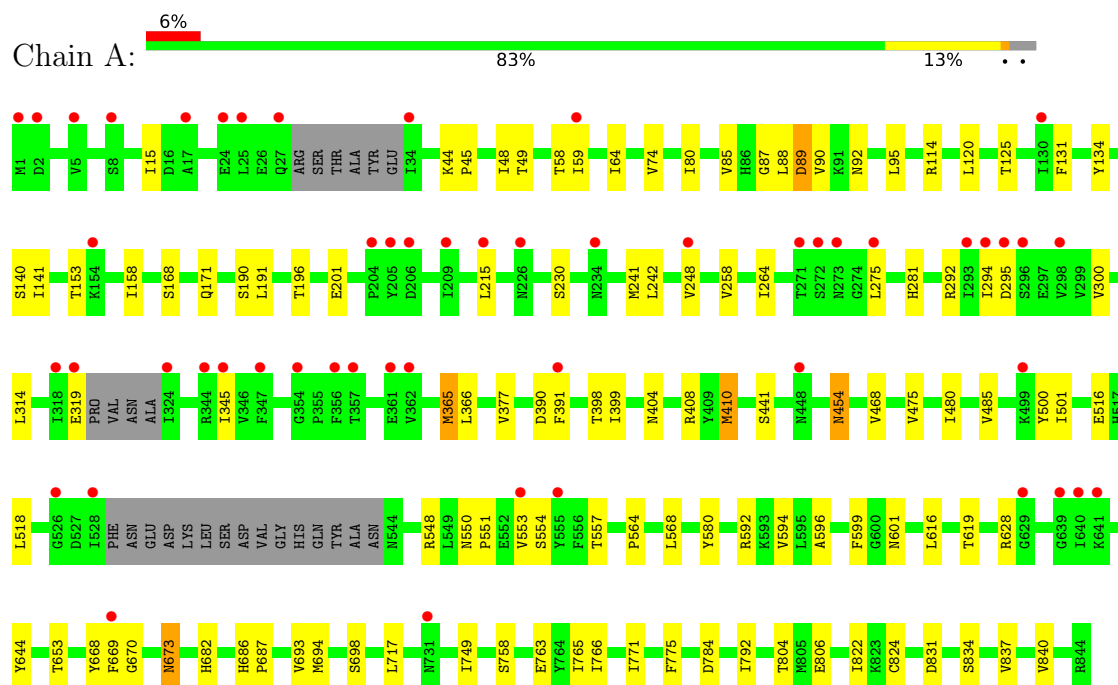
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	27	Total	O	0	0
			27	27		
4	B	5	Total	O	0	0
			5	5		
4	D	30	Total	O	0	0
			30	30		
4	E	20	Total	O	0	0
			20	20		

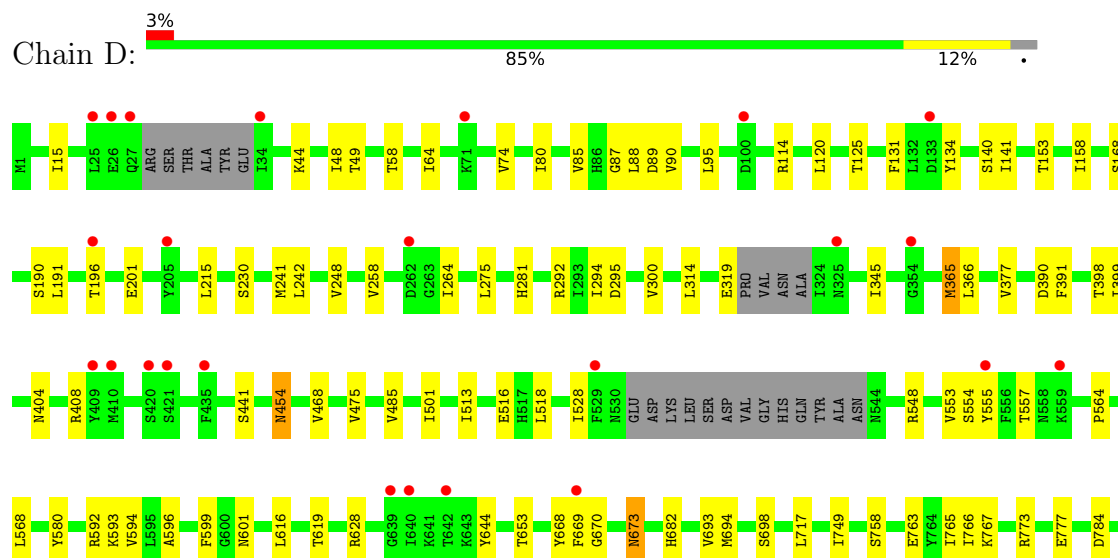
3 Residue-property plots [i](#)

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: MRNA-CAPPING ENZYME CATALYTIC SUBUNIT

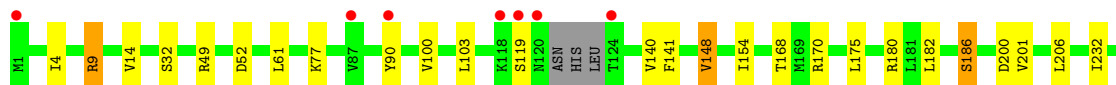
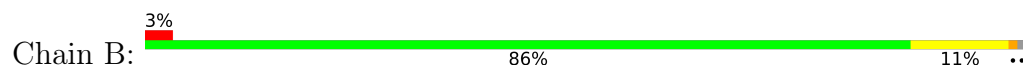


• Molecule 1: MRNA-CAPPING ENZYME CATALYTIC SUBUNIT

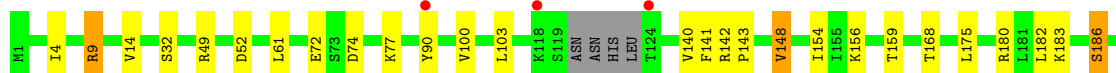
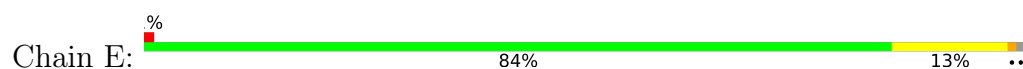




• Molecule 2: MRNA-CAPPING ENZYME REGULATORY SUBUNIT



• Molecule 2: MRNA-CAPPING ENZYME REGULATORY SUBUNIT



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	61.71Å 200.26Å 128.10Å 90.00° 102.18° 90.00°	Depositor
Resolution (Å)	125.00 – 2.90 125.00 – 2.90	Depositor EDS
% Data completeness (in resolution range)	93.5 (125.00-2.90) 93.6 (125.00-2.90)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.84 (at 2.90Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.255 , 0.297 0.256 , 0.298	Depositor DCC
R_{free} test set	3174 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	32.1	Xtriage
Anisotropy	0.881	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 38.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	0.000 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	18065	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 52.92 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 4.5373e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: SAH

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	0.43	0/6774	0.74	4/9157 (0.0%)
1	D	0.45	0/6794	0.74	3/9184 (0.0%)
2	B	0.44	0/2365	0.74	0/3189
2	E	0.46	0/2357	0.75	0/3178
All	All	0.44	0/18290	0.74	7/24708 (0.0%)

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	44	LYS	CA-C-N	5.64	126.19	120.38
1	A	44	LYS	C-N-CA	5.64	126.19	120.38
1	D	44	LYS	CA-C-N	5.49	126.03	120.38
1	D	44	LYS	C-N-CA	5.49	126.03	120.38
1	D	555	TYR	N-CA-C	5.45	116.91	111.07
1	A	45	PRO	CA-C-N	5.12	125.03	119.76
1	A	45	PRO	C-N-CA	5.12	125.03	119.76

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	6640	0	6723	44	1
1	D	6659	0	6738	41	1
2	B	2320	0	2363	12	2
2	E	2312	0	2357	16	1
3	A	26	0	19	0	0
3	D	26	0	19	0	0
4	A	27	0	0	1	1
4	B	5	0	0	1	0
4	D	30	0	0	3	0
4	E	20	0	0	1	0
All	All	18065	0	18219	113	3

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 3.

All (113) 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:410:MET:HE2	1:A:410:MET:HA	1.73	0.70
1:D:513:ILE:CD1	4:D:2014:HOH:O	2.41	0.69
1:D:513:ILE:HD12	4:D:2014:HOH:O	1.95	0.67
1:A:85:VAL:HG13	1:A:90:VAL:HG21	1.79	0.64
1:D:85:VAL:HG13	1:D:90:VAL:HG21	1.80	0.62
1:A:594:VAL:HB	1:A:616:LEU:HD12	1.82	0.61
1:A:158:ILE:HD11	1:A:215:LEU:HD11	1.82	0.60
2:B:263:LEU:HG	2:B:267:MET:HE3	1.84	0.60
1:D:594:VAL:HB	1:D:616:LEU:HD12	1.83	0.60
2:B:61:LEU:HD12	2:B:267:MET:HE1	1.83	0.60
2:E:61:LEU:HD12	2:E:267:MET:HE1	1.82	0.60
1:D:158:ILE:HD11	1:D:215:LEU:HD11	1.83	0.59
2:B:170:ARG:NH2	4:B:2003:HOH:O	2.36	0.59
2:E:263:LEU:HG	2:E:267:MET:HE3	1.87	0.57
2:B:52:ASP:HB3	2:B:264:TYR:CE1	2.41	0.56
1:A:365:MET:HE1	1:A:377:VAL:HG11	1.87	0.56
1:D:365:MET:HE1	1:D:377:VAL:HG11	1.87	0.56
2:E:52:ASP:HB3	2:E:264:TYR:CE1	2.41	0.55
2:E:268:THR:O	2:E:269:SER:C	2.51	0.54
1:A:59:ILE:HA	4:A:2001:HOH:O	2.06	0.53
1:D:580:TYR:O	1:D:592:ARG:NH1	2.43	0.52
1:A:749:ILE:HD11	1:A:765:ILE:HG12	1.90	0.52
1:D:485:VAL:HG11	1:D:518:LEU:HA	1.93	0.51
1:D:749:ILE:HD11	1:D:765:ILE:HG12	1.91	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:268:THR:O	2:B:269:SER:C	2.54	0.50
1:A:365:MET:HE2	1:A:365:MET:C	2.37	0.50
1:A:717:LEU:HD13	1:A:749:ILE:CD1	2.42	0.50
1:A:485:VAL:HG11	1:A:518:LEU:HA	1.94	0.49
1:D:717:LEU:HD13	1:D:749:ILE:CD1	2.42	0.49
1:A:580:TYR:O	1:A:592:ARG:NH1	2.46	0.49
1:D:767:LYS:HE2	4:D:2026:HOH:O	2.12	0.49
1:D:49:THR:HG21	1:D:191:LEU:HD22	1.95	0.48
1:D:682:HIS:HA	1:D:766:ILE:HD11	1.96	0.48
1:A:49:THR:HG21	1:A:191:LEU:HD22	1.95	0.48
1:D:365:MET:HE2	1:D:365:MET:C	2.39	0.47
2:B:9:ARG:NH2	2:B:285:LEU:O	2.45	0.47
1:A:454:ASN:OD1	1:A:454:ASN:N	2.48	0.46
1:A:601:ASN:OD1	1:A:628:ARG:NH1	2.48	0.46
1:A:831:ASP:O	1:A:834:SER:OG	2.25	0.46
1:D:454:ASN:OD1	1:D:454:ASN:N	2.47	0.45
1:D:64:ILE:HD12	1:D:95:LEU:HD22	1.97	0.45
1:A:131:PHE:CE2	1:A:134:TYR:HB3	2.51	0.45
1:D:131:PHE:CE2	1:D:134:TYR:HB3	2.51	0.45
2:E:9:ARG:NH2	2:E:285:LEU:O	2.47	0.45
2:B:141:PHE:CZ	2:B:168:THR:HG23	2.52	0.45
1:D:593:LYS:N	1:D:673:ASN:OD1	2.49	0.45
2:E:277:LYS:HA	2:E:280:ARG:HD2	1.99	0.45
1:A:64:ILE:HD12	1:A:95:LEU:HD22	1.98	0.44
1:D:784:ASP:HB3	1:D:840:VAL:HB	1.99	0.44
1:D:601:ASN:OD1	1:D:628:ARG:NH1	2.47	0.44
1:A:264:ILE:CD1	1:A:281:HIS:HB3	2.47	0.44
1:A:264:ILE:HD13	1:A:281:HIS:HB3	1.99	0.44
1:A:693:VAL:HG12	1:A:694:MET:HE2	1.99	0.43
2:E:141:PHE:CZ	2:E:168:THR:HG23	2.53	0.43
2:B:77:LYS:HG3	2:B:186:SER:OG	2.18	0.43
2:B:277:LYS:HA	2:B:280:ARG:HD2	2.00	0.43
1:D:264:ILE:HD13	1:D:281:HIS:HB3	2.00	0.43
1:A:241:MET:HE1	1:A:516:GLU:HG3	1.99	0.43
1:A:153:THR:HG21	1:A:215:LEU:HD13	2.01	0.43
1:D:264:ILE:CD1	1:D:281:HIS:HB3	2.48	0.43
1:D:300:VAL:HG11	1:D:314:LEU:CD2	2.49	0.43
2:B:4:ILE:HG21	2:B:254:GLU:HG2	2.01	0.43
2:E:159:THR:HB	4:E:2014:HOH:O	2.19	0.43
1:A:294:ILE:O	1:A:295:ASP:C	2.62	0.43
1:A:682:HIS:HA	1:A:766:ILE:HD11	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:300:VAL:HG11	1:A:314:LEU:CD2	2.49	0.43
2:E:4:ILE:HG21	2:E:254:GLU:HG2	2.01	0.42
1:A:141:ILE:O	1:A:168:SER:HA	2.18	0.42
1:A:242:LEU:HB2	1:A:391:PHE:HB3	2.01	0.42
1:D:141:ILE:O	1:D:168:SER:HA	2.18	0.42
1:D:669[A]:PHE:CG	1:D:670:GLY:N	2.87	0.42
2:B:175:LEU:HD22	2:B:206:LEU:HA	2.00	0.42
1:D:153:THR:HG21	1:D:215:LEU:HD13	2.02	0.42
2:E:175:LEU:HD22	2:E:206:LEU:HA	2.00	0.42
1:D:553:VAL:O	1:D:554:SER:C	2.61	0.42
1:D:693:VAL:HG12	1:D:694:MET:HE2	2.01	0.42
2:E:100:VAL:HA	2:E:154:ILE:O	2.19	0.42
1:D:242:LEU:HB2	1:D:391:PHE:HB3	2.02	0.42
1:D:399:ILE:HD11	1:D:485:VAL:HG13	2.02	0.42
1:A:92:ASN:HB2	1:A:171:GLN:O	2.19	0.42
1:D:241:MET:HE1	1:D:516:GLU:HG3	2.01	0.42
1:A:114:ARG:CB	1:A:120:LEU:HD23	2.50	0.42
1:A:399:ILE:HD11	1:A:485:VAL:HG13	2.02	0.42
1:A:480:ILE:HD11	1:A:500:TYR:HB2	2.02	0.42
1:A:550:ASN:N	1:A:551:PRO:HD3	2.35	0.41
1:A:673:ASN:OD1	1:A:673:ASN:N	2.50	0.41
1:D:275:LEU:HD23	1:D:345:ILE:HD12	2.02	0.41
2:E:72:GLU:OE2	2:E:191:LYS:NZ	2.43	0.41
1:A:365:MET:HE2	1:A:366:LEU:N	2.36	0.41
1:A:669[A]:PHE:CG	1:A:670:GLY:N	2.89	0.41
1:A:784:ASP:HB3	1:A:840:VAL:HB	2.02	0.41
2:E:74:ASP:OD2	2:E:156:LYS:NZ	2.45	0.41
1:D:114:ARG:CB	1:D:120:LEU:HD23	2.50	0.41
1:D:596:ALA:HB1	1:D:599:PHE:CD1	2.55	0.41
1:D:831:ASP:O	1:D:834:SER:OG	2.30	0.41
2:B:100:VAL:HA	2:B:154:ILE:O	2.20	0.41
1:D:564:PRO:O	1:D:568:LEU:HG	2.20	0.41
1:A:275:LEU:HD23	1:A:345:ILE:HD12	2.03	0.41
2:E:183:LYS:HG2	2:E:189:TRP:CZ2	2.56	0.41
2:E:142:ARG:N	2:E:143:PRO:CD	2.84	0.41
1:A:87:GLY:O	1:A:89:ASP:N	2.53	0.40
1:A:564:PRO:O	1:A:568:LEU:HG	2.21	0.40
1:A:792:ILE:CG2	1:A:822:ILE:HD11	2.51	0.40
1:D:87:GLY:O	1:D:89:ASP:N	2.54	0.40
1:D:365:MET:HE2	1:D:366:LEU:N	2.37	0.40
1:A:553:VAL:O	1:A:554:SER:C	2.63	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:596:ALA:HB1	1:A:599:PHE:CD1	2.56	0.40
1:D:773:ARG:O	1:D:777:GLU:HG3	2.21	0.40
1:D:784:ASP:CB	1:D:840:VAL:HB	2.52	0.40
1:A:686:HIS:CG	1:A:687:PRO:HD2	2.57	0.40
1:D:294:ILE:O	1:D:295:ASP:C	2.64	0.40
2:E:77:LYS:HG3	2:E:186:SER:OG	2.21	0.40
1:A:771:ILE:HD11	1:A:775:PHE:CE1	2.56	0.40

All (3) 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)
2:B:272:LYS:NZ	4:A:2014:HOH:O[2_556]	2.00	0.20
1:D:668:TYR:O	2:E:180:ARG:NH1[1_455]	2.14	0.06
1:A:668:TYR:O	2:B:180:ARG:NH1[1_455]	2.18	0.02

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	812/844 (96%)	764 (94%)	45 (6%)	3 (0%)	30	59
1	D	814/844 (96%)	767 (94%)	45 (6%)	2 (0%)	43	72
2	B	280/287 (98%)	263 (94%)	13 (5%)	4 (1%)	9	30
2	E	279/287 (97%)	263 (94%)	13 (5%)	3 (1%)	11	36
All	All	2185/2262 (97%)	2057 (94%)	116 (5%)	12 (0%)	24	54

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	119	SER

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Mol	Chain	Res	Type
1	A	88	LEU
1	A	248	VAL
1	D	88	LEU
1	D	248	VAL
2	B	200	ASP
2	B	269	SER
2	E	200	ASP
2	E	269	SER
1	A	89	ASP
2	B	148	VAL
2	E	148	VAL

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	754/774 (97%)	716 (95%)	38 (5%)	22	53
1	D	756/774 (98%)	719 (95%)	37 (5%)	22	54
2	B	269/272 (99%)	257 (96%)	12 (4%)	24	58
2	E	268/272 (98%)	256 (96%)	12 (4%)	24	58
All	All	2047/2092 (98%)	1948 (95%)	99 (5%)	23	55

All (99) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	15	ILE
1	A	48	ILE
1	A	58	THR
1	A	74	VAL
1	A	80	ILE
1	A	125	THR
1	A	140	SER
1	A	190	SER
1	A	196	THR
1	A	201	GLU

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Mol	Chain	Res	Type
1	A	230	SER
1	A	258	VAL
1	A	292	ARG
1	A	319	GLU
1	A	365	MET
1	A	390	ASP
1	A	398	THR
1	A	404	ASN
1	A	408	ARG
1	A	410	MET
1	A	441	SER
1	A	454	ASN
1	A	468	VAL
1	A	475	VAL
1	A	501	ILE
1	A	548	ARG
1	A	557	THR
1	A	619	THR
1	A	644	TYR
1	A	653	THR
1	A	673	ASN
1	A	698	SER
1	A	758	SER
1	A	763	GLU
1	A	804	THR
1	A	806	GLU
1	A	824	CYS
1	A	837	VAL
2	B	9	ARG
2	B	14	VAL
2	B	32	SER
2	B	49	ARG
2	B	90	TYR
2	B	103	LEU
2	B	140	VAL
2	B	148	VAL
2	B	182	LEU
2	B	186	SER
2	B	201	VAL
2	B	232	ILE
1	D	15	ILE
1	D	48	ILE

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Mol	Chain	Res	Type
1	D	58	THR
1	D	74	VAL
1	D	80	ILE
1	D	125	THR
1	D	140	SER
1	D	190	SER
1	D	196	THR
1	D	201	GLU
1	D	230	SER
1	D	258	VAL
1	D	292	ARG
1	D	319	GLU
1	D	365	MET
1	D	390	ASP
1	D	398	THR
1	D	404	ASN
1	D	408	ARG
1	D	441	SER
1	D	454	ASN
1	D	468	VAL
1	D	475	VAL
1	D	501	ILE
1	D	528	ILE
1	D	548	ARG
1	D	557	THR
1	D	619	THR
1	D	644	TYR
1	D	653	THR
1	D	673	ASN
1	D	698	SER
1	D	758	SER
1	D	763	GLU
1	D	806	GLU
1	D	824	CYS
1	D	837	VAL
2	E	9	ARG
2	E	14	VAL
2	E	32	SER
2	E	49	ARG
2	E	90	TYR
2	E	103	LEU
2	E	140	VAL

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Mol	Chain	Res	Type
2	E	148	VAL
2	E	182	LEU
2	E	186	SER
2	E	201	VAL
2	E	232	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	101	ASN
1	A	113	ASN
1	D	4	ASN
1	D	113	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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
3	SAH	A	1845	-	27,28,28	1.43	5 (18%)	36,40,40	2.08	9 (25%)
3	SAH	D	1845	-	27,28,28	1.46	5 (18%)	36,40,40	2.15	11 (30%)

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	SAH	A	1845	-	-	3/15/31/31	0/3/3/3
3	SAH	D	1845	-	-	4/15/31/31	0/3/3/3

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1845	SAH	C5-C4	4.52	1.47	1.39
3	D	1845	SAH	C5-C4	4.52	1.47	1.39
3	D	1845	SAH	C5-N7	-2.91	1.33	1.39
3	A	1845	SAH	C5-N7	-2.75	1.34	1.39
3	A	1845	SAH	C8-N7	2.50	1.36	1.31
3	D	1845	SAH	C5-C6	2.36	1.47	1.41
3	A	1845	SAH	OXT-C	-2.31	1.23	1.30
3	D	1845	SAH	C8-N7	2.28	1.36	1.31
3	A	1845	SAH	C5-C6	2.19	1.47	1.41
3	D	1845	SAH	OXT-C	-2.10	1.23	1.30

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1845	SAH	C5-C4-N3	-6.59	117.65	126.72
3	D	1845	SAH	C5-C4-N3	-6.39	117.92	126.72
3	A	1845	SAH	N3-C4-N9	5.27	136.13	127.17
3	D	1845	SAH	N3-C4-N9	5.24	136.08	127.17
3	D	1845	SAH	C2-N3-C4	4.09	121.83	111.83
3	A	1845	SAH	C2-N3-C4	4.05	121.72	111.83
3	D	1845	SAH	N3-C2-N1	-3.95	122.60	128.58
3	A	1845	SAH	N3-C2-N1	-3.52	123.26	128.58
3	D	1845	SAH	C4-C5-N7	-3.15	106.98	110.58
3	D	1845	SAH	OXT-C-O	-3.00	117.26	124.08
3	A	1845	SAH	C4-C5-N7	-2.97	107.18	110.58
3	A	1845	SAH	OXT-C-O	-2.72	117.91	124.08

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	1845	SAH	C5-N7-C8	2.69	107.69	103.45
3	A	1845	SAH	C5-N7-C8	2.42	107.25	103.45
3	D	1845	SAH	C4-N9-C8	2.35	108.21	105.74
3	A	1845	SAH	C4-N9-C8	2.20	108.05	105.74
3	D	1845	SAH	C2-N1-C6	2.17	122.30	118.73
3	D	1845	SAH	C1'-N9-C8	-2.12	122.38	127.09
3	D	1845	SAH	O4'-C1'-N9	2.02	111.97	108.09
3	A	1845	SAH	C1'-N9-C8	-2.02	122.62	127.09

There are no chirality outliers.

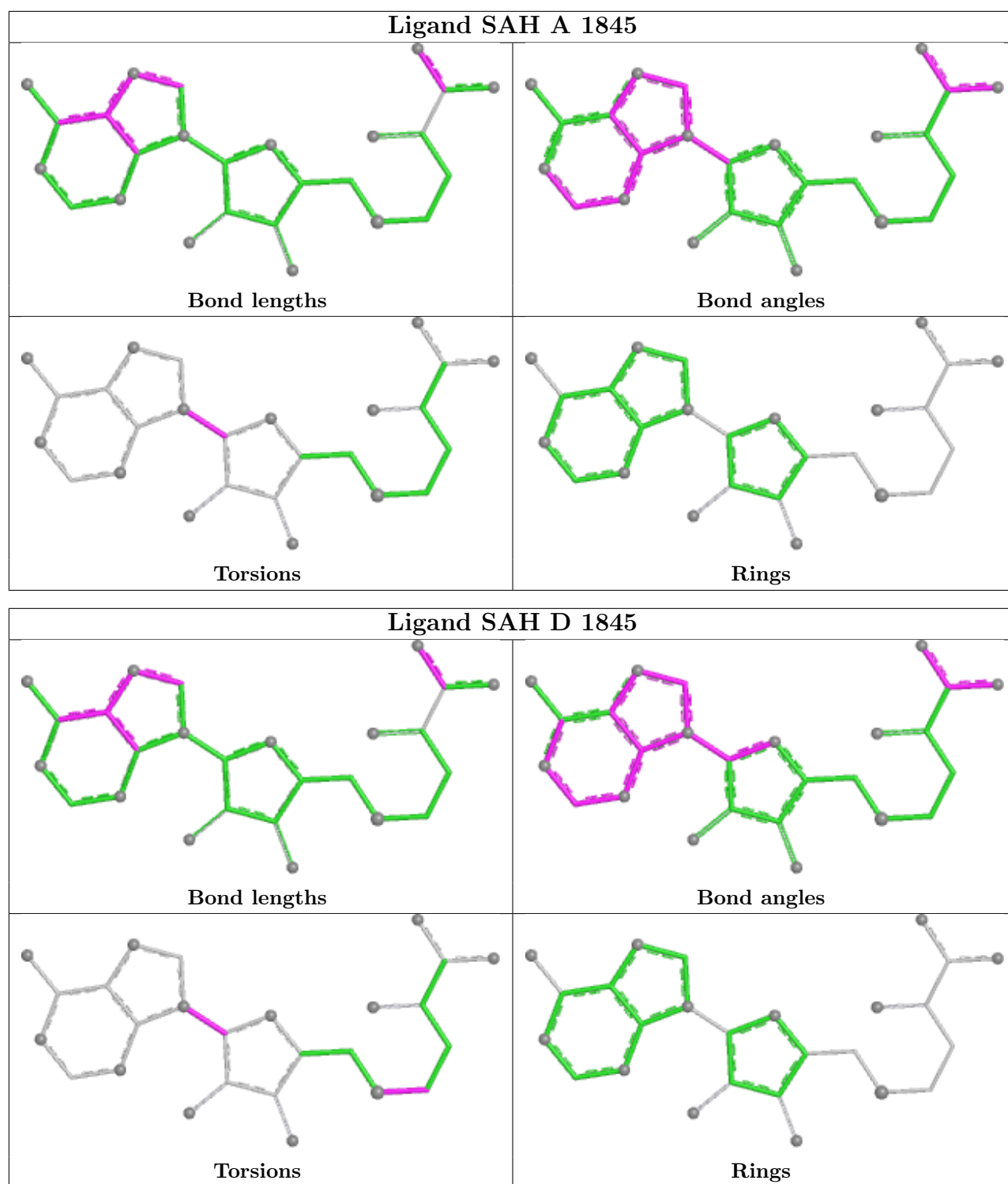
All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	D	1845	SAH	CB-CG-SD-C5'
3	D	1845	SAH	C2'-C1'-N9-C4
3	A	1845	SAH	C2'-C1'-N9-C4
3	A	1845	SAH	C2'-C1'-N9-C8
3	D	1845	SAH	C2'-C1'-N9-C8
3	A	1845	SAH	O4'-C1'-N9-C8
3	D	1845	SAH	O4'-C1'-N9-C8

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 ⓘ

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2		OWAB(Å ²)	Q<0.9
1	A	819/844 (97%)	0.58	53 (6%)	25 20	18, 46, 98, 115	1 (0%)
1	D	821/844 (97%)	0.32	24 (2%)	53 45	12, 36, 77, 111	1 (0%)
2	B	284/287 (98%)	0.29	8 (2%)	55 46	19, 39, 74, 94	0
2	E	283/287 (98%)	0.07	3 (1%)	78 71	16, 24, 57, 82	0
All	All	2207/2262 (97%)	0.38	88 (3%)	42 34	12, 38, 84, 115	2 (0%)

All (88) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	640	ILE	4.6
1	D	409	TYR	4.4
1	A	526	GLY	4.2
1	A	354	GLY	3.7
1	A	272	SER	3.5
2	B	90	TYR	3.5
1	A	234	ASN	3.5
1	A	669[A]	PHE	3.4
1	A	34	ILE	3.4
1	A	344	ARG	3.3
1	D	27	GLN	3.3
1	D	669[A]	PHE	3.3
1	A	154	LYS	3.3
1	D	25	LEU	3.2
1	A	640	ILE	3.2
1	A	391	PHE	3.1
1	A	324	ILE	3.1
1	A	295	ASP	3.0
1	A	293	ILE	3.0
1	D	354	GLY	2.9
1	D	435	PHE	2.9

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Mol	Chain	Res	Type	RSRZ
1	D	529	PHE	2.9
1	A	204	PRO	2.9
1	D	34	ILE	2.9
2	B	118	LYS	2.9
1	A	130	ILE	2.9
2	B	119	SER	2.8
2	B	1	MET	2.7
1	D	262	ASP	2.7
1	D	555	TYR	2.6
1	A	345	ILE	2.6
1	A	226	ASN	2.6
1	A	319	GLU	2.6
1	A	205	TYR	2.6
2	E	124	THR	2.6
1	A	731	ASN	2.5
2	E	90	TYR	2.5
1	A	357	THR	2.5
1	A	294	ILE	2.5
1	A	206	ASP	2.4
1	A	248	VAL	2.4
1	D	26	GLU	2.4
1	D	639	GLY	2.4
1	A	641	LYS	2.4
1	D	410	MET	2.4
2	B	87	VAL	2.4
1	D	205	TYR	2.4
1	A	25	LEU	2.4
1	A	553	VAL	2.4
1	A	555	TYR	2.4
1	A	298	VAL	2.3
1	A	2	ASP	2.3
1	A	362	VAL	2.3
1	A	27	GLN	2.3
1	A	275	LEU	2.3
1	A	8	SER	2.3
1	D	133	ASP	2.2
1	D	642	THR	2.2
2	B	124	THR	2.2
1	D	420	SER	2.2
1	A	209	ILE	2.2
2	B	271	SER	2.2
1	A	215	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
1	D	325	ASN	2.2
1	D	100	ASP	2.2
1	A	499	LYS	2.2
1	A	356	PHE	2.2
1	A	273	ASN	2.1
1	D	559	LYS	2.1
1	A	528	ILE	2.1
1	A	448	ASN	2.1
1	D	71	LYS	2.1
1	A	17	ALA	2.1
1	A	59	ILE	2.1
1	A	318	ILE	2.1
1	A	271	THR	2.1
1	D	196	THR	2.1
1	D	421	SER	2.1
1	A	1	MET	2.1
1	A	5	VAL	2.1
1	A	24	GLU	2.1
1	A	361	GLU	2.1
2	E	118	LYS	2.0
2	B	120	ASN	2.0
1	A	347	PHE	2.0
1	A	296	SER	2.0
1	A	629	GLY	2.0
1	A	639	GLY	2.0

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

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

6.3 Carbohydrates ⓘ

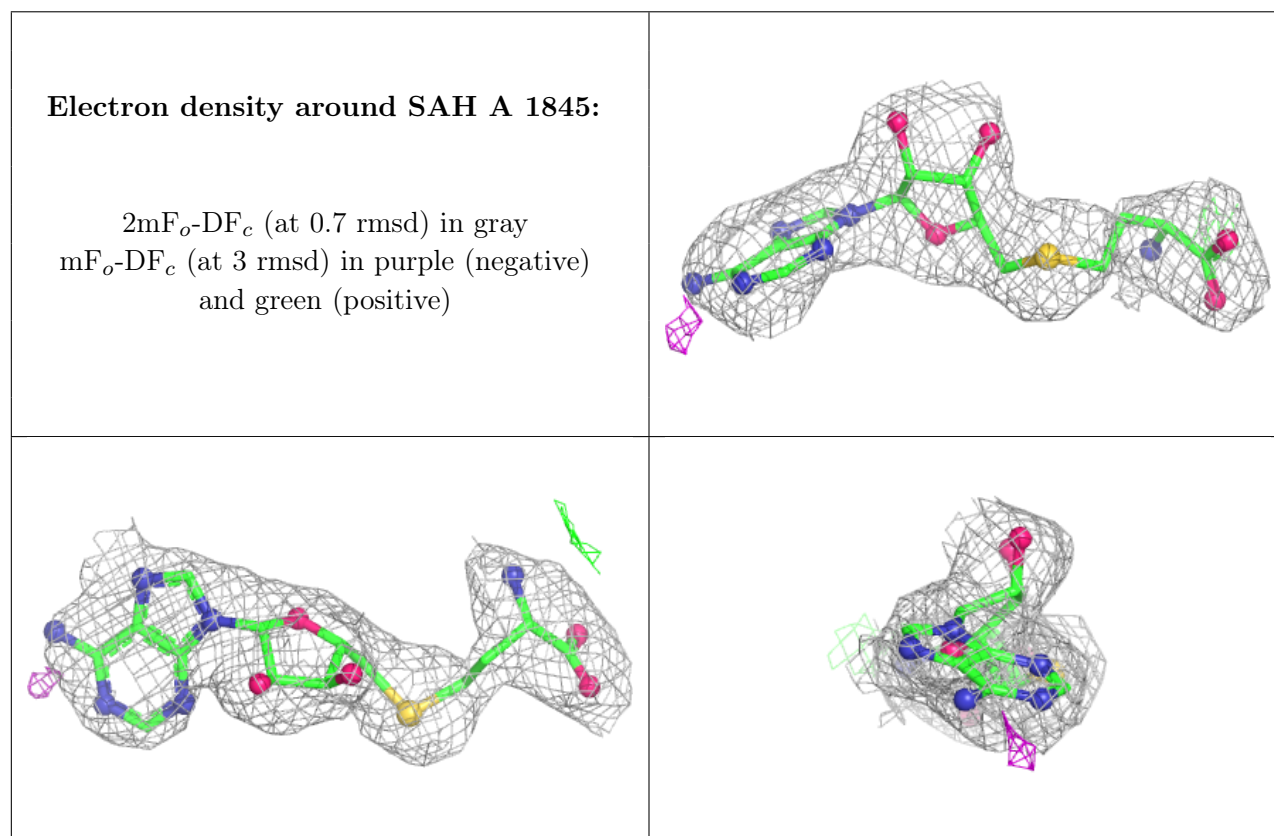
There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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

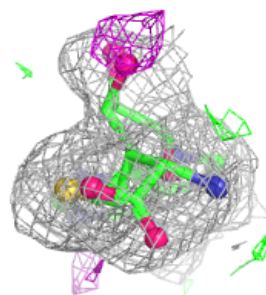
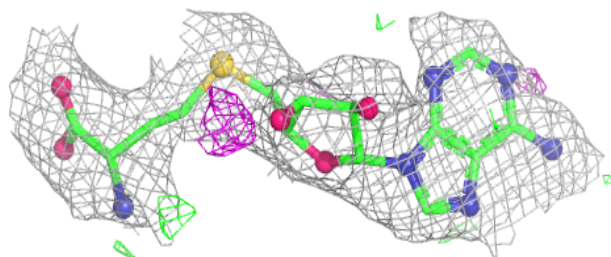
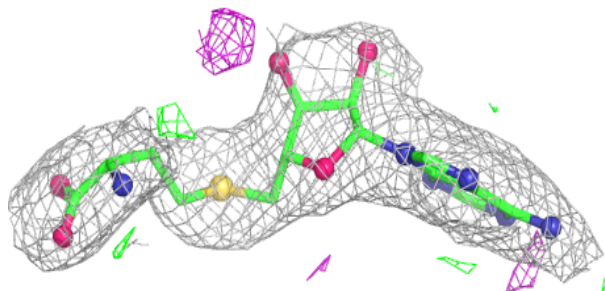
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	SAH	A	1845	26/26	0.94	0.09	22,24,25,25	0
3	SAH	D	1845	26/26	0.94	0.09	18,19,20,21	0

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 SAH D 1845:

$2mF_o-DF_c$ (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.