



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 01:40 AM GMT

PDB ID : 2DUB
Title : ENOYL-COA HYDRATASE COMPLEXED WITH OCTANOYL-COA
Authors : Engel, C.K.; Wierenga, R.K.
Deposited on : 1997-04-28
Resolution : 2.40 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
<http://wwpdb.org/validation/2016/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 : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

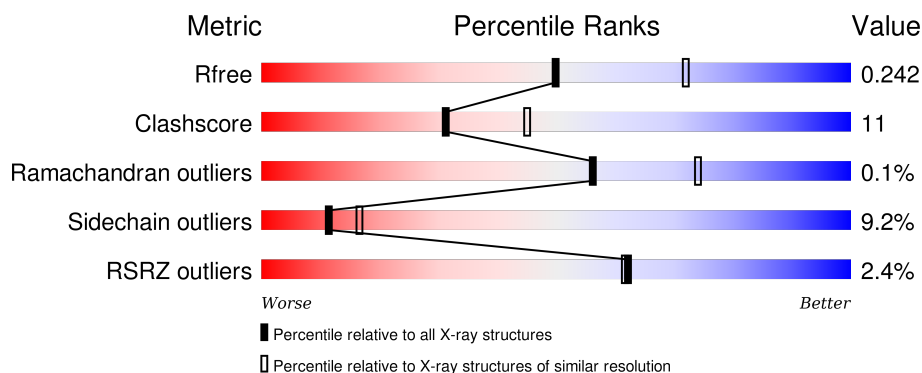
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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}	91344	2919 (2.40-2.40)
Clashscore	102246	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)
RSRZ outliers	91569	2928 (2.40-2.40)

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. 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	261	<div> <div>3%</div> <div> <div></div> <div>74%</div> <div>21%</div> <div>• •</div> </div> </div>
1	B	261	<div> <div>2%</div> <div> <div></div> <div>74%</div> <div>20%</div> <div>• •</div> </div> </div>
1	C	261	<div> <div>5%</div> <div> <div></div> <div>72%</div> <div>23%</div> <div>• •</div> </div> </div>
1	D	261	<div> <div>2%</div> <div> <div></div> <div>76%</div> <div>19%</div> <div>• •</div> </div> </div>
1	E	261	<div> <div>2%</div> <div> <div></div> <div>72%</div> <div>23%</div> <div>• •</div> </div> </div>

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Mol	Chain	Length	Quality of chain
1	F	261	<div> <div></div> <div>%</div> <div>76%</div> <div>20%</div> <div>• •</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	CO8	B	300	-	-	-	X
2	CO8	F	300	-	-	-	X

2 Entry composition [i](#)

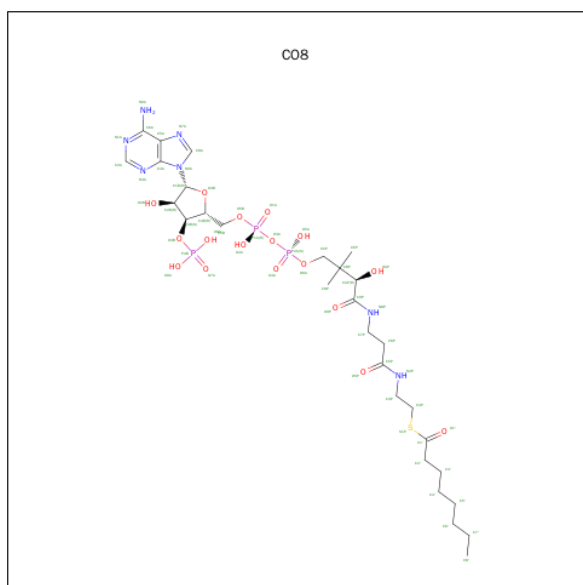
There are 3 unique types of molecules in this entry. The entry contains 12205 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 2-ENOYL-COA HYDRATASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	254	Total	C	N	O	S	0	0	0
			1921	1206	330	372	13			
1	B	254	Total	C	N	O	S	0	0	0
			1918	1205	329	371	13			
1	C	254	Total	C	N	O	S	0	0	0
			1918	1205	329	371	13			
1	D	257	Total	C	N	O	S	0	0	0
			1935	1214	333	375	13			
1	E	255	Total	C	N	O	S	0	0	0
			1923	1208	330	372	13			
1	F	255	Total	C	N	O	S	0	0	0
			1926	1209	331	373	13			

- Molecule 2 is OCTANOYL-COENZYME A (three-letter code: CO8) (formula: $C_{29}H_{50}N_7O_{17}P_3S$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	S	0	0
			57	29	7	17	3	1		
2	B	1	Total	C	N	O	P	S	0	0
			57	29	7	17	3	1		
2	E	1	Total	C	N	O	P	S	0	0
			57	29	7	17	3	1		
2	F	1	Total	C	N	O	P	S	0	0
			57	29	7	17	3	1		

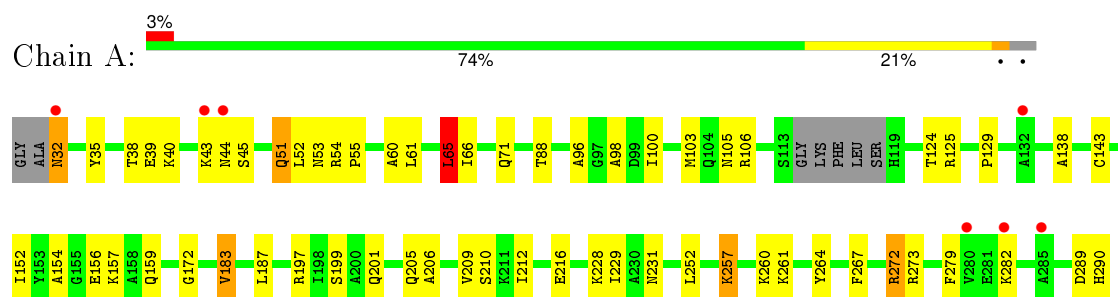
- Molecule 3 is water.

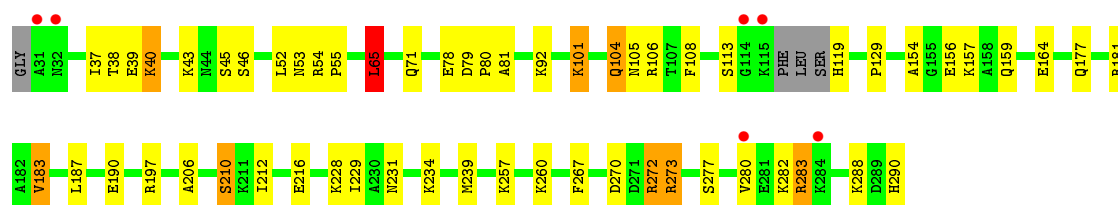
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	76	Total	O	0	0
			76	76		
3	B	73	Total	O	0	0
			73	73		
3	C	59	Total	O	0	0
			59	59		
3	D	73	Total	O	0	0
			73	73		
3	E	82	Total	O	0	0
			82	82		
3	F	73	Total	O	0	0
			73	73		

3 Residue-property plots [i](#)

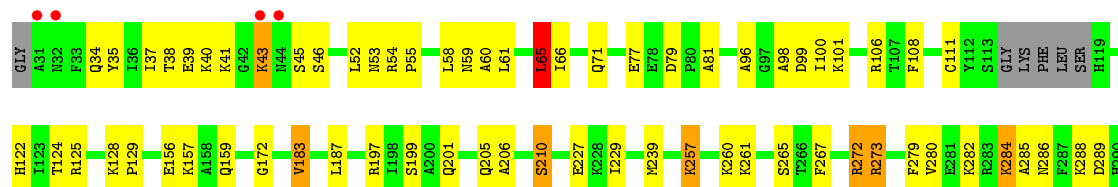
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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: 2-ENOYL-COA HYDRATASE

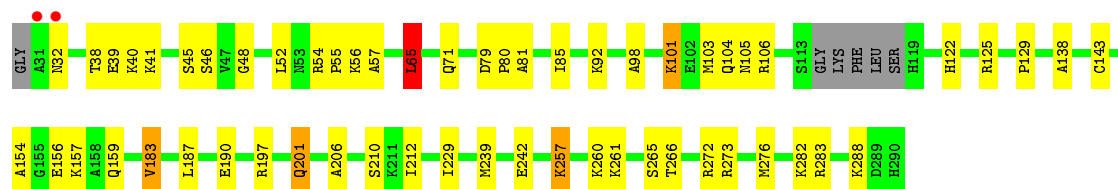
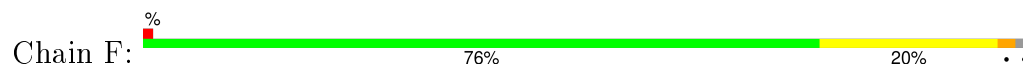




• Molecule 1: 2-ENOYL-COA HYDRATASE



• Molecule 1: 2-ENOYL-COA HYDRATASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	76.24Å 93.93Å 246.98Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.40 19.99 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.6 (20.00-2.40) 99.6 (19.99-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.37 (at 2.41Å)	Xtriage
Refinement program	REFMAC, X-PLOR	Depositor
R, R_{free}	0.204 , 0.260 0.199 , 0.242	Depositor DCC
R_{free} test set	3525 reflections (5.31%)	DCC
Wilson B-factor (Å ²)	28.1	Xtriage
Anisotropy	0.501	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 66.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	1 of 69852 reflections (0.001%)	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	12205	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.68% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: CO8

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	0.37	0/1943	0.58	1/2608 (0.0%)
1	B	0.37	0/1940	0.58	1/2604 (0.0%)
1	C	0.38	0/1940	0.58	1/2604 (0.0%)
1	D	0.37	0/1957	0.59	1/2627 (0.0%)
1	E	0.38	0/1945	0.58	1/2611 (0.0%)
1	F	0.36	0/1948	0.59	1/2615 (0.0%)
All	All	0.37	0/11673	0.58	6/15669 (0.0%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	65	LEU	CA-CB-CG	5.60	128.18	115.30
1	A	65	LEU	CA-CB-CG	5.31	127.51	115.30
1	E	65	LEU	CA-CB-CG	5.25	127.39	115.30
1	D	65	LEU	CA-CB-CG	5.20	127.27	115.30
1	B	65	LEU	CA-CB-CG	5.10	127.03	115.30
1	C	65	LEU	CA-CB-CG	5.01	126.82	115.30

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	1921	0	1942	46	0
1	B	1918	0	1938	49	0
1	C	1918	0	1938	44	0
1	D	1935	0	1952	42	0
1	E	1923	0	1943	54	0
1	F	1926	0	1947	43	0
2	A	57	0	46	6	0
2	B	57	0	46	7	0
2	E	57	0	46	6	0
2	F	57	0	46	1	0
3	A	76	0	0	2	0
3	B	73	0	0	3	0
3	C	59	0	0	3	0
3	D	73	0	0	2	0
3	E	82	0	0	3	0
3	F	73	0	0	2	0
All	All	12205	0	11844	253	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 11.

All (253) 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:104:GLN:HG3	1:C:280:VAL:HG21	1.42	1.01
1:F:201:GLN:H	1:F:201:GLN:NE2	1.66	0.94
1:A:267:PHE:HA	1:A:272:ARG:HD3	1.51	0.94
1:B:172:GLY:H	2:B:300:CO8:H3'1	1.34	0.90
1:D:101:LYS:H	1:D:101:LYS:HD2	1.41	0.86
1:C:108:PHE:HZ	1:D:239:MET:HE1	1.39	0.85
1:A:172:GLY:H	2:A:300:CO8:H3'1	1.41	0.84
1:E:172:GLY:H	2:E:300:CO8:H3'1	1.46	0.81
1:E:183:VAL:HG22	1:E:206:ALA:HB1	1.64	0.80
1:B:183:VAL:HG22	1:B:206:ALA:HB1	1.65	0.78
1:C:183:VAL:HG22	1:C:206:ALA:HB1	1.68	0.76
1:D:183:VAL:HG22	1:D:206:ALA:HB1	1.69	0.74
1:A:201:GLN:H	1:A:201:GLN:NE2	1.86	0.73
1:F:183:VAL:HG22	1:F:206:ALA:HB1	1.71	0.73
1:F:52:LEU:HD22	1:F:65:LEU:HD11	1.71	0.72
1:B:100:ILE:HG23	1:C:276:MET:HG2	1.70	0.71
1:C:52:LEU:HD22	1:C:65:LEU:HD11	1.71	0.71
1:A:183:VAL:HG22	1:A:206:ALA:HB1	1.72	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:273:ARG:HH22	1:F:105:ASN:ND2	1.92	0.68
1:E:100:ILE:HG23	1:F:276:MET:HG2	1.76	0.67
1:B:201:GLN:NE2	1:B:201:GLN:H	1.93	0.67
1:D:280:VAL:HG21	1:F:104:GLN:NE2	2.09	0.66
1:B:183:VAL:HG13	1:B:187:LEU:HB3	1.78	0.66
1:B:239:MET:HE1	1:B:265:SER:HB3	1.77	0.66
3:C:303:HOH:O	1:E:257:LYS:HB3	1.95	0.66
1:B:129:PRO:HG2	1:B:229:ILE:HG21	1.77	0.65
1:D:78:GLU:O	1:D:80:PRO:HD3	1.95	0.65
1:C:267:PHE:HA	1:C:272:ARG:HD3	1.79	0.65
1:A:52:LEU:HD22	1:A:65:LEU:HD11	1.78	0.65
1:E:267:PHE:HA	1:E:272:ARG:HD2	1.78	0.64
1:D:183:VAL:CG1	1:D:187:LEU:HB3	2.27	0.64
1:F:183:VAL:HG13	1:F:187:LEU:HB3	1.78	0.64
1:E:183:VAL:HG13	1:E:187:LEU:HB3	1.80	0.64
1:C:201:GLN:H	1:C:201:GLN:NE2	1.96	0.63
1:A:183:VAL:HG13	1:A:187:LEU:HB3	1.80	0.63
1:D:183:VAL:HG13	1:D:187:LEU:HB3	1.79	0.63
1:E:43:LYS:HB2	1:E:227:GLU:OE2	1.98	0.63
1:C:183:VAL:HG13	1:C:187:LEU:HB3	1.82	0.62
1:A:199:SER:HB2	1:A:201:GLN:NE2	2.14	0.62
1:E:201:GLN:NE2	1:E:201:GLN:H	1.98	0.62
1:A:124:THR:HG22	3:A:360:HOH:O	1.99	0.62
1:C:78:GLU:O	1:C:80:PRO:HD3	2.00	0.62
1:E:183:VAL:CG1	1:E:187:LEU:HB3	2.30	0.61
1:F:239:MET:HE1	1:F:265:SER:HB3	1.82	0.61
1:D:283:ARG:HH11	1:D:283:ARG:HG3	1.65	0.61
1:A:228:LYS:O	1:A:231:ASN:HB2	2.01	0.61
1:A:183:VAL:CG1	1:A:187:LEU:HB3	2.30	0.61
1:E:272:ARG:HD3	3:E:361:HOH:O	2.02	0.60
1:F:129:PRO:HG2	1:F:229:ILE:HG21	1.83	0.60
1:A:129:PRO:HG2	1:A:229:ILE:HG21	1.83	0.60
1:C:183:VAL:CG1	1:C:187:LEU:HB3	2.32	0.60
1:E:122:HIS:HA	1:E:125:ARG:HD3	1.84	0.60
1:A:289:ASP:O	1:A:290:HIS:HB3	2.02	0.59
1:C:65:LEU:HD12	1:C:65:LEU:C	2.23	0.59
1:B:183:VAL:CG1	1:B:187:LEU:HB3	2.33	0.58
1:C:65:LEU:HD12	1:C:65:LEU:O	2.03	0.58
1:D:129:PRO:HG2	1:D:229:ILE:HG21	1.85	0.58
1:C:216:GLU:CD	1:C:216:GLU:H	2.06	0.58
1:B:183:VAL:HG22	1:B:206:ALA:CB	2.33	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:65:LEU:C	1:D:65:LEU:HD12	2.25	0.57
1:B:99:ASP:HA	2:B:300:CO8:N1A	2.20	0.57
1:E:183:VAL:HG22	1:E:206:ALA:CB	2.34	0.57
1:D:52:LEU:HD22	1:D:65:LEU:HD11	1.86	0.57
1:B:65:LEU:C	1:B:65:LEU:HD12	2.24	0.57
1:A:65:LEU:HD12	1:A:65:LEU:C	2.25	0.57
1:F:183:VAL:CG1	1:F:187:LEU:HB3	2.35	0.56
1:B:54:ARG:HH12	1:B:62:CYS:HB3	1.69	0.56
1:C:183:VAL:HG22	1:C:206:ALA:CB	2.34	0.56
1:D:183:VAL:HG22	1:D:206:ALA:CB	2.36	0.56
1:E:284:LYS:HE2	1:E:285:ALA:O	2.04	0.56
1:E:65:LEU:C	1:E:65:LEU:HD12	2.26	0.56
1:E:54:ARG:N	1:E:55:PRO:CD	2.68	0.56
1:B:283:ARG:HD2	3:B:305:HOH:O	2.04	0.56
1:D:105:ASN:HA	1:E:273:ARG:HH12	1.71	0.55
1:E:108:PHE:O	1:E:111:CYS:HB2	2.06	0.55
1:E:46:SER:HA	1:E:81:ALA:O	2.06	0.55
1:F:183:VAL:HG22	1:F:206:ALA:CB	2.36	0.55
1:A:201:GLN:H	1:A:201:GLN:CD	2.09	0.55
1:F:122:HIS:HA	1:F:125:ARG:NH1	2.20	0.55
1:A:61:LEU:HB3	1:A:66:ILE:HD11	1.88	0.55
1:F:65:LEU:C	1:F:65:LEU:HD12	2.28	0.54
1:D:210:SER:HB2	1:F:190:GLU:OE1	2.08	0.54
1:B:190:GLU:OE1	1:C:210:SER:HB2	2.07	0.54
1:E:100:ILE:CG2	1:F:276:MET:HG2	2.37	0.54
1:A:43:LYS:HG3	1:A:44:ASN:ND2	2.23	0.54
1:C:129:PRO:HG2	1:C:229:ILE:HG21	1.90	0.53
1:A:40:LYS:HE2	1:A:45:SER:HB3	1.91	0.53
1:A:199:SER:HB2	1:A:201:GLN:HE22	1.74	0.53
1:D:280:VAL:HG11	1:F:104:GLN:HE22	1.74	0.53
1:B:52:LEU:HD22	1:B:65:LEU:HD11	1.90	0.53
1:D:119:HIS:N	3:D:338:HOH:O	2.41	0.52
1:C:108:PHE:HZ	1:D:239:MET:CE	2.16	0.52
1:C:49:LEU:HD12	1:C:50:ILE:N	2.25	0.52
1:A:183:VAL:HG22	1:A:206:ALA:CB	2.39	0.52
1:E:52:LEU:HD22	1:E:65:LEU:HD11	1.92	0.52
1:B:96:ALA:O	2:B:300:CO8:H62	2.10	0.52
1:C:35:TYR:O	1:C:53:ASN:N	2.40	0.52
1:D:283:ARG:HH11	1:D:283:ARG:CG	2.23	0.51
1:B:54:ARG:NH1	1:B:62:CYS:HB3	2.25	0.51
1:E:54:ARG:HH12	1:E:60:ALA:HB3	1.75	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:75:THR:HB	3:C:326:HOH:O	2.10	0.51
1:E:61:LEU:HB3	1:E:66:ILE:HD11	1.90	0.51
1:D:65:LEU:HD12	1:D:65:LEU:O	2.10	0.51
1:A:65:LEU:HD12	1:A:65:LEU:O	2.10	0.51
1:F:40:LYS:HE2	1:F:79:ASP:OD2	2.09	0.51
1:A:172:GLY:N	2:A:300:CO8:H3'1	2.20	0.51
1:F:201:GLN:HE21	1:F:201:GLN:H	1.54	0.51
1:C:225:CYS:HA	1:C:228:LYS:HE3	1.93	0.50
1:F:159:GLN:NE2	1:F:197:ARG:HD2	2.25	0.50
1:B:164:GLU:HG2	2:B:300:CO8:H2'2	1.92	0.50
1:A:279:PHE:O	1:A:282:LYS:HE2	2.12	0.50
1:B:40:LYS:HE2	1:B:79:ASP:OD2	2.12	0.50
1:A:159:GLN:NE2	1:A:197:ARG:HD2	2.27	0.50
1:B:172:GLY:N	2:B:300:CO8:H3'1	2.15	0.49
1:D:280:VAL:HG21	1:F:104:GLN:HE21	1.77	0.49
1:E:267:PHE:HA	1:E:272:ARG:CD	2.42	0.49
1:C:54:ARG:NH1	1:C:60:ALA:HB3	2.27	0.49
1:B:104:GLN:CG	1:C:280:VAL:HG21	2.29	0.49
1:E:239:MET:HE1	1:E:265:SER:HB3	1.94	0.49
1:A:51:GLN:HG3	1:A:88:THR:HG23	1.95	0.49
1:E:129:PRO:HG2	1:E:229:ILE:HG21	1.95	0.49
1:C:108:PHE:CZ	1:D:239:MET:HE1	2.31	0.48
1:D:104:GLN:O	1:E:272:ARG:NH2	2.41	0.48
1:B:54:ARG:N	1:B:55:PRO:CD	2.76	0.48
1:B:201:GLN:CD	1:B:201:GLN:H	2.16	0.48
1:E:65:LEU:O	1:E:65:LEU:HD12	2.12	0.48
1:E:54:ARG:NH1	1:E:60:ALA:HB3	2.28	0.48
1:B:46:SER:HA	1:B:81:ALA:O	2.13	0.48
1:D:177:GLN:O	1:D:181:ARG:HG3	2.13	0.48
1:C:234:LYS:NZ	3:C:343:HOH:O	2.30	0.48
1:E:159:GLN:NE2	1:E:197:ARG:HD2	2.29	0.48
1:B:65:LEU:O	1:B:65:LEU:HD12	2.13	0.48
1:D:267:PHE:CD2	1:D:272:ARG:HD2	2.49	0.48
1:C:33:PHE:HE1	1:C:71:GLN:OE1	1.95	0.48
1:B:104:GLN:HE21	1:B:104:GLN:C	2.17	0.48
1:E:199:SER:HB2	1:E:201:GLN:NE2	2.29	0.48
1:D:40:LYS:CE	1:D:79:ASP:OD2	2.62	0.47
1:A:54:ARG:NH1	1:A:60:ALA:HB3	2.29	0.47
1:C:239:MET:HE1	1:D:108:PHE:HZ	1.79	0.47
1:F:201:GLN:NE2	1:F:201:GLN:N	2.49	0.47
1:E:40:LYS:NZ	1:E:79:ASP:OD2	2.37	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:99:ASP:OD2	1:B:102:GLU:HG3	2.14	0.47
1:D:280:VAL:HG11	1:F:104:GLN:NE2	2.30	0.47
1:D:272:ARG:HD3	3:D:291:HOH:O	2.15	0.47
1:A:100:ILE:HD11	2:A:300:CO8:H22	1.97	0.47
1:D:46:SER:HA	1:D:81:ALA:O	2.14	0.47
1:C:289:ASP:O	1:C:290:HIS:HB3	2.15	0.47
1:F:38:THR:O	1:F:39:GLU:HB3	2.15	0.47
1:F:98:ALA:O	2:F:300:CO8:N6A	2.48	0.47
1:F:283:ARG:HD2	3:F:361:HOH:O	2.14	0.47
1:B:60:ALA:HA	1:B:97:GLY:O	2.15	0.46
1:C:46:SER:HA	1:C:81:ALA:O	2.15	0.46
1:E:96:ALA:O	2:E:300:CO8:H62	2.15	0.46
1:F:101:LYS:HG2	1:F:101:LYS:H	1.46	0.46
1:B:159:GLN:NE2	1:B:197:ARG:HD2	2.30	0.46
1:E:100:ILE:HG12	2:E:300:CO8:N1A	2.31	0.46
1:E:289:ASP:OD1	3:E:307:HOH:O	2.21	0.46
1:F:46:SER:HA	1:F:81:ALA:O	2.15	0.46
1:E:257:LYS:O	1:E:261:LYS:HG2	2.15	0.45
1:B:284:LYS:HD2	3:B:351:HOH:O	2.15	0.45
1:F:65:LEU:O	1:F:65:LEU:HD12	2.16	0.45
1:D:190:GLU:OE1	1:E:210:SER:HB2	2.16	0.45
1:F:40:LYS:CE	1:F:79:ASP:OD2	2.64	0.45
1:C:54:ARG:N	1:C:55:PRO:CD	2.80	0.45
1:C:284:LYS:HG3	1:C:285:ALA:N	2.31	0.45
1:E:98:ALA:O	2:E:300:CO8:N6A	2.49	0.45
1:F:54:ARG:N	1:F:55:PRO:CD	2.78	0.45
1:A:35:TYR:O	1:A:53:ASN:N	2.44	0.45
1:A:105:ASN:HB2	3:A:330:HOH:O	2.17	0.45
1:D:228:LYS:O	1:D:231:ASN:HB2	2.16	0.45
1:F:138:ALA:O	1:F:143:CYS:HB2	2.17	0.45
1:A:138:ALA:O	1:A:143:CYS:HB2	2.17	0.44
1:F:266:THR:O	1:F:272:ARG:HB2	2.18	0.44
1:F:48:GLY:O	1:F:85:ILE:HG23	2.17	0.44
1:A:106:ARG:NE	1:A:106:ARG:HA	2.32	0.44
1:D:283:ARG:NH1	1:D:283:ARG:CG	2.81	0.44
1:B:257:LYS:O	1:B:261:LYS:HG2	2.18	0.44
1:A:96:ALA:O	2:A:300:CO8:H62	2.18	0.44
1:B:106:ARG:HD3	1:B:106:ARG:HA	1.74	0.44
1:A:257:LYS:O	1:A:261:LYS:HG2	2.17	0.44
1:F:122:HIS:HA	1:F:125:ARG:HH11	1.83	0.43
1:E:124:THR:HG22	3:E:351:HOH:O	2.17	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:267:PHE:HA	1:B:272:ARG:HD2	1.99	0.43
1:E:41:LYS:O	1:E:45:SER:HA	2.18	0.43
1:B:54:ARG:NH1	1:B:62:CYS:SG	2.92	0.43
1:C:61:LEU:HB3	1:C:66:ILE:HD11	2.00	0.43
1:B:48:GLY:O	1:B:85:ILE:HG23	2.18	0.43
1:B:154:ALA:O	1:B:212:ILE:HA	2.18	0.43
1:D:104:GLN:OE1	1:E:280:VAL:HG21	2.18	0.43
1:F:79:ASP:HA	1:F:80:PRO:HD2	1.83	0.43
1:F:54:ARG:HG2	1:F:57:ALA:HB3	2.00	0.43
1:D:54:ARG:N	1:D:55:PRO:CD	2.81	0.43
1:B:216:GLU:CD	1:B:216:GLU:H	2.22	0.43
1:B:205:GLN:HE21	1:B:205:GLN:HA	1.83	0.43
1:C:159:GLN:NE2	1:C:197:ARG:HD2	2.34	0.43
1:E:37:ILE:HD11	1:E:53:ASN:ND2	2.34	0.42
1:E:201:GLN:CD	1:E:201:GLN:H	2.21	0.42
1:C:38:THR:O	1:C:39:GLU:HB3	2.19	0.42
1:F:257:LYS:O	1:F:261:LYS:HG2	2.19	0.42
1:B:57:ALA:HA	2:B:300:CO8:O4B	2.18	0.42
1:E:43:LYS:N	1:E:227:GLU:OE2	2.31	0.42
1:C:132:ALA:HB2	1:C:149:CYS:SG	2.59	0.42
1:A:38:THR:O	1:A:39:GLU:HB3	2.18	0.42
1:B:172:GLY:H	2:B:300:CO8:C3'	2.19	0.42
1:A:54:ARG:HH12	1:A:60:ALA:HB3	1.85	0.42
1:D:270:ASP:OD2	1:D:290:HIS:ND1	2.50	0.42
1:D:37:ILE:HD11	1:D:53:ASN:ND2	2.34	0.42
1:B:100:ILE:HD12	1:B:167:LEU:HD13	2.02	0.42
1:B:65:LEU:CD1	1:B:65:LEU:C	2.86	0.42
3:B:329:HOH:O	1:F:257:LYS:HB3	2.20	0.42
1:C:257:LYS:O	1:C:261:LYS:HG2	2.20	0.42
1:D:159:GLN:NE2	1:D:197:ARG:HD2	2.34	0.42
1:E:286:ASN:ND2	1:E:288:LYS:HE3	2.34	0.42
1:A:54:ARG:N	1:A:55:PRO:CD	2.82	0.42
1:C:154:ALA:O	1:C:212:ILE:HA	2.20	0.42
1:C:65:LEU:C	1:C:65:LEU:CD1	2.89	0.42
1:C:54:ARG:H	1:C:59:ASN:ND2	2.17	0.42
1:A:264:TYR:CD1	1:C:111:CYS:HB3	2.55	0.42
1:B:99:ASP:HB3	1:B:102:GLU:HG3	2.02	0.41
1:E:59:ASN:O	1:E:60:ALA:C	2.58	0.41
1:D:38:THR:O	1:D:39:GLU:HB3	2.19	0.41
1:C:49:LEU:HD22	1:C:223:ILE:HD11	2.01	0.41
1:E:239:MET:CE	1:E:265:SER:HB3	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:154:ALA:O	1:D:212:ILE:HA	2.21	0.41
1:A:32:ASN:HA	1:A:32:ASN:HD22	1.51	0.41
1:A:154:ALA:O	1:A:212:ILE:HA	2.21	0.41
1:F:154:ALA:O	1:F:212:ILE:HA	2.21	0.41
1:A:205:GLN:HE21	1:B:205:GLN:NE2	2.18	0.41
1:A:152:ILE:HG21	1:A:209:VAL:HG12	2.03	0.41
1:F:40:LYS:HB3	1:F:45:SER:HB3	2.01	0.41
1:A:54:ARG:NH1	1:A:60:ALA:O	2.53	0.41
1:B:78:GLU:O	1:B:80:PRO:HD3	2.21	0.41
1:F:201:GLN:H	1:F:201:GLN:CD	2.19	0.41
1:D:65:LEU:C	1:D:65:LEU:CD1	2.88	0.41
1:A:264:TYR:CG	1:C:111:CYS:HB3	2.56	0.41
1:F:56:LYS:HB2	3:F:344:HOH:O	2.20	0.41
1:B:261:LYS:NZ	1:F:242:GLU:OE2	2.50	0.41
1:A:98:ALA:O	2:A:300:CO8:N6A	2.53	0.41
1:E:77:GLU:OE1	1:E:128:LYS:HG3	2.20	0.41
1:B:38:THR:O	1:B:39:GLU:HB3	2.21	0.41
1:B:82:VAL:O	1:B:234:LYS:HE2	2.21	0.41
1:C:122:HIS:HA	1:C:125:ARG:HD3	2.02	0.41
1:E:99:ASP:HA	2:E:300:CO8:N1A	2.36	0.41
1:A:252:LEU:HB3	1:C:249:GLU:OE2	2.21	0.41
1:E:172:GLY:N	2:E:300:CO8:H3'1	2.25	0.40
1:E:35:TYR:CD1	1:E:54:ARG:HD2	2.55	0.40
1:E:38:THR:O	1:E:39:GLU:HB3	2.21	0.40
1:E:58:LEU:HD23	1:E:58:LEU:HA	1.90	0.40
1:A:267:PHE:HA	1:A:272:ARG:CD	2.37	0.40
1:D:164:GLU:OE1	1:D:164:GLU:N	2.43	0.40
1:E:279:PHE:O	1:E:282:LYS:HD2	2.21	0.40
1:A:100:ILE:HG12	2:A:300:CO8:N1A	2.37	0.40
1:E:65:LEU:C	1:E:65:LEU:CD1	2.90	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	250/261 (96%)	240 (96%)	10 (4%)	0	100	100
1	B	250/261 (96%)	241 (96%)	9 (4%)	0	100	100
1	C	250/261 (96%)	239 (96%)	11 (4%)	0	100	100
1	D	253/261 (97%)	245 (97%)	7 (3%)	1 (0%)	39	56
1	E	251/261 (96%)	246 (98%)	5 (2%)	0	100	100
1	F	251/261 (96%)	246 (98%)	5 (2%)	0	100	100
All	All	1505/1566 (96%)	1457 (97%)	47 (3%)	1 (0%)	56	74

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	113	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	199/205 (97%)	184 (92%)	15 (8%)	17	26
1	B	198/205 (97%)	177 (89%)	21 (11%)	8	12
1	C	198/205 (97%)	182 (92%)	16 (8%)	15	22
1	D	199/205 (97%)	176 (88%)	23 (12%)	7	9
1	E	198/205 (97%)	182 (92%)	16 (8%)	15	22
1	F	199/205 (97%)	181 (91%)	18 (9%)	12	17
All	All	1191/1230 (97%)	1082 (91%)	109 (9%)	11	16

All (109) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	32	ASN

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Mol	Chain	Res	Type
1	A	51	GLN
1	A	65	LEU
1	A	71	GLN
1	A	103	MET
1	A	125	ARG
1	A	156	GLU
1	A	157	LYS
1	A	183	VAL
1	A	210	SER
1	A	216	GLU
1	A	257	LYS
1	A	260	LYS
1	A	272	ARG
1	A	273	ARG
1	B	40	LYS
1	B	41	LYS
1	B	65	LEU
1	B	71	GLN
1	B	101	LYS
1	B	103	MET
1	B	104	GLN
1	B	106	ARG
1	B	125	ARG
1	B	127	LYS
1	B	156	GLU
1	B	157	LYS
1	B	183	VAL
1	B	205	GLN
1	B	210	SER
1	B	216	GLU
1	B	257	LYS
1	B	260	LYS
1	B	272	ARG
1	B	282	LYS
1	B	288	LYS
1	C	56	LYS
1	C	65	LEU
1	C	71	GLN
1	C	101	LYS
1	C	103	MET
1	C	156	GLU
1	C	157	LYS

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Mol	Chain	Res	Type
1	C	183	VAL
1	C	205	GLN
1	C	210	SER
1	C	227	GLU
1	C	257	LYS
1	C	260	LYS
1	C	272	ARG
1	C	273	ARG
1	C	288	LYS
1	D	40	LYS
1	D	43	LYS
1	D	45	SER
1	D	65	LEU
1	D	71	GLN
1	D	92	LYS
1	D	101	LYS
1	D	104	GLN
1	D	106	ARG
1	D	156	GLU
1	D	157	LYS
1	D	183	VAL
1	D	210	SER
1	D	216	GLU
1	D	234	LYS
1	D	257	LYS
1	D	260	LYS
1	D	272	ARG
1	D	273	ARG
1	D	277	SER
1	D	282	LYS
1	D	283	ARG
1	D	288	LYS
1	E	34	GLN
1	E	43	LYS
1	E	65	LEU
1	E	71	GLN
1	E	101	LYS
1	E	106	ARG
1	E	156	GLU
1	E	157	LYS
1	E	183	VAL
1	E	205	GLN

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Mol	Chain	Res	Type
1	E	210	SER
1	E	257	LYS
1	E	260	LYS
1	E	272	ARG
1	E	273	ARG
1	E	284	LYS
1	F	32	ASN
1	F	41	LYS
1	F	65	LEU
1	F	71	GLN
1	F	92	LYS
1	F	101	LYS
1	F	103	MET
1	F	106	ARG
1	F	156	GLU
1	F	157	LYS
1	F	183	VAL
1	F	201	GLN
1	F	210	SER
1	F	257	LYS
1	F	260	LYS
1	F	273	ARG
1	F	282	LYS
1	F	288	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (37) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	32	ASN
1	A	34	GLN
1	A	44	ASN
1	A	104	GLN
1	A	159	GLN
1	A	162	GLN
1	A	201	GLN
1	A	205	GLN
1	A	224	GLN
1	B	104	GLN
1	B	159	GLN
1	B	162	GLN
1	B	201	GLN
1	B	205	GLN

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Mol	Chain	Res	Type
1	B	224	GLN
1	C	59	ASN
1	C	159	GLN
1	C	162	GLN
1	C	201	GLN
1	C	205	GLN
1	D	159	GLN
1	D	162	GLN
1	D	205	GLN
1	D	224	GLN
1	E	122	HIS
1	E	159	GLN
1	E	162	GLN
1	E	201	GLN
1	E	224	GLN
1	F	32	ASN
1	F	104	GLN
1	F	105	ASN
1	F	159	GLN
1	F	162	GLN
1	F	201	GLN
1	F	205	GLN
1	F	224	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

4 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	CO8	A	300	-	49,59,59	1.34	4 (8%)	61,85,85	2.31	15 (24%)
2	CO8	B	300	1	49,59,59	1.34	4 (8%)	61,85,85	2.31	15 (24%)
2	CO8	E	300	-	49,59,59	1.34	4 (8%)	61,85,85	2.31	15 (24%)
2	CO8	F	300	-	49,59,59	1.32	4 (8%)	61,85,85	2.35	15 (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
2	CO8	A	300	-	-	0/54/74/74	0/3/3/3
2	CO8	B	300	1	-	0/54/74/74	0/3/3/3
2	CO8	E	300	-	-	0/54/74/74	0/3/3/3
2	CO8	F	300	-	-	0/54/74/74	0/3/3/3

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	300	CO8	O4B-C1B	2.85	1.44	1.41
2	A	300	CO8	O4B-C1B	2.88	1.44	1.41
2	E	300	CO8	O4B-C1B	2.93	1.44	1.41
2	B	300	CO8	O4B-C1B	2.94	1.44	1.41
2	A	300	CO8	C4A-N3A	3.44	1.40	1.35
2	E	300	CO8	C4A-N3A	3.46	1.40	1.35
2	B	300	CO8	C4A-N3A	3.47	1.40	1.35
2	F	300	CO8	C4A-N3A	3.84	1.41	1.35
2	F	300	CO8	C5P-N4P	3.90	1.42	1.33
2	B	300	CO8	C9P-N8P	4.26	1.42	1.33
2	A	300	CO8	C9P-N8P	4.28	1.42	1.33
2	E	300	CO8	C9P-N8P	4.28	1.42	1.33
2	B	300	CO8	C5P-N4P	4.28	1.43	1.33
2	E	300	CO8	C5P-N4P	4.29	1.43	1.33
2	F	300	CO8	C9P-N8P	4.30	1.42	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	300	CO8	C5P-N4P	4.30	1.43	1.33

All (60) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	300	CO8	O1'-C1'-S1P	-7.58	116.82	122.83
2	A	300	CO8	O1'-C1'-S1P	-7.29	117.05	122.83
2	B	300	CO8	O1'-C1'-S1P	-7.28	117.05	122.83
2	E	300	CO8	O1'-C1'-S1P	-7.26	117.08	122.83
2	E	300	CO8	O1'-C1'-C2'	-6.29	119.61	123.94
2	B	300	CO8	O1'-C1'-C2'	-6.29	119.61	123.94
2	A	300	CO8	O1'-C1'-C2'	-6.22	119.66	123.94
2	F	300	CO8	O1'-C1'-C2'	-5.70	120.02	123.94
2	F	300	CO8	C2B-C1B-N9A	-4.37	107.61	114.29
2	E	300	CO8	C2B-C1B-N9A	-3.62	108.76	114.29
2	A	300	CO8	C2B-C1B-N9A	-3.62	108.77	114.29
2	B	300	CO8	C2B-C1B-N9A	-3.61	108.78	114.29
2	F	300	CO8	OAP-CAP-C9P	-2.53	104.59	110.38
2	F	300	CO8	C6P-C5P-N4P	-2.50	112.11	116.46
2	A	300	CO8	C6P-C5P-N4P	-2.31	112.44	116.46
2	B	300	CO8	C6P-C5P-N4P	-2.29	112.48	116.46
2	E	300	CO8	C6P-C5P-N4P	-2.29	112.49	116.46
2	A	300	CO8	C2B-C3B-C4B	-2.18	99.20	103.29
2	B	300	CO8	OAP-CAP-C9P	-2.17	105.40	110.38
2	B	300	CO8	C2B-C3B-C4B	-2.17	99.22	103.29
2	E	300	CO8	C2B-C3B-C4B	-2.17	99.22	103.29
2	A	300	CO8	OAP-CAP-C9P	-2.16	105.41	110.38
2	E	300	CO8	OAP-CAP-C9P	-2.15	105.44	110.38
2	F	300	CO8	C2B-C3B-C4B	-2.10	99.35	103.29
2	F	300	CO8	C5'-C4'-C3'	2.01	124.93	114.53
2	B	300	CO8	C5'-C4'-C3'	2.02	124.98	114.53
2	E	300	CO8	C5'-C4'-C3'	2.02	124.98	114.53
2	A	300	CO8	C5'-C4'-C3'	2.03	125.00	114.53
2	B	300	CO8	O4B-C1B-N9A	2.04	112.37	108.10
2	A	300	CO8	O4B-C1B-N9A	2.04	112.38	108.10
2	E	300	CO8	O4B-C1B-N9A	2.05	112.39	108.10
2	A	300	CO8	O3A-P1A-O5B	2.06	108.39	102.94
2	E	300	CO8	O3A-P1A-O5B	2.06	108.40	102.94
2	B	300	CO8	O3A-P1A-O5B	2.06	108.41	102.94
2	A	300	CO8	O3A-P2A-O6A	2.18	108.72	102.94
2	E	300	CO8	O3A-P2A-O6A	2.18	108.72	102.94
2	B	300	CO8	O3A-P2A-O6A	2.19	108.76	102.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	300	CO8	C4A-C5A-N7A	2.22	111.52	109.48
2	E	300	CO8	C4A-C5A-N7A	2.22	111.52	109.48
2	A	300	CO8	C4A-C5A-N7A	2.24	111.54	109.48
2	A	300	CO8	O6A-CCP-CBP	2.27	114.19	110.55
2	B	300	CO8	O6A-CCP-CBP	2.27	114.19	110.55
2	E	300	CO8	O6A-CCP-CBP	2.28	114.21	110.55
2	F	300	CO8	C4A-C5A-N7A	2.28	111.58	109.48
2	F	300	CO8	O3A-P1A-O5B	2.34	109.15	102.94
2	F	300	CO8	O4B-C1B-N9A	2.40	113.12	108.10
2	F	300	CO8	O6A-CCP-CBP	2.48	114.54	110.55
2	F	300	CO8	O3A-P2A-O6A	2.63	109.90	102.94
2	F	300	CO8	C4'-C3'-C2'	2.87	123.81	113.29
2	E	300	CO8	C4'-C3'-C2'	2.95	124.12	113.29
2	A	300	CO8	C4'-C3'-C2'	2.96	124.13	113.29
2	B	300	CO8	C4'-C3'-C2'	2.96	124.14	113.29
2	A	300	CO8	C3'-C2'-C1'	6.22	122.30	113.12
2	B	300	CO8	C3'-C2'-C1'	6.24	122.33	113.12
2	E	300	CO8	C3'-C2'-C1'	6.25	122.34	113.12
2	F	300	CO8	C3'-C2'-C1'	6.39	122.55	113.12
2	F	300	CO8	C2'-C1'-S1P	9.94	122.30	113.36
2	A	300	CO8	C2'-C1'-S1P	10.14	122.48	113.36
2	E	300	CO8	C2'-C1'-S1P	10.16	122.50	113.36
2	B	300	CO8	C2'-C1'-S1P	10.17	122.50	113.36

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 20 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	300	CO8	6	0
2	B	300	CO8	7	0
2	E	300	CO8	6	0
2	F	300	CO8	1	0

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 ⓘ

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	254/261 (97%)	-0.14	7 (2%) 56 55	12, 29, 52, 71	0
1	B	254/261 (97%)	-0.18	5 (1%) 68 68	12, 28, 52, 62	0
1	C	254/261 (97%)	-0.05	12 (4%) 35 36	12, 32, 54, 64	0
1	D	257/261 (98%)	-0.23	6 (2%) 64 63	10, 29, 53, 84	0
1	E	255/261 (97%)	-0.23	4 (1%) 74 74	11, 28, 53, 76	0
1	F	255/261 (97%)	-0.24	2 (0%) 87 87	13, 27, 53, 74	0
All	All	1529/1566 (97%)	-0.18	36 (2%) 62 61	10, 29, 53, 84	0

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	31	ALA	6.3
1	D	31	ALA	5.5
1	D	115	LYS	4.7
1	F	31	ALA	4.6
1	A	43	LYS	4.5
1	B	56	LYS	3.8
1	D	114	GLY	3.3
1	C	43	LYS	3.2
1	D	32	ASN	3.0
1	A	44	ASN	3.0
1	A	280	VAL	3.0
1	C	113	SER	2.9
1	C	280	VAL	2.9
1	C	45	SER	2.9
1	E	43	LYS	2.8
1	B	125	ARG	2.8
1	A	282	LYS	2.8
1	A	32	ASN	2.7
1	B	106	ARG	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	32	ASN	2.6
1	C	112	TYR	2.5
1	C	55	PRO	2.5
1	C	57	ALA	2.5
1	B	34	GLN	2.4
1	C	56	LYS	2.4
1	D	284	LYS	2.4
1	D	280	VAL	2.4
1	E	44	ASN	2.4
1	F	32	ASN	2.2
1	C	284	LYS	2.2
1	A	285	ALA	2.2
1	A	132	ALA	2.2
1	C	286	ASN	2.1
1	C	137	TYR	2.1
1	E	32	ASN	2.1
1	C	33	PHE	2.0

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
2	CO8	B	300	57/57	0.51	0.55	4.66	40,60,76,79	0
2	CO8	F	300	57/57	0.50	0.40	4.05	45,65,89,92	0
2	CO8	A	300	57/57	0.87	0.21	0.99	40,60,76,79	0
2	CO8	E	300	57/57	0.89	0.18	0.74	40,60,76,79	0

6.5 Other polymers [i](#)

There are no such residues in this entry.