



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 12:37 AM GMT

PDB ID : 2B37
Title : Crystal structure of Mycobacterium tuberculosis enoyl reductase (InhA) inhibited by 5-octyl-2-phenoxyphenol
Authors : Sullivan, T.J.; Truglio, J.J.; Novichenok, P.; Stratton, C.; Zhang, X.; Kaur, T.; Johnson, F.; Boyne, M.S.; Amin, A.
Deposited on : 2005-09-19
Resolution : 2.60 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

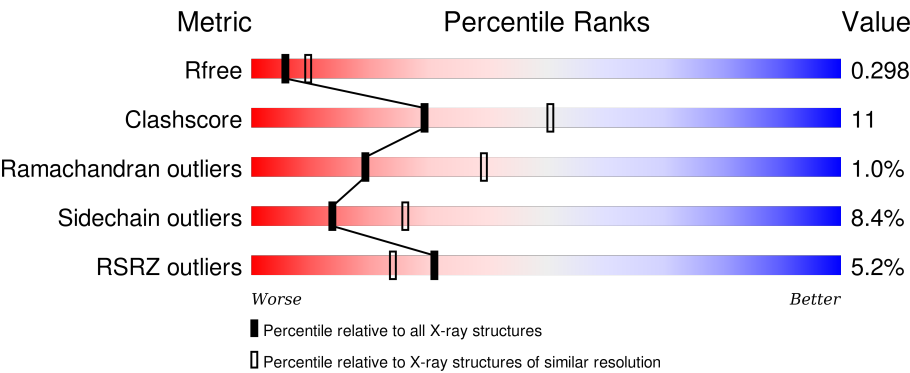
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	2328 (2.60-2.60)
Clashscore	102246	2679 (2.60-2.60)
Ramachandran outliers	100387	2635 (2.60-2.60)
Sidechain outliers	100360	2635 (2.60-2.60)
RSRZ outliers	91569	2334 (2.60-2.60)

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 <=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	269	<div><div>4%</div><div>67%23%8%</div></div>
1	B	269	<div><div>7%</div><div>68%21%9%</div></div>
1	C	269	<div><div>2%</div><div>74%19%</div></div>
1	D	269	<div><div>%</div><div>75%18%</div></div>
1	E	269	<div><div>7%</div><div>68%24%</div></div>

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Mol	Chain	Length	Quality of chain
1	F	269	 A horizontal bar chart showing the quality of chain F. The bar is divided into four segments: a small red segment at the beginning labeled '7%', a large green segment labeled '62%', a yellow segment labeled '27%', and a small grey segment at the end labeled '8%'. There is a small black dot on the yellow segment.

2 Entry composition [i](#)

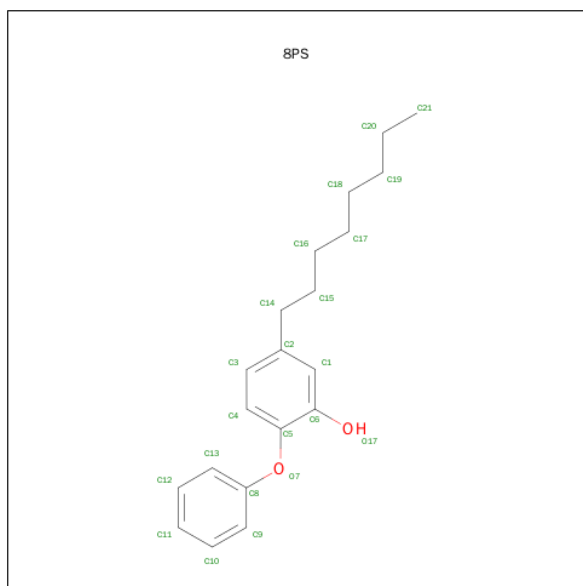
There are 4 unique types of molecules in this entry. The entry contains 11736 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Enoyl-[acyl-carrier-protein] reductase [NADH].

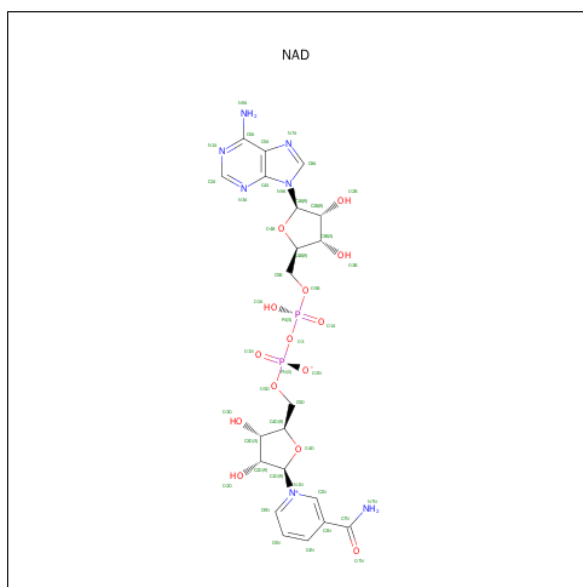
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	248	Total	C	N	O	S	0	0	0
			1865	1184	326	346	9			
1	B	245	Total	C	N	O	S	0	0	0
			1840	1167	323	341	9			
1	C	260	Total	C	N	O	S	0	0	0
			1944	1231	340	364	9			
1	D	260	Total	C	N	O	S	0	0	0
			1944	1231	340	364	9			
1	E	260	Total	C	N	O	S	0	0	0
			1944	1231	340	364	9			
1	F	248	Total	C	N	O	S	0	0	0
			1861	1181	326	344	10			

- Molecule 2 is 5-OCTYL-2-PHENOXYPHENOL (three-letter code: 8PS) (formula: $C_{20}H_{26}O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	C	1	Total	C	O	0	0
			22	20	2		
2	D	1	Total	C	O	0	0
			22	20	2		

- Molecule 3 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: C₂₁H₂₇N₇O₁₄P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
3	B	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
3	C	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
3	D	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
3	E	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
3	F	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	7	Total	O	0	0
			7	7		

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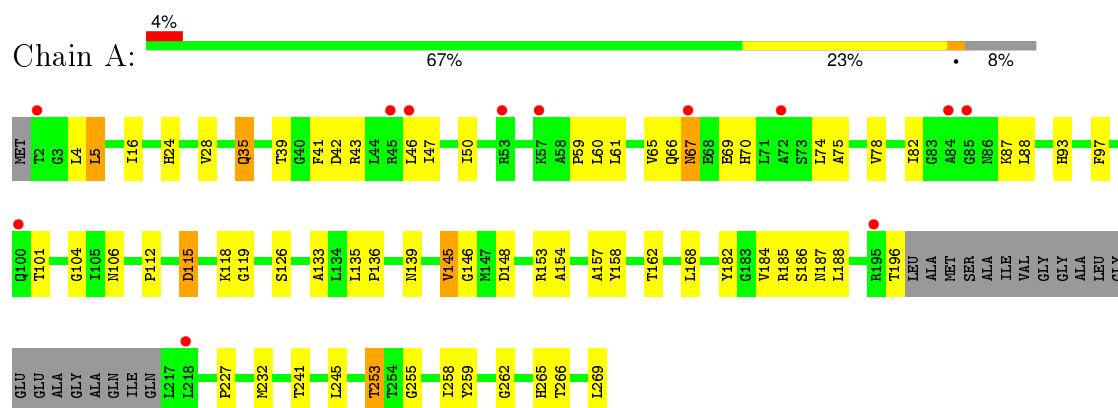
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	2	Total 2	O 2	0	0
4	C	7	Total 7	O 7	0	0
4	D	6	Total 6	O 6	0	0
4	E	3	Total 3	O 3	0	0
4	F	5	Total 5	O 5	0	0

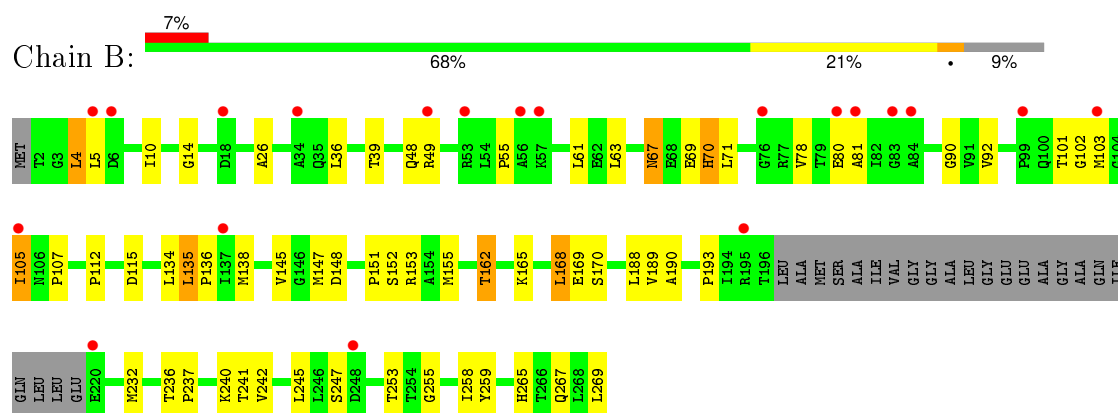
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

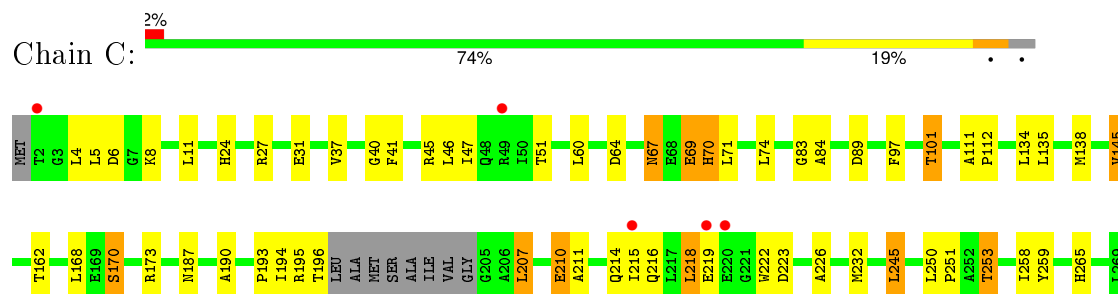
- Molecule 1: Enoyl-[acyl-carrier-protein] reductase [NADH]




- Molecule 1: Enoyl-[acyl-carrier-protein] reductase [NADH]

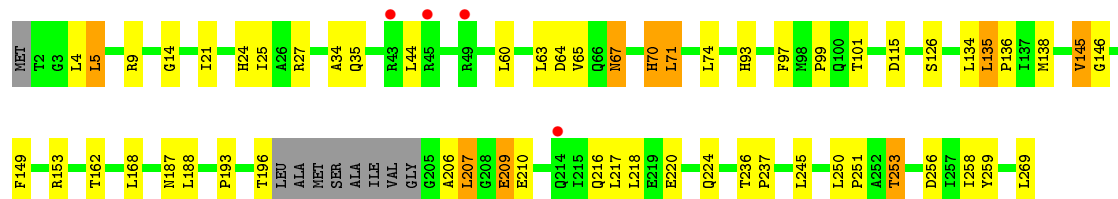


- Molecule 1: Enoyl-[acyl-carrier-protein] reductase [NADH]



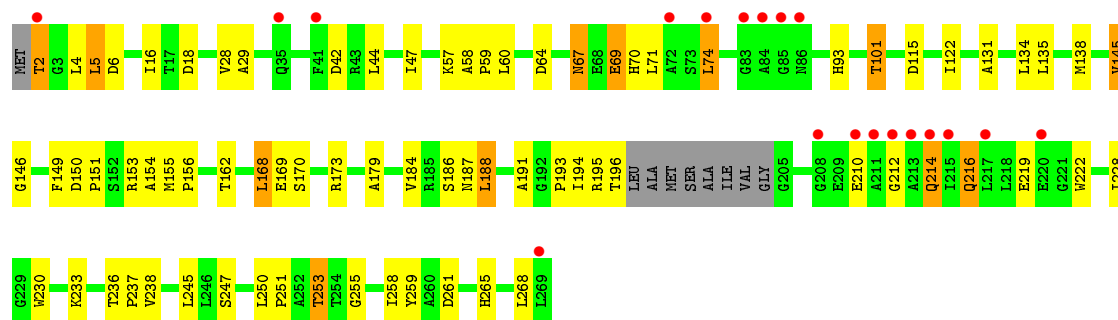
- Molecule 1: Enoyl-[acyl-carrier-protein] reductase [NADH]

Chain D: 



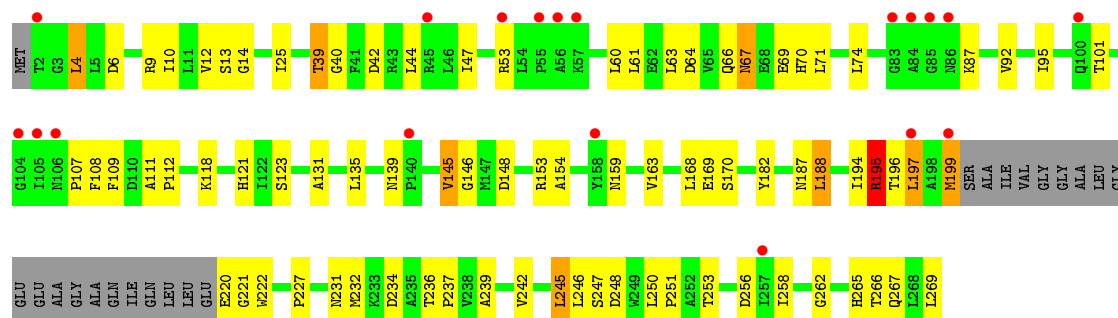
- Molecule 1: Enoyl-[acyl-carrier-protein] reductase [NADH]

Chain E: 



- Molecule 1: Enoyl-[acyl-carrier-protein] reductase [NADH]

Chain F: 



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	82.34Å 100.29Å 379.27Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.00 – 2.60 39.29 – 2.60	Depositor EDS
% Data completeness (in resolution range)	96.4 (15.00-2.60) 96.3 (39.29-2.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.97 (at 2.61Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.225 , 0.294 0.230 , 0.298	Depositor DCC
R_{free} test set	2370 reflections (5.34%)	DCC
Wilson B-factor (Å ²)	34.7	Xtriage
Anisotropy	0.817	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 50.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.42$, $\langle L^2 \rangle = 0.25$	Xtriage
Outliers	0 of 47028 reflections	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	11736	wwPDB-VP
Average B, all atoms (Å ²)	40.0	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 30.88 % 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 1.2140e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical 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: 8PS, NAD

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	0.67	0/1902	0.62	0/2583
1	B	0.70	0/1877	0.62	0/2549
1	C	0.84	1/1981 (0.1%)	0.70	0/2689
1	D	0.85	0/1981	0.70	0/2689
1	E	0.78	0/1981	0.66	0/2689
1	F	0.73	0/1898	0.64	1/2577 (0.0%)
All	All	0.77	1/11620 (0.0%)	0.65	1/15776 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	222	TRP	CB-CG	-5.46	1.40	1.50

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	195	ARG	NE-CZ-NH1	5.11	122.86	120.30

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	1865	0	1880	48	0
1	B	1840	0	1852	44	0
1	C	1944	0	1954	40	0
1	D	1944	0	1954	43	0
1	E	1944	0	1954	55	0
1	F	1861	0	1877	59	0
2	C	22	0	25	3	0
2	D	22	0	25	7	0
3	A	44	0	26	1	0
3	B	44	0	26	1	0
3	C	44	0	26	2	0
3	D	44	0	26	5	0
3	E	44	0	26	4	0
3	F	44	0	26	2	0
4	A	7	0	0	0	0
4	B	2	0	0	0	0
4	C	7	0	0	3	0
4	D	6	0	0	0	0
4	E	3	0	0	1	0
4	F	5	0	0	0	0
All	All	11736	0	11677	265	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 (265) 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:F:101:THR:HG21	1:F:112:PRO:HD2	1.57	0.85
1:F:195:ARG:HH11	1:F:195:ARG:HG2	1.47	0.79
1:E:168:LEU:HD22	1:E:188:LEU:HD21	1.65	0.78
1:A:35:GLN:HE22	1:A:82:ILE:HG22	1.51	0.76
1:E:4:LEU:HD13	1:E:5:LEU:HD13	1.67	0.75
1:D:101:THR:HG21	1:D:115:ASP:OD1	1.85	0.74
1:A:241:THR:HG23	1:E:250:LEU:HD23	1.68	0.73
1:E:245:LEU:HD11	1:E:258:ILE:HD12	1.70	0.73
1:F:195:ARG:CG	1:F:195:ARG:HH11	2.02	0.71
1:C:218:LEU:HB2	2:C:300:8PS:H212	1.72	0.71
1:F:101:THR:CG2	1:F:111:ALA:HA	2.21	0.70
1:A:101:THR:CG2	1:A:112:PRO:HD2	2.23	0.68
1:A:101:THR:HG21	1:A:112:PRO:HD2	1.73	0.68
1:D:67:ASN:HD22	1:D:67:ASN:C	1.98	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:153:ARG:NH1	1:F:265:HIS:O	2.28	0.66
1:A:153:ARG:HH21	1:F:153:ARG:NE	1.94	0.65
1:D:218:LEU:HD11	2:D:301:8PS:H192	1.78	0.65
1:D:206:ALA:HA	1:D:209:GLU:HG2	1.80	0.64
1:F:168:LEU:HD23	1:F:188:LEU:HD21	1.80	0.63
1:F:168:LEU:CD2	1:F:188:LEU:HD21	2.29	0.63
1:B:101:THR:HG21	1:B:112:PRO:HD2	1.81	0.62
1:D:218:LEU:HD11	2:D:301:8PS:C19	2.30	0.62
1:E:212:GLY:O	1:E:216:GLN:HB2	1.99	0.62
1:F:196:THR:HG23	1:F:197:LEU:HD23	1.80	0.62
1:A:259:TYR:O	1:E:253:THR:HB	2.00	0.62
1:E:173:ARG:HB3	1:F:154:ALA:HB2	1.82	0.62
1:A:253:THR:HB	1:E:259:TYR:O	2.00	0.62
1:D:67:ASN:ND2	1:D:70:HIS:H	1.98	0.61
4:C:408:HOH:O	1:D:253:THR:HG22	2.01	0.61
1:E:67:ASN:HD21	1:E:69:GLU:HB2	1.64	0.61
1:A:153:ARG:HH21	1:F:153:ARG:HE	1.47	0.61
1:D:99:PRO:HB2	1:D:101:THR:HG22	1.82	0.61
1:A:158:TYR:HD2	1:A:162:THR:HG1	1.44	0.61
1:B:148:ASP:O	1:B:190:ALA:HA	2.01	0.60
1:F:123:SER:O	1:F:168:LEU:HD12	2.01	0.60
1:E:210:GLU:O	1:E:214:GLN:NE2	2.34	0.60
1:F:253:THR:HG21	1:F:258:ILE:HD11	1.83	0.60
1:B:255:GLY:O	1:F:266:THR:HG21	2.02	0.60
1:D:218:LEU:HG	2:D:301:8PS:H213	1.83	0.59
1:C:245:LEU:HD11	1:C:258:ILE:HD13	1.83	0.59
1:A:60:LEU:HD23	1:A:61:LEU:N	2.17	0.59
1:A:265:HIS:O	1:F:153:ARG:NH1	2.35	0.59
1:D:245:LEU:HD11	1:D:258:ILE:HD13	1.84	0.59
1:E:101:THR:HG21	1:E:115:ASP:OD1	2.02	0.59
1:D:218:LEU:HD21	2:D:301:8PS:H201	1.84	0.59
1:B:14:GLY:O	3:B:402:NAD:O3B	2.20	0.59
1:A:241:THR:HG23	1:E:250:LEU:CD2	2.33	0.58
1:F:131:ALA:O	1:F:135:LEU:HB2	2.03	0.58
1:B:101:THR:HG21	1:B:115:ASP:OD1	2.03	0.58
1:E:145:VAL:HA	1:E:187:ASN:O	2.04	0.58
1:F:40:GLY:HA3	1:F:47:ILE:CD1	2.33	0.58
1:F:145:VAL:HA	1:F:187:ASN:O	2.04	0.57
1:B:67:ASN:HD22	1:B:67:ASN:C	2.07	0.57
1:A:101:THR:O	1:A:106:ASN:ND2	2.38	0.57
1:A:66:GLN:HE22	1:A:118:LYS:HG3	1.69	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:195:ARG:N	1:F:232:MET:O	2.32	0.57
1:C:134:LEU:O	1:C:138:MET:HG3	2.05	0.57
1:F:145:VAL:HG21	1:F:242:VAL:HG13	1.87	0.56
1:D:4:LEU:HB3	1:D:5:LEU:HD13	1.86	0.56
1:A:24:HIS:O	1:A:28:VAL:HG23	2.05	0.56
1:E:131:ALA:O	1:E:135:LEU:HB2	2.05	0.56
1:A:87:LYS:HB2	1:A:139:ASN:ND2	2.21	0.56
1:D:21:ILE:O	1:D:25:ILE:HG13	2.06	0.56
1:B:67:ASN:ND2	1:B:70:HIS:H	2.04	0.55
1:F:95:ILE:HD13	3:F:406:NAD:C2A	2.36	0.55
1:B:267:GLN:OE1	1:E:154:ALA:HB3	2.07	0.55
1:F:14:GLY:O	3:F:406:NAD:O3B	2.24	0.55
1:F:60:LEU:O	1:F:60:LEU:HD13	2.07	0.55
1:C:259:TYR:O	1:D:253:THR:HB	2.07	0.55
1:C:253:THR:HB	1:D:259:TYR:O	2.07	0.55
1:A:158:TYR:HD2	1:A:162:THR:OG1	1.90	0.54
1:B:193:PRO:O	1:B:232:MET:HB3	2.08	0.54
1:F:220:GLU:O	1:F:222:TRP:N	2.40	0.54
1:B:147:MET:HE1	1:B:189:VAL:HB	1.89	0.54
1:B:78:VAL:O	1:B:81:ALA:N	2.40	0.54
1:C:196:THR:HG21	3:C:403:NAD:PN	2.47	0.54
1:C:64:ASP:H	1:C:70:HIS:HD2	1.54	0.54
1:E:196:THR:O	4:E:407:HOH:O	2.18	0.54
1:D:60:LEU:HD13	1:D:60:LEU:O	2.08	0.53
1:A:154:ALA:HB3	1:F:267:GLN:OE1	2.08	0.53
1:F:4:LEU:HD22	1:F:247:SER:HB2	1.89	0.52
1:B:265:HIS:O	1:E:153:ARG:NH1	2.42	0.52
1:A:182:TYR:HB2	1:A:184:VAL:HG23	1.91	0.52
1:E:214:GLN:N	1:E:214:GLN:HE21	2.07	0.52
1:E:214:GLN:HE21	1:E:214:GLN:H	1.58	0.52
1:D:63:LEU:HD11	1:D:71:LEU:CD1	2.39	0.52
1:E:16:ILE:HD12	1:E:47:ILE:CG2	2.39	0.52
1:B:147:MET:HE3	1:B:242:VAL:HG22	1.92	0.52
1:C:8:LYS:HA	1:C:89:ASP:OD2	2.10	0.51
1:E:191:ALA:HB1	1:E:238:VAL:HG21	1.92	0.51
1:F:195:ARG:NH1	1:F:195:ARG:HG2	2.16	0.51
1:E:134:LEU:O	1:E:138:MET:HG3	2.11	0.51
1:D:14:GLY:O	3:D:404:NAD:O3B	2.23	0.51
1:F:135:LEU:HG	1:F:182:TYR:CD1	2.45	0.51
1:E:2:THR:HG22	1:E:6:ASP:OD2	2.11	0.50
1:C:210:GLU:CD	1:C:210:GLU:H	2.15	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:223:ASP:O	4:C:407:HOH:O	2.19	0.50
1:C:216:GLN:HA	1:C:219:GLU:CD	2.32	0.50
1:C:64:ASP:H	1:C:70:HIS:CD2	2.30	0.50
1:C:67:ASN:HD22	1:C:69:GLU:H	1.58	0.50
1:B:241:THR:HG23	1:F:250:LEU:HD23	1.94	0.49
1:A:145:VAL:HA	1:A:187:ASN:O	2.12	0.49
1:C:101:THR:HG23	1:C:111:ALA:HA	1.93	0.49
1:B:190:ALA:HB3	1:B:259:TYR:CD2	2.48	0.49
1:C:101:THR:HG21	1:C:112:PRO:HD2	1.95	0.49
1:D:24:HIS:CD2	1:D:27:ARG:HH21	2.30	0.49
1:D:269:LEU:C	1:D:269:LEU:HD13	2.33	0.49
1:B:4:LEU:HD11	1:B:247:SER:HB3	1.95	0.49
1:F:231:ASN:HD22	1:F:234:ASP:HB2	1.78	0.49
1:F:107:PRO:O	1:F:109:PHE:N	2.46	0.49
1:F:25:ILE:HG12	1:F:239:ALA:HA	1.95	0.48
1:B:147:MET:O	1:B:165:LYS:NZ	2.46	0.48
1:A:227:PRO:HD2	1:A:262:GLY:O	2.14	0.48
1:B:26:ALA:HB1	1:B:36:LEU:CD1	2.44	0.48
1:A:67:ASN:HD22	1:A:70:HIS:H	1.62	0.48
1:C:27:ARG:HD2	1:C:31:GLU:OE2	2.13	0.48
1:E:155:MET:HB2	1:E:156:PRO:HD2	1.95	0.48
1:D:196:THR:HG21	3:D:404:NAD:O1N	2.13	0.48
1:D:65:VAL:HG22	3:D:404:NAD:C6A	2.44	0.48
1:E:210:GLU:CD	1:E:210:GLU:H	2.16	0.48
1:D:65:VAL:HG22	3:D:404:NAD:N1A	2.28	0.48
1:E:236:THR:HB	1:E:237:PRO:HD3	1.96	0.48
1:A:186:SER:O	1:A:255:GLY:N	2.46	0.48
1:B:4:LEU:HD13	1:B:5:LEU:CD1	2.44	0.47
1:D:145:VAL:HA	1:D:187:ASN:O	2.14	0.47
1:E:193:PRO:HA	3:E:405:NAD:O7N	2.13	0.47
1:B:148:ASP:OD2	1:B:169:GLU:OE2	2.33	0.47
1:B:105:ILE:O	1:B:105:ILE:HD13	2.14	0.47
1:C:218:LEU:CB	2:C:300:8PS:H212	2.43	0.47
1:E:250:LEU:N	1:E:251:PRO:HD3	2.30	0.47
1:A:241:THR:O	1:A:245:LEU:HD23	2.14	0.47
1:B:153:ARG:NH1	1:E:265:HIS:O	2.48	0.47
1:C:250:LEU:N	1:C:251:PRO:HD3	2.30	0.47
1:E:93:HIS:O	1:E:146:GLY:HA2	2.14	0.47
1:C:207:LEU:HB2	1:C:211:ALA:CB	2.44	0.47
1:C:245:LEU:CD1	1:C:250:LEU:HD13	2.44	0.47
1:D:149:PHE:CE2	2:D:301:8PS:H142	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:265:HIS:NE2	1:D:256:ASP:OD2	2.28	0.47
1:B:236:THR:O	1:B:240:LYS:N	2.39	0.47
1:F:101:THR:HG21	1:F:112:PRO:CD	2.35	0.46
1:A:245:LEU:HD11	1:A:258:ILE:HD13	1.97	0.46
1:E:67:ASN:ND2	1:E:69:GLU:HB2	2.31	0.46
1:D:24:HIS:HD2	1:D:27:ARG:HH21	1.62	0.46
1:C:207:LEU:HB2	1:C:211:ALA:HB3	1.98	0.46
1:F:66:GLN:HE22	1:F:118:LYS:HG3	1.80	0.46
1:A:16:ILE:HD12	1:A:47:ILE:CG2	2.44	0.46
1:D:93:HIS:O	1:D:146:GLY:HA2	2.15	0.46
1:E:67:ASN:HD22	1:E:67:ASN:C	2.18	0.46
1:E:58:ALA:HB1	1:E:59:PRO:HD2	1.97	0.46
1:E:28:VAL:O	1:E:29:ALA:C	2.55	0.45
1:B:151:PRO:CB	1:B:162:THR:HG22	2.46	0.45
1:A:185:ARG:HG2	1:E:228:ILE:HD13	1.97	0.45
1:F:197:LEU:HD23	1:F:197:LEU:H	1.82	0.45
1:A:93:HIS:O	1:A:146:GLY:HA2	2.16	0.45
1:B:153:ARG:HH21	1:E:153:ARG:NE	2.13	0.45
1:C:24:HIS:HD2	1:C:27:ARG:HH21	1.65	0.45
1:B:39:THR:HB	1:B:63:LEU:HB3	1.99	0.45
1:D:5:LEU:HB3	1:D:34:ALA:HB2	1.97	0.45
1:A:60:LEU:C	1:A:60:LEU:HD23	2.37	0.45
1:A:66:GLN:NE2	1:A:118:LYS:HG3	2.31	0.45
1:E:74:LEU:HD13	1:E:134:LEU:HD21	1.98	0.45
1:A:97:PHE:O	1:A:119:GLY:HA2	2.16	0.45
1:A:148:ASP:O	3:A:401:NAD:H6N	2.17	0.45
1:C:40:GLY:HA3	1:C:47:ILE:HD11	1.98	0.45
1:B:253:THR:HG21	1:B:258:ILE:HD11	1.99	0.45
1:F:227:PRO:HD2	1:F:262:GLY:O	2.16	0.44
1:F:12:VAL:HA	1:F:92:VAL:HB	1.98	0.44
1:E:122:ILE:HD13	3:E:405:NAD:H61A	1.82	0.44
1:B:236:THR:N	1:B:237:PRO:HD2	2.32	0.44
1:D:64:ASP:H	1:D:70:HIS:HD2	1.64	0.44
1:A:153:ARG:HD3	1:F:265:HIS:O	2.17	0.44
1:D:206:ALA:O	1:D:207:LEU:C	2.55	0.44
1:E:64:ASP:H	1:E:70:HIS:CD2	2.35	0.44
1:F:168:LEU:HD23	1:F:168:LEU:C	2.38	0.44
1:B:134:LEU:O	1:B:138:MET:HG3	2.16	0.44
1:F:39:THR:HA	1:F:61:LEU:O	2.17	0.44
1:C:218:LEU:HD23	1:C:219:GLU:N	2.33	0.44
1:F:66:GLN:HG2	1:F:121:HIS:CE1	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:245:LEU:HD11	1:E:258:ILE:CD1	2.42	0.44
1:A:265:HIS:CD2	1:A:266:THR:HG23	2.52	0.44
1:A:16:ILE:O	1:A:50:ILE:HD12	2.17	0.44
1:B:102:GLY:O	1:B:103:MET:HG2	2.18	0.44
1:E:179:ALA:HB1	1:E:184:VAL:HB	2.00	0.44
1:D:65:VAL:HB	1:D:126:SER:HB2	2.00	0.44
1:C:219:GLU:HG2	1:C:232:MET:SD	2.58	0.43
1:D:218:LEU:HD11	2:D:301:8PS:C20	2.48	0.43
1:C:194:ILE:O	1:C:196:THR:N	2.51	0.43
1:A:75:ALA:HB2	1:A:133:ALA:O	2.18	0.43
1:B:10:ILE:HG13	1:B:90:GLY:HA3	2.00	0.43
1:C:97:PHE:HA	2:C:300:8PS:H11	2.00	0.43
1:B:92:VAL:HG22	1:B:145:VAL:CG1	2.47	0.43
1:D:236:THR:N	1:D:237:PRO:CD	2.80	0.43
1:C:83:GLY:O	1:C:84:ALA:C	2.56	0.43
1:F:195:ARG:CG	1:F:195:ARG:NH1	2.71	0.43
1:F:194:ILE:HA	1:F:232:MET:O	2.18	0.43
1:C:190:ALA:HB3	1:C:259:TYR:CD2	2.54	0.43
1:F:64:ASP:H	1:F:70:HIS:HD2	1.65	0.43
1:A:43:ARG:HE	1:A:46:LEU:HD22	1.82	0.43
1:D:135:LEU:N	1:D:136:PRO:CD	2.81	0.43
1:C:170:SER:HA	1:C:173:ARG:NH1	2.33	0.43
1:B:67:ASN:ND2	1:B:69:GLU:H	2.17	0.43
1:C:41:PHE:HB2	3:C:403:NAD:C2A	2.48	0.43
1:D:250:LEU:N	1:D:251:PRO:HD3	2.34	0.43
1:C:145:VAL:HA	1:C:187:ASN:O	2.18	0.43
1:A:41:PHE:CG	1:A:42:ASP:N	2.87	0.43
1:F:236:THR:N	1:F:237:PRO:CD	2.82	0.43
1:A:65:VAL:HB	1:A:126:SER:HB2	1.99	0.43
1:B:168:LEU:HD22	1:B:168:LEU:O	2.19	0.43
1:D:153:ARG:HG2	1:D:153:ARG:HH11	1.83	0.43
1:E:4:LEU:CD1	1:E:5:LEU:HD13	2.44	0.42
1:B:67:ASN:ND2	1:B:69:GLU:N	2.66	0.42
1:C:193:PRO:O	1:C:194:ILE:HD13	2.19	0.42
1:E:67:ASN:HD22	1:E:69:GLU:H	1.67	0.42
1:A:158:TYR:CD2	1:A:162:THR:OG1	2.61	0.42
1:C:67:ASN:HD22	1:C:69:GLU:N	2.17	0.42
1:C:47:ILE:O	1:C:51:THR:HG23	2.19	0.42
1:F:146:GLY:O	1:F:188:LEU:HA	2.19	0.42
1:B:4:LEU:CD1	1:B:247:SER:HB3	2.50	0.42
1:E:194:ILE:N	3:E:405:NAD:O7N	2.42	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:67:ASN:ND2	1:B:67:ASN:C	2.72	0.42
1:F:148:ASP:OD2	1:F:169:GLU:OE2	2.38	0.42
1:F:42:ASP:O	1:F:44:LEU:HD23	2.20	0.42
1:F:196:THR:HB	1:F:199:MET:HA	2.00	0.42
1:E:151:PRO:HD2	1:E:169:GLU:OE2	2.19	0.42
1:E:149:PHE:HB2	3:E:405:NAD:C5N	2.50	0.42
1:B:67:ASN:HD21	1:B:69:GLU:HB2	1.85	0.42
1:F:245:LEU:CD1	1:F:250:LEU:HD13	2.49	0.42
1:A:104:GLY:O	1:A:157:ALA:HA	2.19	0.42
1:E:42:ASP:OD1	1:E:42:ASP:C	2.58	0.42
1:E:250:LEU:N	1:E:251:PRO:CD	2.83	0.42
1:C:11:LEU:HA	1:C:37:VAL:O	2.20	0.42
1:E:186:SER:O	1:E:255:GLY:N	2.50	0.42
1:B:135:LEU:N	1:B:136:PRO:CD	2.83	0.42
1:A:101:THR:HB	1:A:115:ASP:OD1	2.19	0.41
1:F:250:LEU:N	1:F:251:PRO:HD3	2.35	0.41
1:B:39:THR:HA	1:B:61:LEU:O	2.20	0.41
1:B:155:MET:CE	1:E:268:LEU:HD22	2.49	0.41
1:C:214:GLN:O	1:C:215:ILE:C	2.54	0.41
1:D:193:PRO:HA	3:D:404:NAD:O7N	2.20	0.41
1:F:87:LYS:HB2	1:F:139:ASN:ND2	2.35	0.41
1:A:78:VAL:HG11	1:A:88:LEU:HD11	2.03	0.41
1:D:220:GLU:O	1:D:224:GLN:HG3	2.20	0.41
1:B:193:PRO:HB2	1:B:232:MET:SD	2.61	0.41
1:E:101:THR:HG21	1:E:115:ASP:CB	2.51	0.41
1:A:4:LEU:HB3	1:A:5:LEU:HD13	2.03	0.41
1:E:170:SER:HB3	1:F:163:VAL:HG22	2.03	0.41
1:E:4:LEU:HD11	1:E:247:SER:HB3	2.01	0.41
1:F:13:SER:HB2	1:F:95:ILE:HD11	2.03	0.41
1:F:220:GLU:C	1:F:222:TRP:N	2.74	0.41
1:A:67:ASN:HD21	1:A:69:GLU:HB2	1.85	0.41
1:F:39:THR:HB	1:F:63:LEU:HB3	2.02	0.41
1:E:230:TRP:CZ3	1:E:261:ASP:HA	2.55	0.41
1:F:10:ILE:HD13	1:F:246:LEU:HD13	2.03	0.41
1:C:226:ALA:O	4:C:407:HOH:O	2.22	0.41
1:B:165:LYS:HA	1:B:165:LYS:HD3	1.95	0.40
1:D:9:ARG:HA	1:D:35:GLN:O	2.21	0.40
1:D:97:PHE:HA	2:D:301:8PS:H11	2.04	0.40
1:B:147:MET:HE3	1:B:242:VAL:CG2	2.51	0.40
1:F:250:LEU:N	1:F:251:PRO:CD	2.85	0.40
1:F:67:ASN:ND2	1:F:69:GLU:H	2.19	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:135:LEU:HD21	1:E:179:ALA:HA	2.03	0.40
1:A:136:PRO:HG2	1:C:45:ARG:NH2	2.37	0.40
1:D:134:LEU:O	1:D:138:MET:HG3	2.22	0.40
1:D:216:GLN:C	1:D:218:LEU:N	2.75	0.40
1:A:39:THR:HA	1:A:61:LEU:O	2.22	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	244/269 (91%)	224 (92%)	18 (7%)	2 (1%)	24	46
1	B	241/269 (90%)	220 (91%)	19 (8%)	2 (1%)	24	46
1	C	256/269 (95%)	237 (93%)	18 (7%)	1 (0%)	39	65
1	D	256/269 (95%)	234 (91%)	18 (7%)	4 (2%)	12	24
1	E	256/269 (95%)	237 (93%)	17 (7%)	2 (1%)	24	46
1	F	244/269 (91%)	217 (89%)	23 (9%)	4 (2%)	12	24
All	All	1497/1614 (93%)	1369 (91%)	113 (8%)	15 (1%)	19	39

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	195	ARG
1	D	207	LEU
1	D	209	GLU
1	E	195	ARG
1	F	108	PHE
1	A	74	LEU
1	D	217	LEU

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Mol	Chain	Res	Type
1	F	221	GLY
1	A	59	PRO
1	F	53	ARG
1	B	55	PRO
1	E	150	ASP
1	F	159	ASN
1	D	210	GLU
1	B	107	PRO

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	193/205 (94%)	181 (94%)	12 (6%)	23	45
1	B	190/205 (93%)	174 (92%)	16 (8%)	14	26
1	C	199/205 (97%)	178 (89%)	21 (11%)	8	15
1	D	199/205 (97%)	187 (94%)	12 (6%)	24	47
1	E	199/205 (97%)	178 (89%)	21 (11%)	8	15
1	F	192/205 (94%)	175 (91%)	17 (9%)	12	23
All	All	1172/1230 (95%)	1073 (92%)	99 (8%)	14	26

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

Mol	Chain	Res	Type
1	A	5	LEU
1	A	35	GLN
1	A	67	ASN
1	A	115	ASP
1	A	135	LEU
1	A	145	VAL
1	A	168	LEU
1	A	188	LEU
1	A	196	THR
1	A	232	MET

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Mol	Chain	Res	Type
1	A	253	THR
1	A	269	LEU
1	B	4	LEU
1	B	48	GLN
1	B	49	ARG
1	B	67	ASN
1	B	70	HIS
1	B	71	LEU
1	B	80	GLU
1	B	105	ILE
1	B	135	LEU
1	B	152	SER
1	B	162	THR
1	B	168	LEU
1	B	170	SER
1	B	188	LEU
1	B	245	LEU
1	B	269	LEU
1	C	4	LEU
1	C	5	LEU
1	C	6	ASP
1	C	46	LEU
1	C	60	LEU
1	C	67	ASN
1	C	69	GLU
1	C	70	HIS
1	C	71	LEU
1	C	74	LEU
1	C	101	THR
1	C	135	LEU
1	C	145	VAL
1	C	162	THR
1	C	168	LEU
1	C	170	SER
1	C	207	LEU
1	C	210	GLU
1	C	218	LEU
1	C	245	LEU
1	C	253	THR
1	D	5	LEU
1	D	44	LEU
1	D	67	ASN

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Mol	Chain	Res	Type
1	D	70	HIS
1	D	71	LEU
1	D	74	LEU
1	D	135	LEU
1	D	145	VAL
1	D	162	THR
1	D	168	LEU
1	D	188	LEU
1	D	253	THR
1	E	2	THR
1	E	5	LEU
1	E	18	ASP
1	E	44	LEU
1	E	57	LYS
1	E	60	LEU
1	E	67	ASN
1	E	69	GLU
1	E	71	LEU
1	E	74	LEU
1	E	101	THR
1	E	145	VAL
1	E	162	THR
1	E	168	LEU
1	E	188	LEU
1	E	214	GLN
1	E	216	GLN
1	E	219	GLU
1	E	222	TRP
1	E	233	LYS
1	E	253	THR
1	F	4	LEU
1	F	6	ASP
1	F	9	ARG
1	F	39	THR
1	F	67	ASN
1	F	71	LEU
1	F	74	LEU
1	F	145	VAL
1	F	170	SER
1	F	188	LEU
1	F	195	ARG
1	F	197	LEU

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Mol	Chain	Res	Type
1	F	199	MET
1	F	245	LEU
1	F	248	ASP
1	F	256	ASP
1	F	269	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	24	HIS
1	A	35	GLN
1	A	67	ASN
1	A	139	ASN
1	B	35	GLN
1	B	66	GLN
1	B	67	ASN
1	B	224	GLN
1	C	24	HIS
1	C	67	ASN
1	C	70	HIS
1	D	35	GLN
1	D	67	ASN
1	D	70	HIS
1	D	139	ASN
1	E	67	ASN
1	E	139	ASN
1	E	214	GLN
1	E	216	GLN
1	F	24	HIS
1	F	67	ASN
1	F	70	HIS
1	F	106	ASN
1	F	139	ASN
1	F	231	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

8 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	NAD	A	401	-	38,48,48	1.49	3 (7%)	47,73,73	2.34	7 (14%)
3	NAD	B	402	-	38,48,48	1.62	3 (7%)	47,73,73	2.09	7 (14%)
2	8PS	C	300	-	23,23,23	1.31	1 (4%)	28,28,28	1.07	4 (14%)
3	NAD	C	403	-	38,48,48	1.88	4 (10%)	47,73,73	2.48	8 (17%)
2	8PS	D	301	-	23,23,23	1.33	1 (4%)	28,28,28	1.24	3 (10%)
3	NAD	D	404	-	38,48,48	1.69	3 (7%)	47,73,73	2.07	9 (19%)
3	NAD	E	405	-	38,48,48	1.71	4 (10%)	47,73,73	2.55	11 (23%)
3	NAD	F	406	-	38,48,48	1.61	3 (7%)	47,73,73	2.63	7 (14%)

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	NAD	A	401	-	-	0/22/62/62	0/5/5/5
3	NAD	B	402	-	-	0/22/62/62	0/5/5/5
2	8PS	C	300	-	-	0/12/12/12	0/2/2/2
3	NAD	C	403	-	-	0/22/62/62	0/5/5/5
2	8PS	D	301	-	-	0/12/12/12	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAD	D	404	-	-	0/22/62/62	0/5/5/5
3	NAD	E	405	-	-	0/22/62/62	0/5/5/5
3	NAD	F	406	-	-	0/22/62/62	0/5/5/5

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	405	NAD	PN-O2N	-2.25	1.45	1.54
3	D	404	NAD	O4D-C1D	2.04	1.43	1.41
3	C	403	NAD	C2A-N1A	2.05	1.37	1.33
3	E	405	NAD	C2A-N1A	2.35	1.38	1.33
3	F	406	NAD	C2A-N1A	2.39	1.38	1.33
3	A	401	NAD	C2A-N1A	2.55	1.38	1.33
3	B	402	NAD	C2A-N1A	2.57	1.38	1.33
3	A	401	NAD	C2A-N3A	2.64	1.36	1.32
3	E	405	NAD	C2A-N3A	3.11	1.37	1.32
3	C	403	NAD	O4D-C1D	3.20	1.45	1.41
3	F	406	NAD	C2A-N3A	3.44	1.38	1.32
3	B	402	NAD	C2A-N3A	3.90	1.39	1.32
3	D	404	NAD	C2A-N3A	4.22	1.39	1.32
2	D	301	8PS	C6-C5	4.57	1.47	1.40
2	C	300	8PS	C6-C5	5.12	1.48	1.40
3	C	403	NAD	C2A-N3A	5.57	1.42	1.32
3	A	401	NAD	O7N-C7N	7.20	1.39	1.24
3	B	402	NAD	O7N-C7N	7.57	1.40	1.24
3	F	406	NAD	O7N-C7N	7.88	1.41	1.24
3	C	403	NAD	O7N-C7N	7.91	1.41	1.24
3	D	404	NAD	O7N-C7N	7.96	1.41	1.24
3	E	405	NAD	O7N-C7N	8.07	1.41	1.24

All (56) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	406	NAD	N3A-C2A-N1A	-12.92	119.00	128.89
3	A	401	NAD	N3A-C2A-N1A	-12.38	119.42	128.89
3	C	403	NAD	N3A-C2A-N1A	-12.36	119.43	128.89
3	E	405	NAD	N3A-C2A-N1A	-12.33	119.45	128.89
3	D	404	NAD	N3A-C2A-N1A	-10.61	120.77	128.89
3	B	402	NAD	N3A-C2A-N1A	-10.58	120.79	128.89
3	F	406	NAD	C2B-C1B-N9A	-5.06	106.56	114.29
3	A	401	NAD	C4B-O4B-C1B	-5.02	104.20	109.72
3	C	403	NAD	C4B-O4B-C1B	-4.25	105.05	109.72

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	403	NAD	PN-O3-PA	-4.10	121.22	132.73
3	B	402	NAD	C4B-O4B-C1B	-4.09	105.22	109.72
3	C	403	NAD	C2B-C1B-N9A	-3.76	108.55	114.29
3	D	404	NAD	PN-O3-PA	-3.76	122.18	132.73
3	E	405	NAD	O7N-C7N-C3N	-3.61	115.64	119.59
3	A	401	NAD	C1B-N9A-C4A	-3.61	121.50	126.94
3	A	401	NAD	PN-O3-PA	-3.53	122.81	132.73
3	F	406	NAD	PN-O3-PA	-3.53	122.83	132.73
3	E	405	NAD	C4B-O4B-C1B	-3.49	105.89	109.72
3	B	402	NAD	PN-O3-PA	-3.45	123.03	132.73
3	F	406	NAD	C4B-O4B-C1B	-3.35	106.03	109.72
3	D	404	NAD	C4A-C5A-N7A	-2.95	106.77	109.48
3	E	405	NAD	C2B-C1B-N9A	-2.77	110.05	114.29
3	E	405	NAD	C3N-C2N-N1N	-2.75	117.20	120.36
3	F	406	NAD	C1B-N9A-C4A	-2.70	122.86	126.94
3	D	404	NAD	C4B-O4B-C1B	-2.51	106.96	109.72
3	A	401	NAD	C2B-C1B-N9A	-2.45	110.55	114.29
3	D	404	NAD	O7N-C7N-C3N	-2.36	117.01	119.59
3	E	405	NAD	C4D-O4D-C1D	-2.23	107.27	109.72
2	D	301	8PS	C15-C14-C2	-2.23	105.23	113.71
2	C	300	8PS	C1-C6-C5	-2.20	117.35	119.79
3	B	402	NAD	C4A-C5A-N7A	-2.18	107.47	109.48
2	C	300	8PS	C15-C14-C2	-2.11	105.66	113.71
3	A	401	NAD	C4A-C5A-N7A	-2.04	107.60	109.48
3	C	403	NAD	C4A-C5A-N7A	-2.02	107.62	109.48
3	E	405	NAD	PN-O3-PA	-2.00	127.10	132.73
2	C	300	8PS	C5-O7-C8	2.02	123.05	117.77
2	D	301	8PS	C3-C2-C1	2.26	121.89	118.55
2	C	300	8PS	C3-C2-C1	2.27	121.91	118.55
3	E	405	NAD	C2N-C3N-C4N	2.46	121.03	118.29
3	D	404	NAD	O4B-C1B-N9A	2.50	113.34	108.10
3	D	404	NAD	C2N-C3N-C4N	2.61	121.20	118.29
3	B	402	NAD	C3N-C7N-N7N	2.70	120.78	117.82
3	D	404	NAD	C3N-C7N-N7N	2.75	120.83	117.82
3	D	404	NAD	O4D-C1D-N1N	2.82	111.23	108.13
3	B	402	NAD	O4D-C1D-N1N	2.86	111.28	108.13
3	C	403	NAD	O4B-C1B-N9A	3.32	115.04	108.10
3	E	405	NAD	C3N-C7N-N7N	3.36	121.49	117.82
3	E	405	NAD	O4B-C1B-N9A	3.37	115.14	108.10
2	D	301	8PS	C5-O7-C8	3.63	127.25	117.77
3	A	401	NAD	O4B-C1B-N9A	3.67	115.78	108.10
3	C	403	NAD	C3N-C7N-N7N	3.86	122.04	117.82

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	402	NAD	O4B-C1B-N9A	4.28	117.06	108.10
3	C	403	NAD	O4D-C1D-N1N	5.18	113.83	108.13
3	F	406	NAD	O4B-C1B-N9A	5.83	120.29	108.10
3	F	406	NAD	O4D-C1D-N1N	6.25	115.00	108.13
3	E	405	NAD	O4D-C1D-N1N	6.57	115.34	108.13

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

8 monomers are involved in 25 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	401	NAD	1	0
3	B	402	NAD	1	0
2	C	300	8PS	3	0
3	C	403	NAD	2	0
2	D	301	8PS	7	0
3	D	404	NAD	5	0
3	E	405	NAD	4	0
3	F	406	NAD	2	0

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 ⓘ

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	248/269 (92%)	0.20	12 (4%) 34 27	28, 44, 66, 77	0
1	B	245/269 (91%)	0.53	20 (8%) 14 10	32, 51, 78, 91	0
1	C	260/269 (96%)	-0.21	5 (1%) 70 64	9, 25, 54, 66	0
1	D	260/269 (96%)	-0.15	4 (1%) 76 71	9, 24, 58, 70	0
1	E	260/269 (96%)	0.23	19 (7%) 18 12	25, 39, 63, 73	0
1	F	248/269 (92%)	0.37	19 (7%) 16 11	25, 44, 64, 76	0
All	All	1521/1614 (94%)	0.16	79 (5%) 31 24	9, 39, 65, 91	0

All (79) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	49	ARG	4.9
1	A	218	LEU	4.3
1	B	80	GLU	4.3
1	A	46	LEU	4.1
1	B	56	ALA	3.9
1	E	2	THR	3.8
1	E	83	GLY	3.8
1	E	208	GLY	3.7
1	F	85	GLY	3.7
1	C	219	GLU	3.6
1	A	84	ALA	3.5
1	E	213	ALA	3.5
1	F	57	LYS	3.5
1	F	86	ASN	3.4
1	B	195	ARG	3.4
1	F	2	THR	3.3
1	F	105	ILE	3.3
1	F	104	GLY	3.3
1	E	217	LEU	3.2

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Mol	Chain	Res	Type	RSRZ
1	B	84	ALA	3.2
1	B	57	LYS	3.0
1	B	137	ILE	3.0
1	C	2	THR	3.0
1	B	76	GLY	3.0
1	A	2	THR	2.9
1	C	215	ILE	2.9
1	B	105	ILE	2.9
1	D	45	ARG	2.9
1	B	53	ARG	2.9
1	B	99	PRO	2.9
1	A	195	ARG	2.8
1	E	212	GLY	2.8
1	F	140	PRO	2.7
1	D	49	ARG	2.7
1	E	220	GLU	2.7
1	E	269	LEU	2.6
1	E	41	PHE	2.6
1	E	74	LEU	2.6
1	B	81	ALA	2.6
1	E	84	ALA	2.6
1	B	103	MET	2.6
1	F	197	LEU	2.5
1	B	220	GLU	2.5
1	F	83	GLY	2.5
1	F	56	ALA	2.5
1	F	158	TYR	2.5
1	B	83	GLY	2.4
1	A	53	ARG	2.4
1	D	43	ARG	2.4
1	A	45	ARG	2.4
1	F	53	ARG	2.4
1	F	55	PRO	2.3
1	E	214	GLN	2.3
1	D	214	GLN	2.3
1	B	18	ASP	2.3
1	A	72	ALA	2.3
1	A	57	LYS	2.3
1	B	248	ASP	2.3
1	F	106	ASN	2.3
1	E	86	ASN	2.2
1	E	215	ILE	2.2

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Mol	Chain	Res	Type	RSRZ
1	E	72	ALA	2.2
1	E	85	GLY	2.2
1	E	211	ALA	2.2
1	B	6	ASP	2.2
1	F	257	ILE	2.1
1	F	45	ARG	2.1
1	A	67	ASN	2.1
1	E	35	GLN	2.1
1	F	199	MET	2.1
1	F	100	GLN	2.1
1	C	220	GLU	2.1
1	E	210	GLU	2.1
1	A	85	GLY	2.1
1	B	5	LEU	2.0
1	C	49	ARG	2.0
1	B	34	ALA	2.0
1	F	84	ALA	2.0
1	A	100	GLN	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	8PS	D	301	22/22	0.89	0.21	1.16	36,39,43,44	0
3	NAD	B	402	44/44	0.90	0.19	0.04	45,53,55,56	0
2	8PS	C	300	22/22	0.93	0.16	-0.40	34,35,37,38	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	NAD	C	403	44/44	0.96	0.15	-0.41	22,27,33,34	0
3	NAD	A	401	44/44	0.93	0.15	-0.61	38,42,44,44	0
3	NAD	F	406	44/44	0.93	0.15	-0.66	24,46,55,55	0
3	NAD	D	404	44/44	0.96	0.13	-0.69	15,26,31,35	0
3	NAD	E	405	44/44	0.96	0.12	-1.12	29,36,40,41	0

6.5 Other polymers [i](#)

There are no such residues in this entry.