



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:49 PM GMT

PDB ID : 4PLY
Title : Crystal structure of ancestral apicomplexan malate dehydrogenase with malate.
Authors : Boucher, J.I.; Jacobowitz, J.R.; Beckett, B.C.; Classen, S.; Theobald, D.L.
Deposited on : 2014-05-19
Resolution : 1.90 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
<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

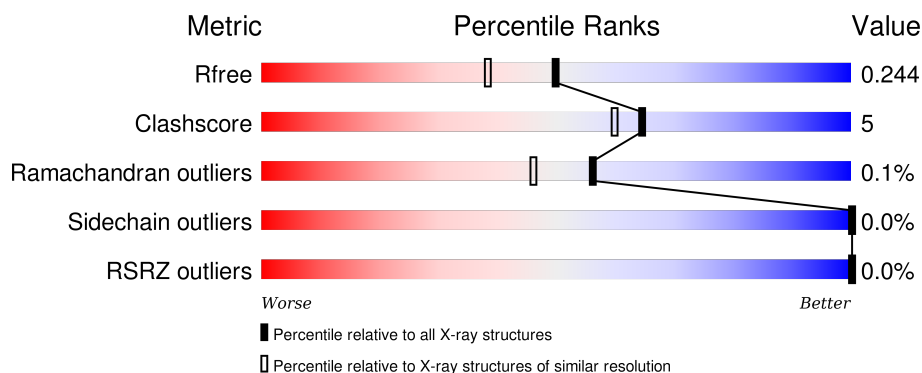
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION




The reported resolution of this entry is 1.90 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}	91344	4755 (1.90-1.90)
Clashscore	102246	5398 (1.90-1.90)
Ramachandran outliers	100387	5338 (1.90-1.90)
Sidechain outliers	100360	5339 (1.90-1.90)
RSRZ outliers	91569	4766 (1.90-1.90)

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 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	331	
1	B	331	
1	C	331	
1	D	331	
1	E	331	

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Mol	Chain	Length	Quality of chain
1	F	331	
1	G	331	
1	H	331	

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
3	PYR	A	402	-	-	-	X
3	PYR	B	402	-	-	-	X
3	PYR	D	402	-	-	-	X
3	PYR	E	402	-	-	-	X
3	PYR	F	402	-	-	-	X
3	PYR	H	402	-	-	-	X

2 Entry composition

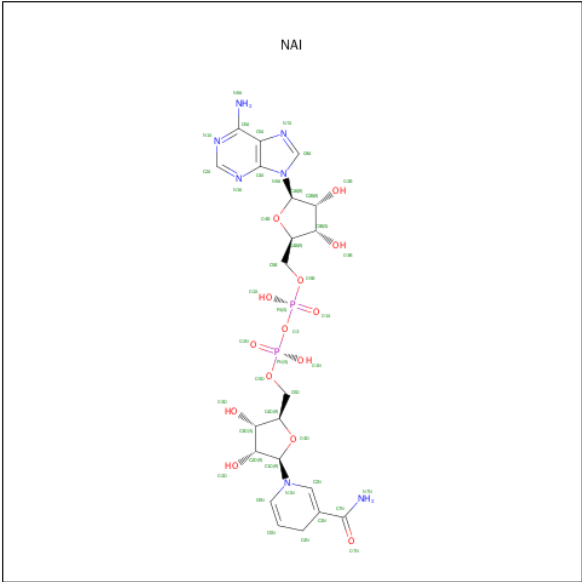
There are 4 unique types of molecules in this entry. The entry contains 19938 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 malate dehydrogenase.

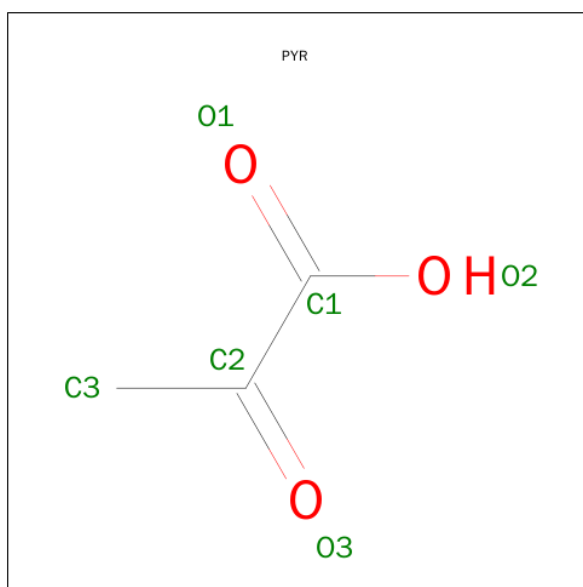
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	318	Total	C	N	O	S	0	2	0
			2434	1550	407	462	15			
1	A	318	Total	C	N	O	S	0	3	0
			2437	1550	406	466	15			
1	C	303	Total	C	N	O	S	0	2	0
			2268	1444	377	432	15			
1	D	303	Total	C	N	O	S	0	2	0
			2294	1463	383	433	15			
1	F	318	Total	C	N	O	S	0	2	0
			2434	1549	407	463	15			
1	E	318	Total	C	N	O	S	0	2	0
			2422	1541	405	461	15			
1	G	303	Total	C	N	O	S	0	2	0
			2264	1442	378	429	15			
1	H	303	Total	C	N	O	S	0	2	0
			2287	1458	381	433	15			

- Molecule 2 is 1,4-DIHYDRONICOTINAMIDE ADENINE DINUCLEOTIDE (three-letter code: NAI) (formula: C₂₁H₂₉N₇O₁₄P₂).



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

- Molecule 3 is PYRUVIC ACID (three-letter code: PYR) (formula: C₃H₄O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	D	1	Total	C	O	0	0
			6	3	3		
3	F	1	Total	C	O	0	0
			6	3	3		
3	E	1	Total	C	O	0	0
			6	3	3		
3	H	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is water.

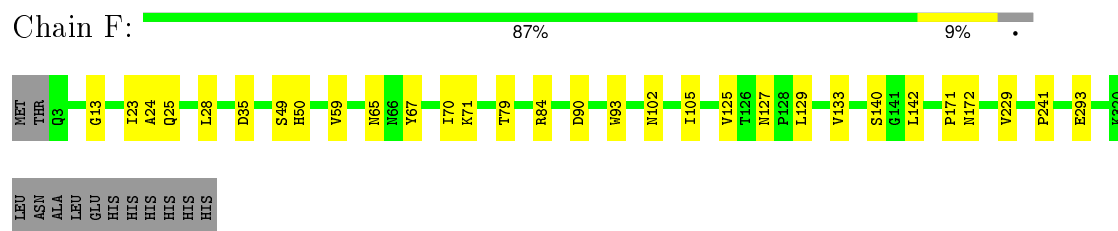
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	111	Total	O	0	0
			111	111		
4	A	80	Total	O	0	0
			80	80		
4	C	96	Total	O	0	0
			96	96		
4	D	78	Total	O	0	0
			78	78		
4	F	117	Total	O	0	0
			117	117		
4	E	85	Total	O	0	0
			85	85		

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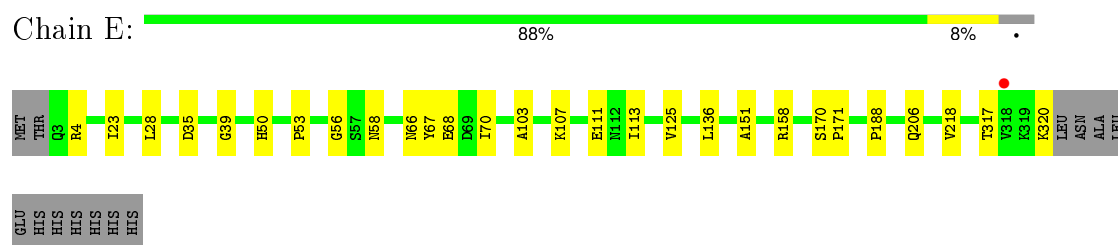
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	G	77	Total 77	O 77	0	0
4	H	66	Total 66	O 66	0	0

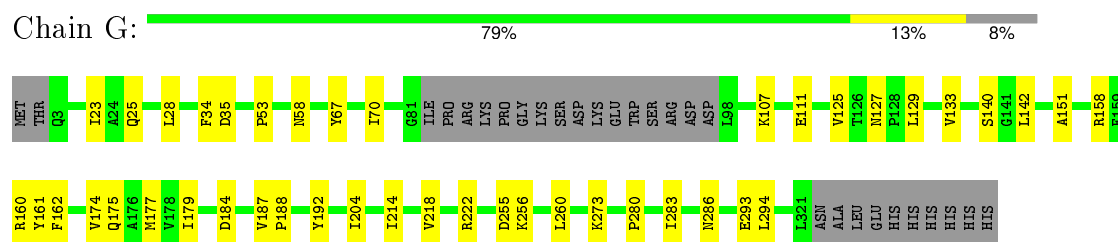
- Molecule 1: malate dehydrogenase



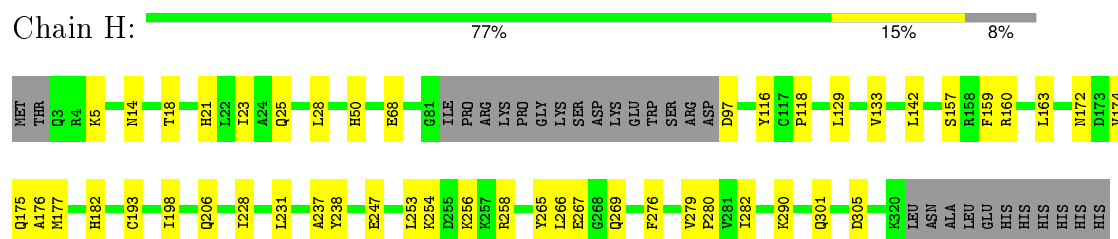
- Molecule 1: malate dehydrogenase



- Molecule 1: malate dehydrogenase



- Molecule 1: malate dehydrogenase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	59.31Å 90.69Å 136.30Å 90.01° 90.01° 71.44°	Depositor
Resolution (Å)	45.44 – 1.90 45.43 – 1.90	Depositor EDS
% Data completeness (in resolution range)	90.2 (45.44-1.90) 76.9 (45.43-1.90)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.20 (at 1.89Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, R_{free}	0.189 , 0.238 0.191 , 0.244	Depositor DCC
R_{free} test set	1515 reflections (0.90%)	DCC
Wilson B-factor (Å ²)	23.0	Xtriage
Anisotropy	0.338	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 41.1	EDS
Estimated twinning fraction	0.500 for -h,-k,l 0.035 for h,h-k,-l 0.477 for -h,-k,l 0.034 for -h,-h+k,-l	Xtriage
Reported twinning fraction	0.500 for -h,-k,l	Depositor
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 197555 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	19938	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.05% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: PYR, NAI

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	0.22	0/2485	0.39	0/3361
1	B	0.21	0/2479	0.39	0/3351
1	C	0.22	0/2308	0.39	0/3128
1	D	0.22	0/2334	0.39	0/3156
1	E	0.21	0/2467	0.39	0/3339
1	F	0.21	0/2479	0.39	0/3352
1	G	0.22	0/2304	0.40	0/3123
1	H	0.21	0/2327	0.39	0/3149
All	All	0.22	0/19183	0.39	0/25959

There are no bond length outliers.

There are no bond angle outliers.

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	2437	0	2483	28	0
1	B	2434	0	2490	23	0
1	C	2268	0	2269	33	0
1	D	2294	0	2337	27	0
1	E	2422	0	2459	21	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	F	2434	0	2485	25	0
1	G	2264	0	2267	32	0
1	H	2287	0	2317	32	0
2	A	44	0	27	2	0
2	B	44	0	27	3	0
2	C	44	0	27	1	0
2	D	44	0	27	4	0
2	E	44	0	27	3	0
2	F	44	0	27	5	0
2	G	44	0	27	3	0
2	H	44	0	27	1	0
3	A	6	0	3	0	0
3	B	6	0	3	3	0
3	D	6	0	3	1	0
3	E	6	0	3	1	0
3	F	6	0	3	3	0
3	H	6	0	3	1	0
4	A	80	0	0	2	0
4	B	111	0	0	0	0
4	C	96	0	0	3	0
4	D	78	0	0	2	0
4	E	85	0	0	2	0
4	F	117	0	0	2	0
4	G	77	0	0	5	0
4	H	66	0	0	1	0
All	All	19938	0	19341	197	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 5.

All (197) 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:D:129:LEU:HD23	1:D:151:ALA:HB2	1.68	0.76
1:G:273:LYS:O	4:G:543:HOH:O	2.07	0.72
1:C:129:LEU:HD23	1:C:151:ALA:HB2	1.73	0.70
1:F:293:GLU:O	1:E:206:GLN:NE2	2.25	0.69
1:B:5:LYS:HD3	1:B:253:LEU:HD22	1.74	0.69
1:A:23:ILE:HG23	1:A:28:LEU:HB2	1.74	0.68
1:G:23:ILE:HG23	1:G:28:LEU:HB2	1.76	0.67
1:H:5:LYS:HD3	1:H:253:LEU:HD22	1.78	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:140:SER:HB2	1:F:142:LEU:HD13	1.79	0.64
1:F:23:ILE:HG23	1:F:28:LEU:HB2	1.79	0.64
1:D:23:ILE:HG23	1:D:28:LEU:HB2	1.79	0.63
1:F:71:LYS:NZ	4:F:501:HOH:O	2.31	0.63
1:A:129:LEU:HD22	1:A:151:ALA:HB2	1.79	0.62
1:H:97:ASP:N	4:H:538:HOH:O	2.32	0.62
1:C:102:ASN:ND2	1:C:127:ASN:O	2.25	0.62
1:A:266:LEU:HD21	1:A:279:VAL:HG21	1.81	0.62
1:H:23:ILE:HG23	1:H:28:LEU:HB2	1.81	0.60
1:H:301:GLN:NE2	1:H:305:ASP:OD1	2.35	0.60
1:G:107:LYS:NZ	1:G:111:GLU:OE2	2.34	0.58
1:E:107:LYS:NZ	1:E:111:GLU:OE2	2.34	0.58
1:A:182:HIS:NE2	2:A:401:NAI:O7N	2.36	0.58
1:E:23:ILE:HG23	1:E:28:LEU:HB2	1.85	0.58
1:E:113:ILE:HD12	1:E:136:LEU:HD21	1.83	0.58
1:G:286:ASN:ND2	4:G:502:HOH:O	2.36	0.58
1:E:317:THR:HA	1:E:320:LYS:HE2	1.86	0.57
1:G:175:GLN:NE2	1:H:175:GLN:HG2	2.19	0.57
1:G:127:ASN:ND2	4:G:542:HOH:O	2.37	0.56
1:H:21:HIS:CD2	1:H:25:GLN:HE21	2.22	0.56
1:D:182:HIS:NE2	2:D:401:NAI:O7N	2.39	0.56
1:E:50:HIS:O	1:H:157:SER:HB3	2.06	0.56
1:G:188:PRO:HG2	1:G:218:VAL:HG13	1.88	0.56
1:C:184:ASP:O	1:C:222:ARG:NH1	2.36	0.55
1:A:247:GLU:OE1	1:A:258:ARG:NH1	2.34	0.55
1:E:53:PRO:O	1:H:172:ASN:ND2	2.37	0.55
1:G:175:GLN:HE22	1:H:175:GLN:HG2	1.72	0.55
1:G:294:LEU:HD21	1:H:198:ILE:HD13	1.89	0.55
1:E:70:ILE:HG13	4:E:533:HOH:O	2.07	0.54
1:B:23:ILE:HG23	1:B:28:LEU:HB2	1.89	0.54
1:C:129:LEU:CD2	1:C:151:ALA:HB2	2.38	0.53
1:H:266:LEU:HD21	1:H:279:VAL:HG21	1.90	0.53
1:E:67:TYR:O	1:E:70:ILE:HG12	2.08	0.53
1:B:35:ASP:OD1	1:B:36:ILE:N	2.43	0.52
1:C:182:HIS:NE2	4:C:590:HOH:O	2.34	0.52
1:F:50:HIS:CG	1:G:158:ARG:HG2	2.44	0.52
1:G:140:SER:HB2	1:G:142:LEU:HD13	1.91	0.52
1:A:256:LYS:HB3	1:C:58:ASN:HD22	1.75	0.52
1:G:125:VAL:O	2:G:401:NAI:H2N	2.09	0.52
1:D:127:ASN:ND2	3:D:402:PYP:O3	2.43	0.52
1:E:56:GLY:HA2	1:H:172:ASN:HB2	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:254:LYS:HD2	1:H:256:LYS:HE3	1.92	0.51
1:A:108:ASP:O	1:A:112:ASN:ND2	2.40	0.51
1:G:35:ASP:OD2	2:G:401:NAI:O2B	2.29	0.51
1:C:266:LEU:HD21	1:C:279:VAL:HG21	1.92	0.51
1:D:8:SER:OG	1:D:73:SER:OG	2.23	0.51
1:F:25:GLN:OE1	1:G:25:GLN:NE2	2.44	0.51
1:C:110:ALA:HB2	1:C:136:LEU:HD22	1.93	0.51
1:G:23:ILE:HG12	1:G:28:LEU:HD12	1.92	0.50
1:A:27:GLU:HG2	1:A:57:SER:HB2	1.93	0.50
1:F:35:ASP:O	1:F:65:ASN:ND2	2.43	0.50
1:C:116:TYR:OH	4:C:501:HOH:O	2.17	0.50
1:A:14:ASN:ND2	4:A:540:HOH:O	2.41	0.50
1:B:99:LEU:HD21	1:B:318:VAL:HG22	1.94	0.50
1:G:204:ILE:HD11	1:G:214:ILE:HD12	1.95	0.49
1:G:160:ARG:NE	1:G:174:VAL:O	2.44	0.49
1:G:260:LEU:HD12	1:G:283:ILE:HD12	1.94	0.49
1:C:294:LEU:HD21	1:D:198:ILE:HD13	1.94	0.48
1:E:4:ARG:HE	1:G:255:ASP:HB3	1.78	0.48
1:E:103:ALA:HB2	4:E:574:HOH:O	2.14	0.48
1:G:184:ASP:O	1:G:222:ARG:NH1	2.46	0.48
1:G:67:TYR:O	1:G:70:ILE:HG12	2.13	0.48
1:D:152:GLY:HA3	1:D:261:PRO:HB2	1.96	0.48
1:D:129:LEU:HD22	2:D:401:NAI:H72N	1.78	0.47
1:B:6:LYS:HE3	1:B:32:VAL:HG21	1.96	0.47
1:B:113:ILE:HD12	1:B:136:LEU:HD21	1.96	0.47
1:E:66:ASN:HD21	1:E:68:GLU:HB2	1.79	0.47
1:D:124:VAL:HB	1:D:133:VAL:HG22	1.95	0.47
1:D:67:TYR:O	1:D:70:ILE:HG12	2.14	0.47
1:F:125:VAL:O	2:F:401:NAI:H2N	2.15	0.47
1:F:129:LEU:O	1:F:133:VAL:HG23	2.15	0.47
1:H:182:HIS:NE2	3:H:402:PYR:O3	2.43	0.47
1:D:270:TYR:OH	1:D:292:ILE:O	2.27	0.47
1:A:140:SER:HB2	1:A:142:LEU:HD13	1.96	0.47
1:C:14:ASN:HB3	1:C:238:TYR:CG	2.50	0.46
1:B:108:ASP:O	1:B:112:ASN:ND2	2.47	0.46
1:H:159:PHE:CD2	1:H:176:ALA:HB1	2.51	0.46
1:A:35:ASP:OD1	2:A:401:NAI:O2B	2.28	0.46
1:C:114:LYS:HB2	1:C:140:SER:HA	1.98	0.46
1:A:186:MET:O	1:A:222:ARG:NH1	2.45	0.46
1:C:151:ALA:HB3	4:C:569:HOH:O	2.16	0.46
1:C:309:GLU:O	1:C:313:GLU:HG3	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:125:VAL:O	2:B:401:NAI:H2N	2.16	0.45
1:A:68:GLU:HG2	1:A:116:TYR:CZ	2.52	0.45
1:H:265:TYR:HB2	1:H:276:PHE:CE2	2.52	0.45
1:B:290:LYS:HE2	1:A:208:TRP:HZ2	1.81	0.45
1:E:58:ASN:ND2	1:G:256:LYS:O	2.45	0.45
1:E:158:ARG:HG2	1:H:50:HIS:CG	2.51	0.45
1:G:293:GLU:O	1:H:206:GLN:NE2	2.49	0.45
1:A:5:LYS:HD3	1:A:253:LEU:HD22	1.97	0.45
1:B:50:HIS:CG	1:C:158:ARG:HG2	2.51	0.45
1:A:157[B]:SER:OG	1:D:50:HIS:O	2.32	0.45
1:F:127:ASN:HD21	3:F:402:PYR:H31	1.82	0.45
1:E:39:GLY:O	1:H:231:LEU:HB3	2.16	0.45
1:F:84:ARG:NH1	4:F:568:HOH:O	2.50	0.45
1:D:107:LYS:NZ	1:D:111:GLU:OE2	2.42	0.45
1:D:131:VAL:HG21	1:D:314:LEU:HB3	1.98	0.45
1:A:266:LEU:HD11	1:A:279:VAL:HG11	1.99	0.44
1:A:14:ASN:ND2	4:A:519:HOH:O	2.50	0.44
1:A:90:ASP:OD2	1:A:233:LYS:HA	2.17	0.44
1:B:35:ASP:OD1	2:B:401:NAI:O2B	2.32	0.44
1:G:177:MET:SD	1:G:280:PRO:HB3	2.57	0.44
1:C:107:LYS:HE2	1:C:139:TYR:CE2	2.52	0.44
1:C:8:SER:OG	1:C:73:SER:OG	2.20	0.44
1:A:88:LYS:NZ	1:A:92:GLU:OE1	2.50	0.44
1:C:35:ASP:OD2	1:C:36:ILE:N	2.50	0.44
1:A:129:LEU:O	1:A:133:VAL:HG23	2.17	0.44
1:H:256:LYS:HB3	1:H:258:ARG:HG2	1.99	0.44
1:D:110:ALA:HB1	1:D:140:SER:HB3	1.99	0.44
1:C:247:GLU:OE1	1:C:258:ARG:NH1	2.46	0.44
1:A:166:LYS:HE3	1:A:166:LYS:HB2	1.88	0.44
1:F:241:PRO:HB3	2:F:401:NAI:H4N	2.00	0.44
1:B:312:ARG:O	1:B:316:GLU:HG2	2.18	0.44
1:H:228:ILE:HG21	1:H:237:ALA:HB2	2.00	0.44
1:F:84:ARG:HG2	1:F:93:TRP:CD1	2.53	0.44
1:C:265:TYR:HB2	1:C:276:PHE:CE2	2.53	0.44
1:C:145:ASN:ND2	1:C:286:ASN:O	2.50	0.43
1:H:177:MET:SD	1:H:280:PRO:HB3	2.58	0.43
1:A:58:ASN:HD22	1:C:256:LYS:HA	1.83	0.43
2:H:401:NAI:H6N	2:H:401:NAI:H2D	1.83	0.43
1:B:158:ARG:HG2	1:C:50:HIS:CG	2.53	0.43
1:F:171:PRO:HB3	1:G:53:PRO:HG3	2.00	0.43
1:E:125:VAL:O	2:E:401:NAI:H2N	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:21:HIS:CD2	1:A:25:GLN:HE21	2.37	0.43
1:H:14:ASN:O	1:H:18:THR:OG1	2.34	0.43
1:C:124:VAL:HB	1:C:133:VAL:HG22	2.00	0.43
1:F:93:TRP:CD2	1:F:229:VAL:HG21	2.53	0.43
1:G:192:TYR:O	4:G:528:HOH:O	2.21	0.43
1:G:179:ILE:HG13	1:G:187:VAL:HB	1.99	0.43
1:D:286:ASN:HB2	1:D:289:GLU:HG2	1.99	0.43
1:D:192:TYR:CE1	1:D:294:LEU:HD22	2.53	0.43
1:F:102:ASN:ND2	1:F:127:ASN:O	2.25	0.43
1:D:35:ASP:OD1	2:D:401:NAI:O2B	2.36	0.42
1:B:95:ARG:HH21	3:B:402:PYR:C1	2.32	0.42
1:C:137:HIS:CE1	1:C:144:HIS:HB3	2.54	0.42
1:H:160:ARG:NE	1:H:174:VAL:O	2.36	0.42
1:H:68:GLU:HG2	1:H:116:TYR:CE1	2.54	0.42
1:F:67:TYR:O	1:F:70:ILE:HG12	2.18	0.42
1:A:204:ILE:HA	1:A:209:ILE:O	2.19	0.42
1:E:188:PRO:HG2	1:E:218:VAL:HG13	2.00	0.42
1:H:247:GLU:OE1	1:H:258:ARG:NH2	2.38	0.42
1:B:265:TYR:HB2	1:B:276:PHE:CE2	2.54	0.42
1:C:125:VAL:O	2:C:401:NAI:H2N	2.18	0.42
1:B:50:HIS:CE1	1:C:228:ILE:HD11	2.54	0.42
1:D:129:LEU:CD2	1:D:151:ALA:HB2	2.45	0.42
1:C:129:LEU:O	1:C:133:VAL:HG23	2.19	0.42
1:B:160:ARG:HH21	1:B:172:ASN:HA	1.84	0.42
1:A:256:LYS:HB3	1:C:58:ASN:ND2	2.34	0.42
1:F:79:THR:HG22	1:F:125:VAL:HB	2.02	0.42
1:F:13:GLY:HA3	2:F:401:NAI:O5B	2.20	0.42
1:H:129:LEU:O	1:H:133:VAL:HG23	2.20	0.42
1:B:129:LEU:O	1:B:133:VAL:HG23	2.20	0.42
1:D:220:ARG:NH1	4:D:553:HOH:O	2.36	0.42
1:B:231:LEU:HB3	1:C:39:GLY:O	2.20	0.41
1:D:301:GLN:NE2	1:D:305:ASP:OD1	2.49	0.41
2:F:401:NAI:H42N	3:F:402:PYR:C2	2.50	0.41
2:E:401:NAI:H42N	3:E:402:PYR:C2	2.51	0.41
1:D:265:TYR:HB2	1:D:276:PHE:CZ	2.55	0.41
1:G:162:PHE:O	4:G:549:HOH:O	2.21	0.41
1:E:66:ASN:ND2	1:E:68:GLU:HB2	2.35	0.41
1:F:127:ASN:ND2	3:F:402:PYR:H31	2.35	0.41
1:F:84:ARG:NH1	1:F:90:ASP:OD1	2.52	0.41
1:B:67:TYR:O	1:B:70:ILE:HG12	2.20	0.41
2:B:401:NAI:H42N	3:B:402:PYR:C2	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:14:ASN:HB3	1:H:238:TYR:CG	2.55	0.41
1:A:173:ASP:O	1:A:196:GLY:N	2.46	0.41
1:G:129:LEU:O	1:G:133:VAL:HG23	2.20	0.41
1:F:105:ILE:HG23	2:F:401:NAI:N6A	2.35	0.41
1:H:267:GLU:N	1:H:269:GLN:OE1	2.48	0.41
1:D:266:LEU:HD21	1:D:279:VAL:HG21	2.02	0.41
1:G:34:PHE:CE2	2:G:401:NAI:H2A	2.56	0.41
1:H:118:PRO:HA	1:H:142:LEU:HD11	2.03	0.41
1:B:266:LEU:HD21	1:B:279:VAL:HG21	2.02	0.41
1:E:35:ASP:OD1	2:E:401:NAI:O2B	2.31	0.41
1:F:24:ALA:HA	1:F:59:VAL:HG21	2.03	0.41
1:D:266:LEU:HD12	1:D:272:VAL:HB	2.03	0.41
1:F:172:ASN:ND2	1:G:58:ASN:HD21	2.18	0.41
1:H:163:LEU:HD13	1:H:193[A]:CYS:SG	2.61	0.41
1:D:151:ALA:HB3	4:D:531:HOH:O	2.21	0.40
1:D:125:VAL:O	2:D:401:NAI:H2N	2.21	0.40
1:A:282:ILE:HD12	1:A:290:LYS:HB3	2.02	0.40
1:B:39:GLY:O	1:C:231:LEU:HB3	2.22	0.40
1:B:95:ARG:NH2	3:B:402:PYR:O2	2.51	0.40
1:D:8:SER:HB2	1:D:76:VAL:HG22	2.04	0.40
1:E:170:SER:HA	1:E:171:PRO:HD3	1.89	0.40
1:F:49:SER:HB3	1:G:161:TYR:CZ	2.57	0.40
1:C:265:TYR:OH	1:C:274:ASP:HA	2.22	0.40
1:H:282:ILE:HD12	1:H:290:LYS:HD2	2.03	0.40
1:C:118:PRO:HA	1:C:142:LEU:HD11	2.04	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	319/331 (96%)	305 (96%)	14 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	318/331 (96%)	307 (96%)	11 (4%)	0	100	100
1	C	301/331 (91%)	293 (97%)	8 (3%)	0	100	100
1	D	301/331 (91%)	290 (96%)	11 (4%)	0	100	100
1	E	318/331 (96%)	308 (97%)	9 (3%)	1 (0%)	46	35
1	F	318/331 (96%)	307 (96%)	11 (4%)	0	100	100
1	G	301/331 (91%)	291 (97%)	9 (3%)	1 (0%)	46	35
1	H	301/331 (91%)	290 (96%)	11 (4%)	0	100	100
All	All	2477/2648 (94%)	2391 (96%)	84 (3%)	2 (0%)	56	46

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	G	151	ALA
1	E	151	ALA

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	269/282 (95%)	269 (100%)	0	100	100
1	B	268/282 (95%)	267 (100%)	1 (0%)	93	94
1	C	242/282 (86%)	242 (100%)	0	100	100
1	D	249/282 (88%)	249 (100%)	0	100	100
1	E	265/282 (94%)	265 (100%)	0	100	100
1	F	268/282 (95%)	268 (100%)	0	100	100
1	G	241/282 (86%)	241 (100%)	0	100	100
1	H	247/282 (88%)	247 (100%)	0	100	100
All	All	2049/2256 (91%)	2048 (100%)	1 (0%)	100	100

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	159	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	14	ASN
1	A	25	GLN
1	A	58	ASN
1	F	25	GLN
1	F	127	ASN
1	F	223	ASN
1	E	127	ASN
1	G	25	GLN
1	G	58	ASN
1	G	223	ASN
1	G	286	ASN
1	H	25	GLN
1	H	206	GLN

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 ⓘ

14 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	NAI	A	401	-	38,48,48	1.01	2 (5%)	48,73,73	1.52	5 (10%)
3	PYR	A	402	-	2,5,5	1.81	1 (50%)	2,6,6	0.03	0
2	NAI	B	401	-	38,48,48	1.01	2 (5%)	48,73,73	1.53	5 (10%)
3	PYR	B	402	-	2,5,5	1.77	1 (50%)	2,6,6	0.08	0
2	NAI	C	401	-	38,48,48	1.01	2 (5%)	48,73,73	1.53	5 (10%)
2	NAI	D	401	-	38,48,48	1.02	2 (5%)	48,73,73	1.51	3 (6%)
3	PYR	D	402	-	2,5,5	1.80	1 (50%)	2,6,6	0.06	0
2	NAI	E	401	-	38,48,48	1.02	2 (5%)	48,73,73	1.53	5 (10%)
3	PYR	E	402	-	2,5,5	1.79	1 (50%)	2,6,6	0.04	0
2	NAI	F	401	-	38,48,48	1.02	2 (5%)	48,73,73	1.59	7 (14%)
3	PYR	F	402	-	2,5,5	1.78	1 (50%)	2,6,6	0.17	0
2	NAI	G	401	-	38,48,48	1.02	2 (5%)	48,73,73	1.54	6 (12%)
2	NAI	H	401	-	38,48,48	1.02	2 (5%)	48,73,73	1.55	5 (10%)
3	PYR	H	402	-	2,5,5	1.80	1 (50%)	2,6,6	0.04	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAI	A	401	-	-	0/25/72/72	0/5/5/5
3	PYR	A	402	-	-	0/0/4/4	0/0/0/0
2	NAI	B	401	-	-	0/25/72/72	0/5/5/5
3	PYR	B	402	-	-	0/0/4/4	0/0/0/0
2	NAI	C	401	-	-	0/25/72/72	0/5/5/5
2	NAI	D	401	-	-	0/25/72/72	0/5/5/5
3	PYR	D	402	-	-	0/0/4/4	0/0/0/0
2	NAI	E	401	-	-	0/25/72/72	0/5/5/5
3	PYR	E	402	-	-	0/0/4/4	0/0/0/0
2	NAI	F	401	-	-	0/25/72/72	0/5/5/5
3	PYR	F	402	-	-	0/0/4/4	0/0/0/0
2	NAI	G	401	-	-	0/25/72/72	0/5/5/5
2	NAI	H	401	-	-	0/25/72/72	0/5/5/5
3	PYR	H	402	-	-	0/0/4/4	0/0/0/0

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	402	PYR	O3-C2	2.49	1.31	1.22
3	F	402	PYR	O3-C2	2.51	1.31	1.22
3	E	402	PYR	O3-C2	2.52	1.31	1.22
3	D	402	PYR	O3-C2	2.53	1.31	1.22
3	H	402	PYR	O3-C2	2.53	1.31	1.22
3	A	402	PYR	O3-C2	2.54	1.31	1.22
2	F	401	NAI	C5A-C4A	2.99	1.47	1.40
2	A	401	NAI	C5A-C4A	3.03	1.47	1.40
2	H	401	NAI	C5A-C4A	3.05	1.47	1.40
2	G	401	NAI	C5A-C4A	3.06	1.47	1.40
2	B	401	NAI	C5A-C4A	3.06	1.47	1.40
2	D	401	NAI	C5A-C4A	3.07	1.47	1.40
2	E	401	NAI	C5A-C4A	3.10	1.47	1.40
2	C	401	NAI	C5A-C4A	3.12	1.47	1.40
2	A	401	NAI	C6N-C5N	3.13	1.39	1.33
2	B	401	NAI	C6N-C5N	3.17	1.39	1.33
2	F	401	NAI	C6N-C5N	3.21	1.39	1.33
2	D	401	NAI	C6N-C5N	3.24	1.39	1.33
2	C	401	NAI	C6N-C5N	3.24	1.39	1.33
2	E	401	NAI	C6N-C5N	3.25	1.39	1.33
2	H	401	NAI	C6N-C5N	3.28	1.39	1.33
2	G	401	NAI	C6N-C5N	3.37	1.39	1.33

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	401	NAI	N3A-C2A-N1A	-7.09	123.47	128.89
2	A	401	NAI	N3A-C2A-N1A	-7.03	123.51	128.89
2	B	401	NAI	N3A-C2A-N1A	-7.00	123.53	128.89
2	C	401	NAI	N3A-C2A-N1A	-6.96	123.56	128.89
2	H	401	NAI	N3A-C2A-N1A	-6.96	123.57	128.89
2	F	401	NAI	N3A-C2A-N1A	-6.93	123.59	128.89
2	D	401	NAI	N3A-C2A-N1A	-6.87	123.63	128.89
2	G	401	NAI	N3A-C2A-N1A	-6.81	123.68	128.89
2	C	401	NAI	PN-O3-PA	-3.64	122.52	132.73
2	D	401	NAI	PN-O3-PA	-3.50	122.91	132.73
2	F	401	NAI	PN-O3-PA	-3.46	123.00	132.73
2	B	401	NAI	PN-O3-PA	-3.43	123.10	132.73
2	H	401	NAI	PN-O3-PA	-3.39	123.20	132.73
2	G	401	NAI	C4A-C5A-N7A	-3.37	106.38	109.48
2	E	401	NAI	PN-O3-PA	-3.32	123.42	132.73
2	H	401	NAI	C4A-C5A-N7A	-3.31	106.43	109.48
2	F	401	NAI	C4A-C5A-N7A	-3.30	106.44	109.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	401	NAI	PN-O3-PA	-3.28	123.51	132.73
2	E	401	NAI	C4A-C5A-N7A	-3.26	106.48	109.48
2	A	401	NAI	C4A-C5A-N7A	-3.23	106.50	109.48
2	D	401	NAI	C4A-C5A-N7A	-3.22	106.52	109.48
2	B	401	NAI	C4A-C5A-N7A	-3.09	106.63	109.48
2	C	401	NAI	C4A-C5A-N7A	-3.07	106.65	109.48
2	A	401	NAI	PN-O3-PA	-2.92	124.53	132.73
2	F	401	NAI	C4B-O4B-C1B	-2.81	106.63	109.72
2	C	401	NAI	C2B-C1B-N9A	-2.58	110.36	114.29
2	A	401	NAI	C2B-C1B-N9A	-2.30	110.77	114.29
2	G	401	NAI	C4B-O4B-C1B	-2.24	107.25	109.72
2	F	401	NAI	C1B-N9A-C4A	-2.24	123.56	126.94
2	F	401	NAI	C1D-N1N-C2N	-2.17	117.13	120.91
2	B	401	NAI	C2B-C1B-N9A	-2.12	111.05	114.29
2	H	401	NAI	C4B-O4B-C1B	-2.09	107.42	109.72
2	C	401	NAI	C1D-N1N-C2N	-2.07	117.30	120.91
2	G	401	NAI	C1B-N9A-C4A	-2.03	123.88	126.94
2	E	401	NAI	C2B-C1B-N9A	-2.02	111.20	114.29
2	B	401	NAI	O4B-C1B-N9A	2.04	112.37	108.10
2	E	401	NAI	O4B-C1B-N9A	2.09	112.46	108.10
2	A	401	NAI	O4B-C1B-N9A	2.09	112.48	108.10
2	F	401	NAI	O4B-C1B-N9A	2.10	112.49	108.10
2	H	401	NAI	O4B-C1B-N9A	2.37	113.05	108.10
2	G	401	NAI	O4B-C1B-N9A	2.55	113.43	108.10

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

13 monomers are involved in 28 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	NAI	2	0
2	B	401	NAI	3	0
3	B	402	PYR	3	0
2	C	401	NAI	1	0
2	D	401	NAI	4	0
3	D	402	PYR	1	0
2	E	401	NAI	3	0
3	E	402	PYR	1	0
2	F	401	NAI	5	0
3	F	402	PYR	3	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	G	401	NAI	3	0
2	H	401	NAI	1	0
3	H	402	PYR	1	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	318/331 (96%)	-0.37	0 100 100	28, 40, 56, 64	0
1	B	318/331 (96%)	-0.45	0 100 100	25, 36, 47, 55	0
1	C	303/331 (91%)	-0.30	0 100 100	25, 41, 57, 64	0
1	D	303/331 (91%)	-0.39	0 100 100	27, 40, 53, 61	0
1	E	318/331 (96%)	-0.37	1 (0%) 94 94	29, 39, 54, 61	0
1	F	318/331 (96%)	-0.44	0 100 100	26, 35, 46, 54	0
1	G	303/331 (91%)	-0.35	0 100 100	27, 40, 57, 64	0
1	H	303/331 (91%)	-0.37	0 100 100	29, 40, 54, 66	0
All	All	2484/2648 (93%)	-0.38	1 (0%) 100 100	25, 39, 54, 66	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	318	VAL	2.6

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
3	PYR	D	402	6/6	0.90	0.19	7.10	43,46,48,49	0
3	PYR	B	402	6/6	0.97	0.11	5.30	28,30,34,35	0
3	PYR	H	402	6/6	0.90	0.20	4.57	34,45,50,50	0
3	PYR	F	402	6/6	0.97	0.12	3.98	28,32,34,36	0
3	PYR	E	402	6/6	0.95	0.14	2.57	37,41,45,45	0
3	PYR	A	402	6/6	0.95	0.13	2.04	37,41,44,46	0
2	NAI	B	401	44/44	0.98	0.10	0.92	27,31,42,48	0
2	NAI	D	401	44/44	0.97	0.09	0.42	34,42,46,49	0
2	NAI	G	401	44/44	0.97	0.09	0.30	33,40,46,49	0
2	NAI	C	401	44/44	0.98	0.10	0.29	34,41,44,46	0
2	NAI	F	401	44/44	0.98	0.09	-0.10	25,30,35,39	0
2	NAI	A	401	44/44	0.97	0.09	-0.14	33,39,44,48	0
2	NAI	H	401	44/44	0.97	0.08	-0.19	37,44,46,53	0
2	NAI	E	401	44/44	0.98	0.08	-0.54	35,40,44,47	0

6.5 Other polymers [i](#)

There are no such residues in this entry.