



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:37 PM GMT

PDB ID : 4M55
Title : Crystal structure of Human UDP-xylose synthase R236H substitution
Authors : Walsh Jr., R.M.; Polizzi, S.J.; Wood, Z.A.
Deposited on : 2013-08-08
Resolution : 2.86 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

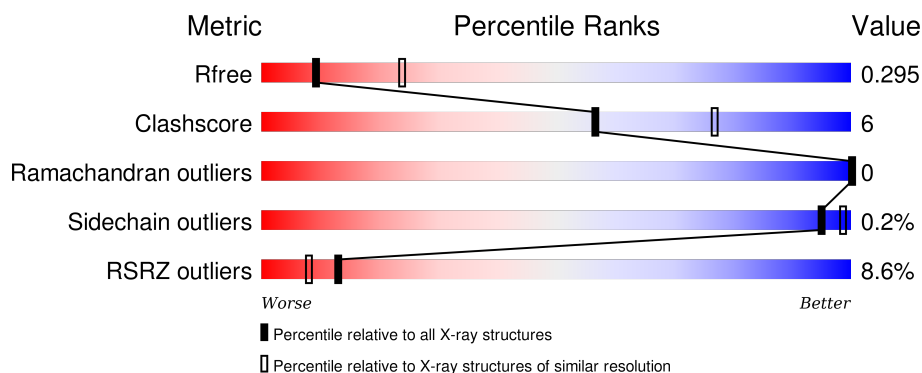
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.86 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.



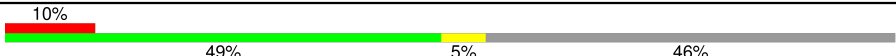
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	2228 (2.90-2.82)
Clashscore	102246	2499 (2.90-2.82)
Ramachandran outliers	100387	2439 (2.90-2.82)
Sidechain outliers	100360	2442 (2.90-2.82)
RSRZ outliers	91569	2236 (2.90-2.82)

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	336	<div> <div>10%</div> <div>66%</div> <div>12%</div> <div>22%</div> </div>
1	B	336	<div> <div>4%</div> <div>71%</div> <div>10%</div> <div>19%</div> </div>
1	C	336	<div> <div>3%</div> <div>69%</div> <div>12%</div> <div>19%</div> </div>
1	D	336	<div> <div>6%</div> <div>62%</div> <div>16%</div> <div>22%</div> </div>
1	E	336	<div> <div>4%</div> <div>39%</div> <div>10%</div> <div>51%</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	336	

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
5	UDP	B	502	-	-	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 11483 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 UDP-glucuronic acid decarboxylase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	263	Total	C	N	O	S	0	0	0
			2086	1332	365	379	10			
1	B	272	Total	C	N	O	S	0	0	0
			2145	1367	374	394	10			
1	C	271	Total	C	N	O	S	0	0	0
			2135	1362	370	393	10			
1	D	261	Total	C	N	O	S	0	0	0
			2070	1323	362	375	10			
1	E	164	Total	C	N	O	S	0	0	0
			1281	816	222	235	8			
1	F	181	Total	C	N	O	S	0	0	0
			1429	917	243	260	9			

There are 6 discrepancies between the modelled and reference sequences:

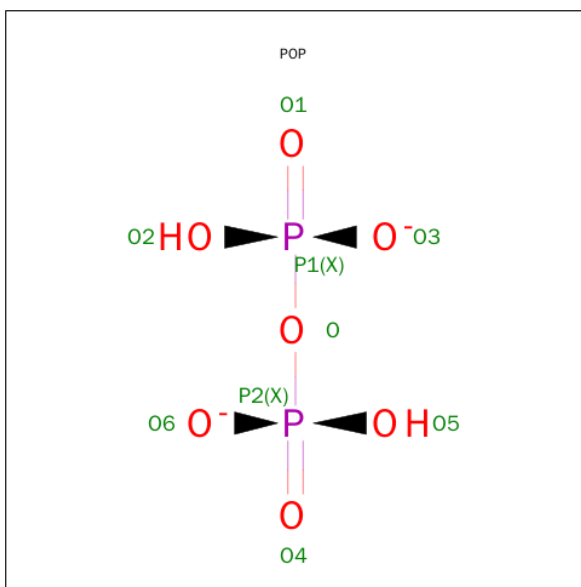
Chain	Residue	Modelled	Actual	Comment	Reference
A	236	HIS	ARG	ENGINEERED MUTATION	UNP Q8NBZ7
B	236	HIS	ARG	ENGINEERED MUTATION	UNP Q8NBZ7
C	236	HIS	ARG	ENGINEERED MUTATION	UNP Q8NBZ7
D	236	HIS	ARG	ENGINEERED MUTATION	UNP Q8NBZ7
E	236	HIS	ARG	ENGINEERED MUTATION	UNP Q8NBZ7
F	236	HIS	ARG	ENGINEERED MUTATION	UNP Q8NBZ7

- Molecule 2 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula: C₂₁H₂₇N₇O₁₄P₂).



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

- Molecule 3 is PYROPHOSPHATE 2- (three-letter code: POP) (formula: H₂O₇P₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	P	0	0
			9	7	2		
3	C	1	Total	O	P	0	0
			9	7	2		

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



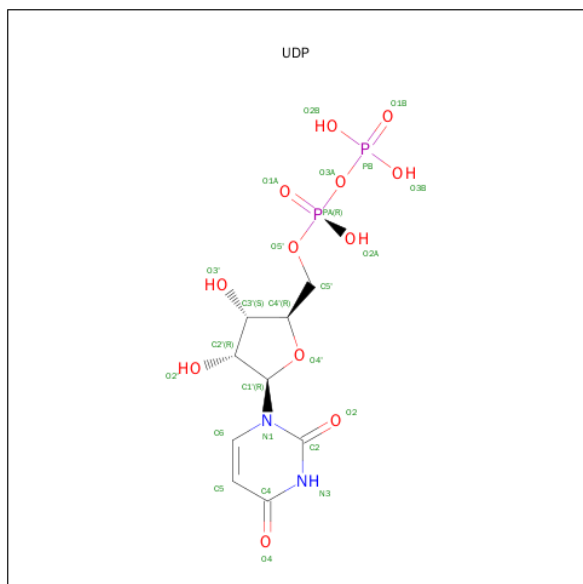
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	C	1	Total	O	S	0	0
			5	4	1		

- Molecule 5 is URIDINE-5'-DIPHOSPHATE (three-letter code: UDP) (formula: $C_9H_{14}N_2O_{12}P_2$).

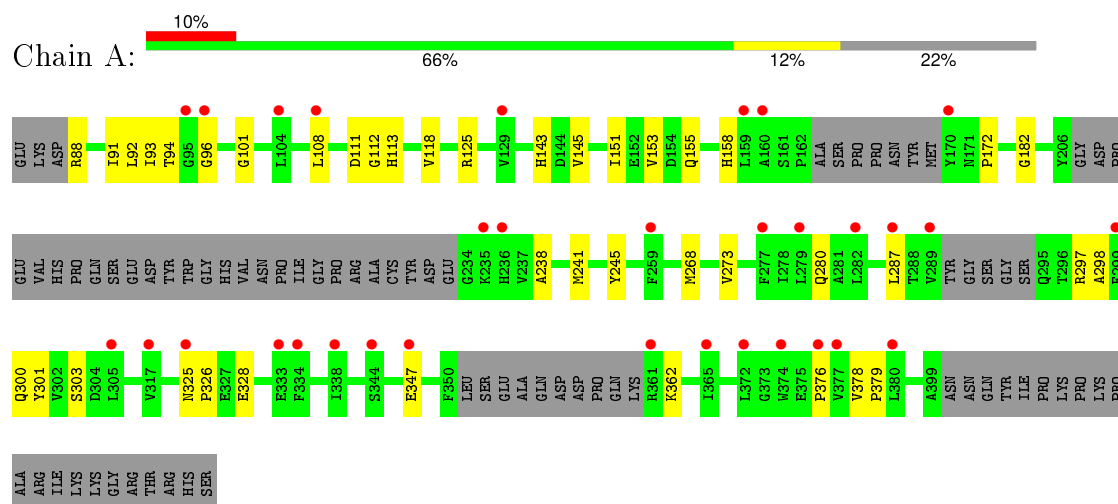


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	B	1	Total	C	N	O	P	0	0
			25	9	2	12	2		
5	D	1	Total	C	O	P		0	0
			15	3	10	2			

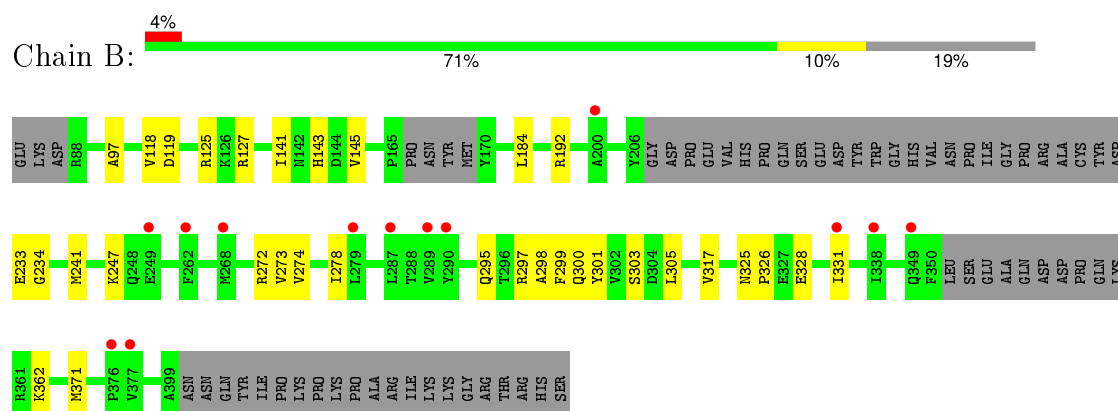
3 Residue-property plots

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.

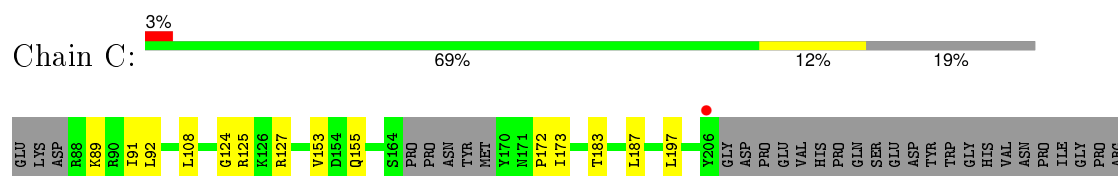
• Molecule 1: UDP-glucuronic acid decarboxylase 1

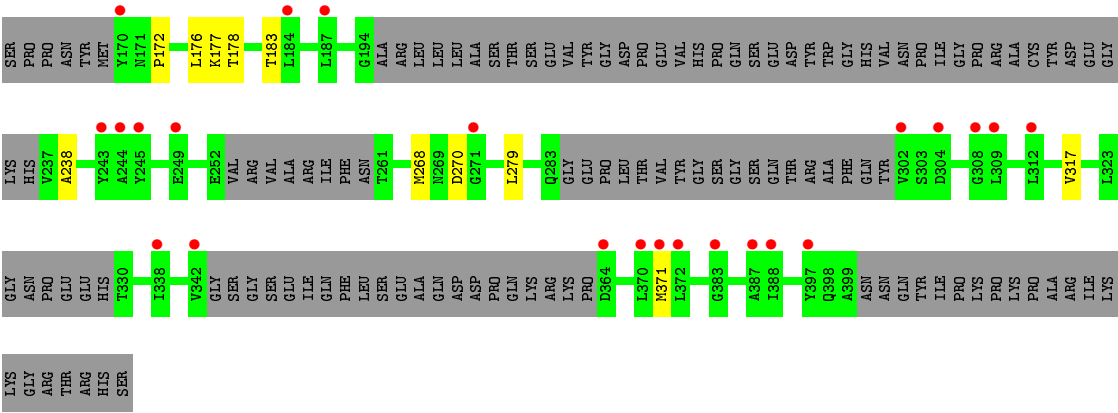


• Molecule 1: UDP-glucuronic acid decarboxylase 1



• Molecule 1: UDP-glucuronic acid decarboxylase 1





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	83.17Å 85.14Å 292.55Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.31 – 2.86 48.22 – 2.86	Depositor EDS
% Data completeness (in resolution range)	99.9 (42.31-2.86) 100.0 (48.22-2.86)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.35 (at 2.86Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.2_1309)	Depositor
R, R_{free}	0.228 , 0.285 0.246 , 0.295	Depositor DCC
R_{free} test set	2450 reflections (5.26%)	DCC
Wilson B-factor (Å ²)	78.4	Xtriage
Anisotropy	0.678	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 79.6	EDS
Estimated twinning fraction	0.156 for k,h,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	2 of 49011 reflections (0.004%)	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	11483	wwPDB-VP
Average B, all atoms (Å ²)	116.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.53% 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: UDP, SO4, NAD, POP

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	0/2125	0.51	0/2868
1	B	0.50	0/2187	0.59	0/2954
1	C	0.51	0/2176	0.59	0/2939
1	D	0.42	0/2108	0.54	0/2844
1	E	0.37	0/1299	0.51	0/1752
1	F	0.34	0/1448	0.46	0/1946
All	All	0.43	0/11343	0.54	0/15303

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	2086	0	2100	26	0
1	B	2145	0	2149	25	0
1	C	2135	0	2140	27	0
1	D	2070	0	2090	36	0
1	E	1281	0	1301	24	0
1	F	1429	0	1445	11	0
2	A	44	0	26	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	44	0	26	1	0
2	C	44	0	26	0	0
2	D	44	0	26	1	0
2	E	44	0	26	1	0
2	F	44	0	26	1	0
3	A	9	0	0	1	0
3	C	9	0	0	0	0
4	A	5	0	0	1	0
4	B	5	0	0	0	0
4	C	5	0	0	0	0
5	B	25	0	11	1	0
5	D	15	0	3	1	0
All	All	11483	0	11395	139	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 6.

All (139) 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:118:VAL:HG13	1:D:143:HIS:HB3	1.58	0.83
1:A:92:LEU:HD22	1:A:151:ILE:HD11	1.69	0.74
1:D:321:VAL:HG22	1:D:367:LYS:HE2	1.71	0.73
1:C:187:LEU:HD23	1:C:197:LEU:HD13	1.70	0.72
1:C:247:LYS:HE3	1:D:172:PRO:HG3	1.74	0.70
1:F:178:THR:OG1	2:F:800:NAD:N6A	2.23	0.69
1:A:172:PRO:HG3	1:B:247:LYS:HE3	1.79	0.64
1:D:238:ALA:HA	1:D:241:MET:HE2	1.79	0.64
1:A:326:PRO:O	1:A:362:LYS:NZ	2.31	0.64
1:C:233:GLU:HG2	1:C:234:GLY:H	1.62	0.63
1:B:272:ARG:HD3	1:B:299:PHE:HE1	1.63	0.63
1:A:301:TYR:HD2	1:A:303:SER:HG	1.47	0.63
1:C:172:PRO:HG3	1:D:247:LYS:HE3	1.82	0.61
1:D:119:ASP:O	1:D:142:ASN:HA	1.99	0.61
1:B:233:GLU:HG2	1:B:234:GLY:H	1.66	0.60
2:A:501:NAD:O2N	2:A:501:NAD:N7N	2.33	0.60
1:C:272:ARG:HD3	1:C:299:PHE:HE1	1.67	0.59
1:A:268:MET:N	4:A:503:SO4:O2	2.36	0.59
1:B:326:PRO:O	1:B:362:LYS:NZ	2.36	0.58
1:D:253:VAL:O	1:D:254:ARG:NH1	2.35	0.58
1:C:297:ARG:HB3	1:C:299:PHE:HE2	1.68	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:93:ILE:HD11	1:A:108:LEU:HD12	1.87	0.57
1:A:238:ALA:HA	1:A:241:MET:HE2	1.85	0.57
1:D:125:ARG:HH21	1:E:127:ARG:HE	1.51	0.57
1:D:295:GLN:O	1:D:331:ILE:HG12	2.04	0.57
1:D:92:LEU:HB2	1:D:153:VAL:HG11	1.86	0.57
1:A:118:VAL:HG13	1:A:143:HIS:HB3	1.85	0.57
1:D:118:VAL:HG21	1:D:149:LEU:HD21	1.86	0.57
1:E:92:LEU:HB2	1:E:153:VAL:HG11	1.85	0.56
1:D:120:ASN:O	1:E:142:ASN:ND2	2.39	0.56
1:C:298:ALA:HB2	1:C:328:GLU:HG2	1.88	0.56
1:D:329:HIS:ND1	1:D:333:GLU:OE1	2.32	0.55
1:B:192:ARG:NE	5:B:502:UDP:O4	2.36	0.55
1:E:248:GLN:OE1	1:F:172:PRO:HD2	2.07	0.54
1:C:297:ARG:HB3	1:C:299:PHE:CE2	2.42	0.54
1:A:325:ASN:HD22	1:A:376:PRO:HG2	1.71	0.54
1:C:92:LEU:HB2	1:C:153:VAL:HG11	1.90	0.54
1:C:233:GLU:HG2	1:C:234:GLY:N	2.22	0.54
1:D:157:TYR:HD2	1:D:198:LEU:HD22	1.73	0.53
1:F:268:MET:HE2	1:F:279:LEU:HD21	1.90	0.53
1:D:105:THR:HG23	1:D:115:VAL:HG11	1.91	0.53
1:D:198:LEU:HD21	1:D:309:LEU:HD22	1.92	0.52
1:B:127:ARG:HE	1:C:127:ARG:HH21	1.57	0.52
1:E:106:ASP:OD1	1:E:132:TRP:NE1	2.39	0.52
1:B:127:ARG:HH21	1:C:127:ARG:HE	1.57	0.52
1:E:259:PHE:HE1	1:E:322:ASN:HB3	1.75	0.52
1:A:298:ALA:HB1	1:A:325:ASN:O	2.09	0.52
1:D:300:GLN:HB3	1:D:325:ASN:HB2	1.92	0.51
1:C:326:PRO:HA	1:C:363:PRO:HG2	1.92	0.51
1:E:259:PHE:CE1	1:E:322:ASN:HB3	2.45	0.51
1:D:366:LYS:HA	1:D:369:LYS:HD2	1.93	0.51
1:A:92:LEU:HB2	1:A:153:VAL:HG11	1.93	0.50
1:E:96:GLY:HA2	1:E:101:GLY:HA3	1.93	0.50
1:B:297:ARG:HB3	1:B:299:PHE:CE2	2.47	0.50
1:A:287:LEU:N	1:A:347:GLU:O	2.44	0.50
1:D:326:PRO:HA	1:D:363:PRO:HG2	1.93	0.50
1:B:298:ALA:HB1	1:B:325:ASN:O	2.12	0.50
1:D:248:GLN:HG2	1:D:249:GLU:HG2	1.93	0.49
1:A:125:ARG:HH22	1:F:127:ARG:H	1.60	0.49
1:E:119:ASP:OD1	1:E:120:ASN:N	2.44	0.49
1:E:162:PRO:HG2	1:E:175:THR:HG23	1.95	0.48
1:A:238:ALA:HA	1:A:241:MET:CE	2.44	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:89:LYS:HE3	1:C:155:GLN:NE2	2.29	0.48
1:B:273:VAL:HG11	1:B:297:ARG:HG3	1.95	0.48
1:B:297:ARG:HB3	1:B:299:PHE:HE2	1.78	0.48
1:A:96:GLY:HA2	1:A:101:GLY:HA3	1.96	0.47
1:B:298:ALA:HB2	1:B:328:GLU:HG2	1.96	0.47
1:E:189:LEU:O	1:E:193:VAL:HG22	2.14	0.47
1:A:245:TYR:OH	3:A:502:POP:O5	2.27	0.47
1:B:273:VAL:HG21	1:B:331:ILE:HD12	1.97	0.47
1:B:145:VAL:HG22	2:B:501:NAD:N1A	2.29	0.47
1:A:94:THR:OG1	1:A:158:HIS:HA	2.14	0.47
1:A:145:VAL:O	1:A:182:GLY:HA2	2.15	0.47
1:C:183:THR:HB	1:C:241:MET:HE1	1.96	0.47
1:F:183:THR:HG21	1:F:238:ALA:HB1	1.97	0.47
1:B:301:TYR:HD2	1:B:303:SER:HG	1.59	0.46
1:E:151:ILE:HG12	1:E:152:GLU:H	1.79	0.46
1:B:274:VAL:O	1:B:278:ILE:HG13	2.14	0.46
1:C:341:LEU:HD21	1:C:381:GLU:HG3	1.96	0.46
1:D:329:HIS:HB2	1:D:334:PHE:CE2	2.51	0.46
1:D:183:THR:HB	1:D:241:MET:HE1	1.98	0.45
1:C:124:GLY:O	1:C:125:ARG:HD3	2.16	0.45
1:E:302:VAL:O	1:E:306:VAL:HG23	2.16	0.45
1:D:157:TYR:CD2	1:D:198:LEU:HD22	2.51	0.45
1:A:273:VAL:HG11	1:A:297:ARG:HG3	1.99	0.45
1:F:270:ASP:N	1:F:270:ASP:OD1	2.48	0.45
1:D:141:ILE:HG22	1:D:143:HIS:HB2	2.00	0.44
1:F:317:VAL:HG21	1:F:371:MET:HB3	2.00	0.44
1:A:298:ALA:HB2	1:A:328:GLU:HG2	2.00	0.44
1:A:280:GLN:OE1	1:A:287:LEU:HA	2.17	0.44
1:E:241:MET:HG2	1:F:176:LEU:HD13	2.00	0.44
1:C:290:TYR:CE1	1:C:351:LEU:HD22	2.52	0.44
1:B:317:VAL:HG21	1:B:371:MET:HB3	1.99	0.44
1:D:100:VAL:HB	2:D:501:NAD:H51N	2.00	0.44
1:A:111:ASP:HB2	1:A:113:HIS:CD2	2.53	0.44
1:B:233:GLU:HG2	1:B:234:GLY:N	2.32	0.44
1:A:300:GLN:HB3	1:A:325:ASN:HB2	1.99	0.44
1:E:254:ARG:NH2	1:E:315:SER:O	2.51	0.43
1:D:118:VAL:HG21	1:D:149:LEU:CD2	2.48	0.43
1:C:298:ALA:HB1	1:C:325:ASN:O	2.18	0.43
1:B:118:VAL:HG13	1:B:143:HIS:HB3	2.00	0.43
1:B:295:GLN:O	1:B:331:ILE:HG12	2.19	0.43
1:C:324:GLY:O	1:C:365:ILE:HG21	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:93:ILE:HD11	1:E:108:LEU:HD12	1.99	0.43
1:C:91:ILE:HD13	1:C:108:LEU:HD13	2.00	0.43
1:D:296:THR:HA	1:D:330:THR:HA	2.00	0.42
1:D:148:PRO:HG3	5:D:502:UDP:O2'	2.19	0.42
1:C:256:ALA:HB2	1:C:312:LEU:HD23	2.00	0.42
1:E:106:ASP:O	1:E:110:MET:HG2	2.19	0.42
1:D:130:GLU:HA	1:D:133:ILE:HG13	2.01	0.42
1:D:328:GLU:HG3	1:D:362:LYS:HG3	2.02	0.42
1:B:125:ARG:NH1	1:B:127:ARG:HH22	2.17	0.42
1:E:184:LEU:HG	1:E:241:MET:SD	2.59	0.42
1:C:273:VAL:HG11	1:C:297:ARG:HG3	2.02	0.42
1:E:119:ASP:OD2	2:E:800:NAD:O3B	2.27	0.42
1:F:177:LYS:HB3	1:F:177:LYS:HE2	1.80	0.42
1:B:300:GLN:OE1	1:B:305:LEU:HB2	2.20	0.42
1:D:106:ASP:OD2	1:D:265:ARG:NH1	2.52	0.42
1:D:182:GLY:O	1:D:186:MET:HG2	2.20	0.42
1:F:102:SER:HB3	1:F:128:ASN:HB3	2.01	0.42
1:B:184:LEU:HG	1:B:241:MET:SD	2.60	0.42
1:C:290:TYR:HD1	1:C:351:LEU:HB2	1.85	0.41
1:F:118:VAL:HG13	1:F:143:HIS:HB3	2.00	0.41
1:C:268:MET:HE2	1:C:391:PHE:CD2	2.55	0.41
1:B:118:VAL:HG22	1:B:141:ILE:HB	2.01	0.41
1:E:122:PHE:HD2	1:E:123:THR:HG23	1.85	0.41
1:E:256:ALA:HB2	1:E:312:LEU:HD23	2.02	0.41
1:C:259:PHE:CD2	1:C:363:PRO:HB3	2.55	0.41
1:D:235:LYS:HB3	1:D:235:LYS:HE3	1.84	0.41
1:B:97:ALA:HB3	1:B:119:ASP:OD2	2.21	0.41
1:A:88:ARG:HD2	1:A:112:GLY:O	2.20	0.41
1:C:173:ILE:HG13	1:D:248:GLN:OE1	2.21	0.41
1:E:151:ILE:HG12	1:E:152:GLU:N	2.35	0.41
1:E:122:PHE:CD2	1:E:123:THR:HG23	2.56	0.41
1:A:91:ILE:HG12	1:A:155:GLN:HB2	2.03	0.41
1:E:202:THR:OG1	1:E:203:SER:N	2.54	0.41
1:D:384:LEU:O	1:D:388:ILE:HG13	2.20	0.40
1:A:378:VAL:HA	1:A:379:PRO:HD2	1.97	0.40
1:D:260:ASN:OD1	1:D:260:ASN:N	2.54	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	253/336 (75%)	244 (96%)	9 (4%)	0	100	100
1	B	264/336 (79%)	258 (98%)	6 (2%)	0	100	100
1	C	263/336 (78%)	255 (97%)	8 (3%)	0	100	100
1	D	251/336 (75%)	238 (95%)	13 (5%)	0	100	100
1	E	154/336 (46%)	146 (95%)	8 (5%)	0	100	100
1	F	163/336 (48%)	158 (97%)	5 (3%)	0	100	100
All	All	1348/2016 (67%)	1299 (96%)	49 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	227/290 (78%)	227 (100%)	0	100	100
1	B	233/290 (80%)	233 (100%)	0	100	100
1	C	232/290 (80%)	231 (100%)	1 (0%)	93	98
1	D	226/290 (78%)	225 (100%)	1 (0%)	93	98
1	E	142/290 (49%)	142 (100%)	0	100	100
1	F	157/290 (54%)	157 (100%)	0	100	100
All	All	1217/1740 (70%)	1215 (100%)	2 (0%)	95	98

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	237	VAL
1	D	237	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 ⓘ

13 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	NAD	A	501	-	38,48,48	0.86	1 (2%)	47,73,73	1.52	4 (8%)
3	POP	A	502	-	8,8,8	0.51	0	13,13,13	1.28	1 (7%)
4	SO4	A	503	-	4,4,4	0.25	0	6,6,6	0.12	0
2	NAD	B	501	-	38,48,48	1.10	3 (7%)	47,73,73	1.54	10 (21%)
5	UDP	B	502	-	18,26,26	1.13	1 (5%)	26,40,40	1.43	2 (7%)
4	SO4	B	503	-	4,4,4	0.15	0	6,6,6	0.16	0
2	NAD	C	501	-	38,48,48	0.85	1 (2%)	47,73,73	1.72	7 (14%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	POP	C	502	-	8,8,8	0.60	0	13,13,13	1.16	1 (7%)
4	SO4	C	503	-	4,4,4	0.22	0	6,6,6	0.23	0
2	NAD	D	501	-	38,48,48	0.85	2 (5%)	47,73,73	1.73	5 (10%)
5	UDP	D	502	-	11,13,26	1.00	0	15,19,40	1.04	1 (6%)
2	NAD	E	800	-	38,48,48	0.98	3 (7%)	47,73,73	1.64	8 (17%)
2	NAD	F	800	-	38,48,48	0.80	1 (2%)	47,73,73	1.68	6 (12%)

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	NAD	A	501	-	-	0/22/62/62	0/5/5/5
3	POP	A	502	-	-	0/6/6/6	0/0/0/0
4	SO4	A	503	-	-	0/0/0/0	0/0/0/0
2	NAD	B	501	-	-	0/22/62/62	0/5/5/5
5	UDP	B	502	-	-	0/12/32/32	0/2/2/2
4	SO4	B	503	-	-	0/0/0/0	0/0/0/0
2	NAD	C	501	-	-	0/22/62/62	0/5/5/5
3	POP	C	502	-	-	0/6/6/6	0/0/0/0
4	SO4	C	503	-	-	0/0/0/0	0/0/0/0
2	NAD	D	501	-	-	0/22/62/62	0/5/5/5
5	UDP	D	502	-	-	0/14/14/32	0/0/0/2
2	NAD	E	800	-	-	0/22/62/62	0/5/5/5
2	NAD	F	800	-	-	0/22/62/62	0/5/5/5

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	501	NAD	O4D-C1D	2.12	1.43	1.41
2	E	800	NAD	C2A-N3A	2.21	1.36	1.32
2	E	800	NAD	O4D-C1D	2.40	1.44	1.41
2	B	501	NAD	O4B-C1B	2.50	1.44	1.41
5	B	502	UDP	C4-N3	2.78	1.38	1.33
2	B	501	NAD	C2A-N3A	2.80	1.37	1.32
2	C	501	NAD	C5A-C4A	2.85	1.46	1.40
2	F	800	NAD	C5A-C4A	2.91	1.47	1.40
2	A	501	NAD	C5A-C4A	3.09	1.47	1.40
2	D	501	NAD	C5A-C4A	3.12	1.47	1.40
2	E	800	NAD	C5A-C4A	3.19	1.47	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	NAD	C5A-C4A	3.84	1.49	1.40

All (45) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	501	NAD	N3A-C2A-N1A	-8.50	122.38	128.89
2	F	800	NAD	N3A-C2A-N1A	-7.74	122.97	128.89
2	A	501	NAD	N3A-C2A-N1A	-7.29	123.31	128.89
2	C	501	NAD	N3A-C2A-N1A	-6.73	123.74	128.89
2	E	800	NAD	N3A-C2A-N1A	-6.55	123.88	128.89
2	C	501	NAD	C1B-N9A-C4A	-3.87	121.10	126.94
2	B	501	NAD	N3A-C2A-N1A	-3.68	126.07	128.89
3	A	502	POP	P2-O-P1	-3.55	122.77	132.73
2	C	501	NAD	C4A-C5A-N7A	-3.35	106.40	109.48
5	D	502	UDP	PA-O3A-PB	-3.16	122.06	132.67
2	B	501	NAD	PN-O3-PA	-3.16	123.85	132.73
2	F	800	NAD	C2B-C1B-N9A	-3.14	109.50	114.29
2	C	501	NAD	PN-O3-PA	-3.13	123.95	132.73
2	A	501	NAD	C4A-C5A-N7A	-3.11	106.62	109.48
2	B	501	NAD	O3-PN-O5D	-3.05	94.84	102.94
3	C	502	POP	P2-O-P1	-3.05	124.17	132.73
2	D	501	NAD	PN-O3-PA	-3.04	124.20	132.73
2	D	501	NAD	C1B-N9A-C4A	-2.98	122.44	126.94
2	F	800	NAD	C4A-C5A-N7A	-2.96	106.75	109.48
2	E	800	NAD	O3-PN-O5D	-2.95	95.11	102.94
2	E	800	NAD	C1B-N9A-C4A	-2.95	122.49	126.94
2	E	800	NAD	PN-O3-PA	-2.83	124.78	132.73
2	F	800	NAD	PN-O3-PA	-2.78	124.94	132.73
2	B	501	NAD	C5D-C4D-C3D	-2.70	104.48	115.21
2	E	800	NAD	C4A-C5A-N7A	-2.66	107.03	109.48
2	B	501	NAD	C4A-C5A-N7A	-2.51	107.17	109.48
5	B	502	UDP	PA-O3A-PB	-2.45	124.46	132.67
2	E	800	NAD	O3-PA-O5B	-2.37	96.66	102.94
2	F	800	NAD	C1B-N9A-C4A	-2.35	123.40	126.94
2	C	501	NAD	O3-PN-O5D	-2.16	97.22	102.94
2	A	501	NAD	PN-O3-PA	-2.11	126.79	132.73
2	B	501	NAD	C5N-C4N-C3N	-2.01	117.81	120.33
2	A	501	NAD	C2A-N1A-C6A	2.01	122.35	118.77
2	B	501	NAD	O4B-C4B-C5B	2.05	116.66	109.32
2	F	800	NAD	C2A-N1A-C6A	2.09	122.50	118.77
2	E	800	NAD	C2D-C3D-C4D	2.10	106.93	102.61
2	D	501	NAD	O4D-C1D-N1N	2.11	110.45	108.13

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	800	NAD	O4D-C1D-N1N	2.19	110.54	108.13
2	B	501	NAD	O2N-PN-O1N	2.35	125.27	112.53
2	C	501	NAD	C3N-C7N-N7N	2.44	120.49	117.82
2	D	501	NAD	C2A-N1A-C6A	2.53	123.28	118.77
2	B	501	NAD	C2N-C3N-C4N	2.64	121.23	118.29
2	C	501	NAD	O4D-C1D-N1N	2.79	111.19	108.13
2	B	501	NAD	O4D-C1D-N1N	3.98	112.50	108.13
5	B	502	UDP	C4-N3-C2	5.60	119.68	114.14

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

9 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	NAD	1	0
3	A	502	POP	1	0
4	A	503	SO4	1	0
2	B	501	NAD	1	0
5	B	502	UDP	1	0
2	D	501	NAD	1	0
5	D	502	UDP	1	0
2	E	800	NAD	1	0
2	F	800	NAD	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

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	263/336 (78%)	0.84	32 (12%) 5 3	52, 131, 173, 208	0
1	B	272/336 (80%)	0.55	13 (4%) 34 28	47, 88, 152, 174	0
1	C	271/336 (80%)	0.55	10 (3%) 45 38	47, 91, 143, 169	0
1	D	261/336 (77%)	0.66	21 (8%) 15 10	47, 110, 182, 212	0
1	E	164/336 (48%)	0.62	13 (7%) 15 10	84, 125, 169, 205	0
1	F	181/336 (53%)	0.97	33 (18%) 2 1	85, 142, 175, 184	0
All	All	1412/2016 (70%)	0.68	122 (8%) 13 8	47, 112, 170, 212	0

All (122) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	290	TYR	7.4
1	F	309	LEU	6.5
1	C	351	LEU	6.2
1	A	376	PRO	6.2
1	F	364	ASP	5.6
1	F	397	TYR	4.6
1	A	334	PHE	4.4
1	F	243	TYR	4.4
1	D	283	GLN	4.4
1	A	338	ILE	4.3
1	F	104	LEU	4.0
1	A	333	GLU	3.9
1	A	289	VAL	3.9
1	E	151	ILE	3.7
1	A	299	PHE	3.7
1	D	277	PHE	3.6
1	F	372	LEU	3.6
1	D	388	ILE	3.5
1	F	371	MET	3.4

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Mol	Chain	Res	Type	RSRZ
1	B	289	VAL	3.4
1	D	392	ARG	3.2
1	F	150	TYR	3.2
1	F	245	TYR	3.2
1	F	244	ALA	3.1
1	A	377	VAL	3.1
1	D	328	GLU	3.1
1	B	349	GLN	3.1
1	F	308	GLY	3.1
1	F	370	LEU	3.0
1	E	111	ASP	3.0
1	D	204	GLU	3.0
1	F	302	VAL	3.0
1	A	365	ILE	3.0
1	E	242	CYS	3.0
1	E	99	PHE	3.0
1	D	338	ILE	3.0
1	A	108	LEU	2.9
1	F	144	ASP	2.9
1	F	338	ILE	2.9
1	A	325	ASN	2.9
1	D	205	VAL	2.8
1	C	299	PHE	2.8
1	B	376	PRO	2.8
1	F	148	PRO	2.8
1	A	317	VAL	2.8
1	E	113	HIS	2.8
1	D	349	GLN	2.8
1	D	299	PHE	2.8
1	B	290	TYR	2.8
1	A	372	LEU	2.8
1	F	187	LEU	2.8
1	A	305	LEU	2.7
1	B	262	PHE	2.7
1	A	170	TYR	2.7
1	E	162	PRO	2.7
1	B	279	LEU	2.7
1	A	96	GLY	2.7
1	A	236	HIS	2.6
1	C	206	TYR	2.6
1	D	268	MET	2.6
1	F	388	ILE	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	160	ALA	2.6
1	F	304	ASP	2.6
1	C	334	PHE	2.6
1	E	155	GLN	2.6
1	A	104	LEU	2.6
1	A	235	LYS	2.6
1	F	312	LEU	2.6
1	B	200	ALA	2.5
1	A	374	TRP	2.5
1	F	145	VAL	2.5
1	D	334	PHE	2.5
1	F	170	TYR	2.5
1	A	361	ARG	2.5
1	A	282	LEU	2.5
1	D	288	THR	2.5
1	D	113	HIS	2.5
1	F	118	VAL	2.5
1	E	110	MET	2.4
1	B	377	VAL	2.4
1	A	344	SER	2.4
1	A	129	VAL	2.4
1	B	287	LEU	2.4
1	F	141	ILE	2.4
1	F	387	ALA	2.4
1	A	277	PHE	2.3
1	D	273	VAL	2.3
1	B	249	GLU	2.3
1	D	170	TYR	2.3
1	F	383	GLY	2.3
1	C	350	PHE	2.3
1	A	347	GLU	2.3
1	B	268	MET	2.3
1	F	129	VAL	2.3
1	C	397	TYR	2.2
1	E	313	MET	2.2
1	D	332	LEU	2.2
1	F	249	GLU	2.2
1	F	121	PHE	2.2
1	F	271	GLY	2.2
1	B	331	ILE	2.2
1	A	380	LEU	2.2
1	F	184	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	338	ILE	2.2
1	C	287	LEU	2.2
1	D	236	HIS	2.2
1	A	259	PHE	2.1
1	A	95	GLY	2.1
1	E	255	VAL	2.1
1	C	362	LYS	2.1
1	F	342	VAL	2.1
1	A	159	LEU	2.1
1	F	143	HIS	2.1
1	C	332	LEU	2.1
1	A	279	LEU	2.1
1	E	302	VAL	2.1
1	A	287	LEU	2.0
1	D	305	LEU	2.0
1	D	309	LEU	2.0
1	E	321	VAL	2.0
1	E	306	VAL	2.0
1	D	274	VAL	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
5	UDP	B	502	25/25	0.93	0.30	2.17	101,130,171,323	0
2	NAD	B	501	44/44	0.97	0.20	-0.58	43,72,87,276	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	NAD	E	800	44/44	0.91	0.20	-0.64	70,125,168,299	0
2	NAD	C	501	44/44	0.97	0.19	-0.75	49,76,95,183	0
2	NAD	F	800	44/44	0.92	0.19	-0.89	82,123,143,153	0
2	NAD	A	501	44/44	0.96	0.18	-0.90	52,107,118,120	0
2	NAD	D	501	44/44	0.95	0.18	-1.02	63,99,118,122	0
5	UDP	D	502	15/25	0.89	0.15	-1.55	100,135,199,289	0
3	POP	A	502	9/9	0.85	0.15	-2.75	111,120,154,274	1
3	POP	C	502	9/9	0.89	0.30	-	103,111,174,175	1
4	SO4	A	503	5/5	0.95	0.17	-	164,165,168,168	0
4	SO4	B	503	5/5	0.91	0.20	-	125,131,133,141	0
4	SO4	C	503	5/5	0.90	0.16	-	131,133,138,146	0

6.5 Other polymers ⓘ

There are no such residues in this entry.