



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 04:42 AM GMT

PDB ID : 2NVB
Title : Contribution of Pro275 to the Thermostability of the Alcohol Dehydrogenases (ADHs)
Authors : Goihberg, E.; Tel-Or, S.; Peretz, M.; Frolov, F.; Dym, O.; Burstein, Y.; Israel Structural Proteomics Center (ISPC)
Deposited on : 2006-11-12
Resolution : 2.80 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

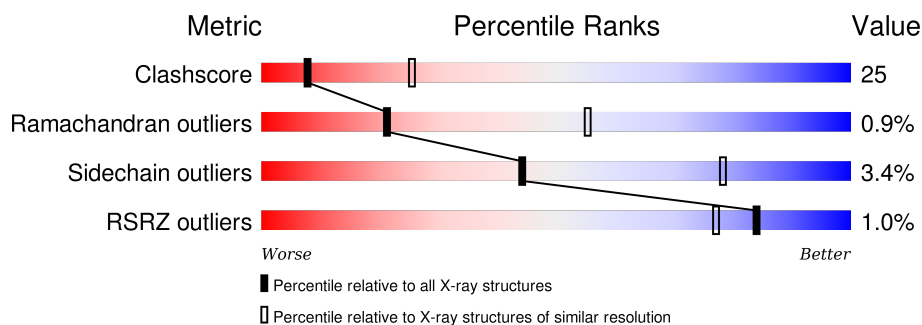
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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))
Clashscore	102246	2827 (2.80-2.80)
Ramachandran outliers	100387	2782 (2.80-2.80)
Sidechain outliers	100360	2784 (2.80-2.80)
RSRZ outliers	91569	2404 (2.80-2.80)

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	352	
1	B	352	
1	C	352	
1	D	352	

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	NAP	A	353	-	-	-	X
3	NAP	C	2353	-	-	-	X

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10772 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 NADP-dependent alcohol dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	352	Total	C	N	O	S	0	0	0
			2644	1692	455	478	19			
1	B	352	Total	C	N	O	S	0	0	0
			2644	1692	455	478	19			
1	C	352	Total	C	N	O	S	0	0	0
			2644	1692	455	478	19			
1	D	352	Total	C	N	O	S	0	0	0
			2644	1692	455	478	19			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	275	ASP	PRO	ENGINEERED	UNP P14941
B	275	ASP	PRO	ENGINEERED	UNP P14941
C	275	ASP	PRO	ENGINEERED	UNP P14941
D	275	ASP	PRO	ENGINEERED	UNP P14941

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Zn	0	0
			1	1		
2	A	1	Total	Zn	0	0
			1	1		
2	D	1	Total	Zn	0	0
			1	1		
2	C	1	Total	Zn	0	0
			1	1		

- Molecule 3 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NAP) (formula: C₂₁H₂₈N₇O₁₇P₃).

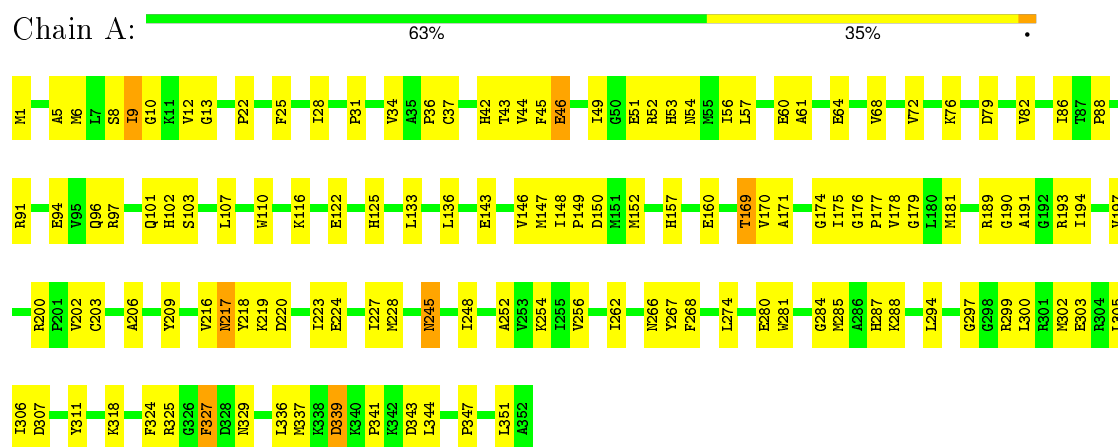


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
3	B	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
3	C	1	Total	C	N	O	P	0	0
			48	21	7	17	3		
3	D	1	Total	C	N	O	P	0	0
			48	21	7	17	3		

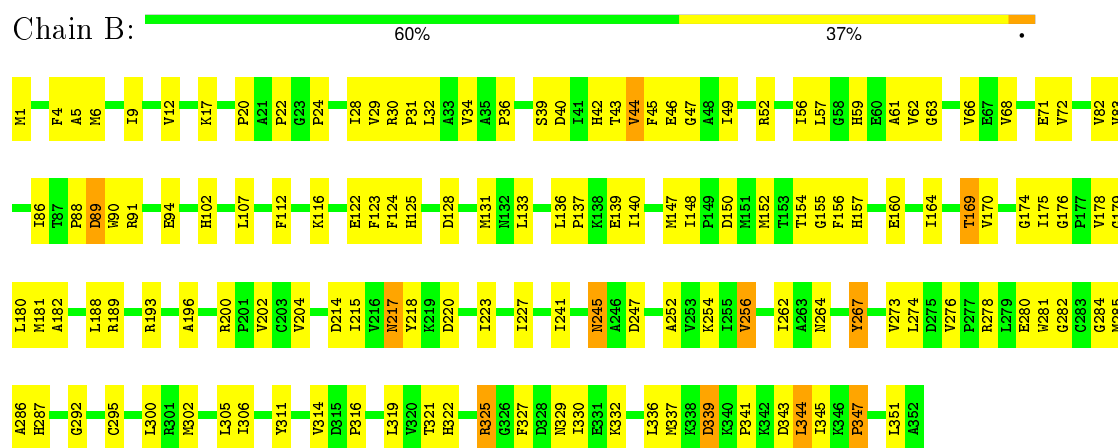
3 Residue-property plots

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 ($\text{RSRZ} > 2$). 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.

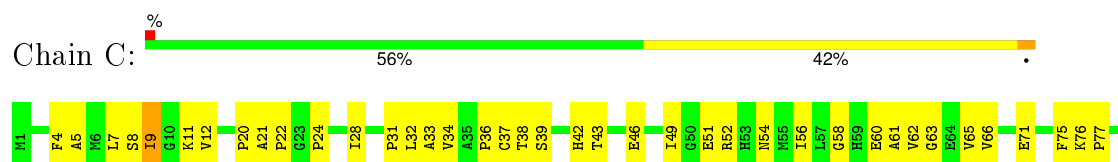
• Molecule 1: NADP-dependent alcohol dehydrogenase

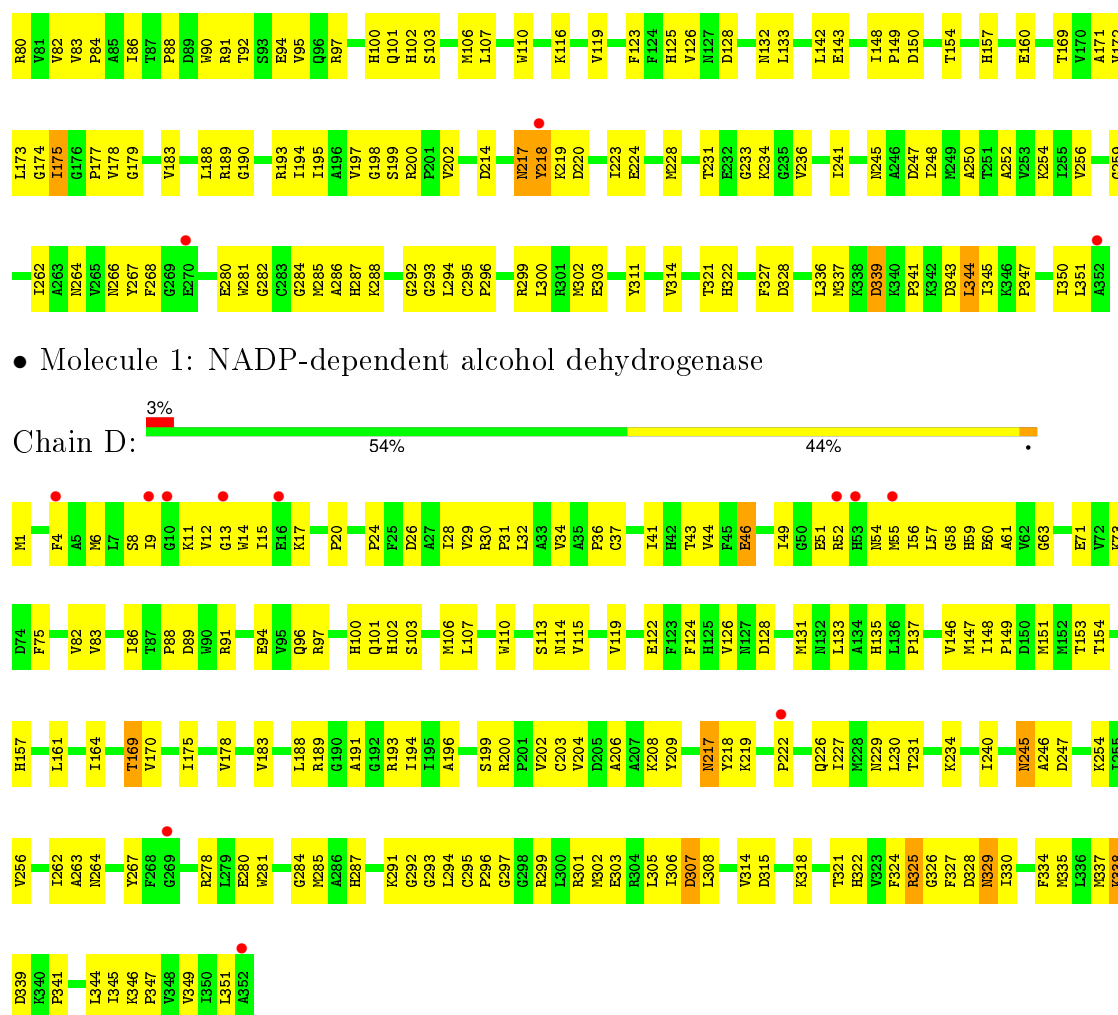


• Molecule 1: NADP-dependent alcohol dehydrogenase



• Molecule 1: NADP-dependent alcohol dehydrogenase





• Molecule 1: NADP-dependent alcohol dehydrogenase

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	79.60Å 125.01Å 167.11Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.60 – 2.80 39.62 – 2.80	Depositor EDS
% Data completeness (in resolution range)	96.4 (39.60-2.80) 92.0 (39.62-2.80)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.90 (at 2.81Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.217 , 0.278 0.234 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	44.5	Xtriage
Anisotropy	0.141	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 30.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 40410 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	10772	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.21% 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: ZN, NAP

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	0.39	0/2701	0.68	1/3653 (0.0%)
1	B	0.38	0/2701	0.68	0/3653
1	C	0.37	0/2701	0.65	0/3653
1	D	0.35	0/2701	0.64	0/3653
All	All	0.37	0/10804	0.66	1/14612 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	176	GLY	N-CA-C	-5.16	100.21	113.10

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	2644	0	2686	132	0
1	B	2644	0	2686	140	0
1	C	2644	0	2686	151	0
1	D	2644	0	2687	183	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	48	0	25	9	0
3	B	48	0	25	6	0
3	C	48	0	25	7	0
3	D	48	0	25	8	0
All	All	10772	0	10845	540	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 25.

All (540) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:189:ARG:NH2	1:C:189:ARG:HH21	1.59	0.98
1:A:49:ILE:HG22	1:A:52:ARG:HH22	1.27	0.98
1:B:102:HIS:HD2	1:B:107:LEU:H	1.05	0.95
1:C:49:ILE:HG22	1:C:52:ARG:HH22	1.34	0.93
1:A:217:ASN:ND2	1:A:219:LYS:H	1.68	0.92
1:D:49:ILE:HG22	1:D:52:ARG:HH22	1.33	0.91
1:C:157:HIS:HE1	1:D:287:HIS:HE1	1.20	0.89
1:B:218:TYR:HB3	1:B:223:ILE:HD11	1.55	0.88
1:B:102:HIS:CD2	1:B:107:LEU:H	1.93	0.86
1:B:178:VAL:HG21	3:B:1353:NAP:H6N	1.56	0.85
1:B:252:ALA:HB1	1:B:262:ILE:HD12	1.59	0.85
1:D:245:ASN:ND2	1:D:246:ALA:H	1.73	0.85
1:A:86:ILE:HG22	1:A:88:PRO:HD3	1.57	0.85
1:C:22:PRO:HD3	1:C:28:ILE:HD11	1.59	0.85
1:A:256:VAL:HG21	1:A:262:ILE:HD11	1.58	0.85
1:C:102:HIS:HE1	1:D:287:HIS:HD2	1.26	0.83
1:B:204:VAL:HG13	1:B:215:ILE:HG12	1.58	0.83
1:D:86:ILE:HB	1:D:297:GLY:HA3	1.58	0.83
1:D:245:ASN:HD22	1:D:246:ALA:H	1.26	0.83
1:A:9:ILE:HG13	1:A:51:GLU:HG3	1.61	0.82
1:C:94:GLU:HG2	1:C:102:HIS:O	1.78	0.82
1:D:59:HIS:NE2	1:D:113:SER:HB3	1.96	0.80
1:D:59:HIS:CD2	1:D:113:SER:HB3	2.15	0.80
1:C:157:HIS:HE1	1:D:287:HIS:CE1	1.99	0.79
1:D:281:TRP:CE3	1:D:284:GLY:HA2	2.16	0.79
1:D:209:TYR:CE2	1:D:318:LYS:HG3	2.17	0.79
1:C:178:VAL:HG21	3:C:2353:NAP:H6N	1.65	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:102:HIS:HE1	1:D:287:HIS:CD2	2.01	0.78
1:A:344:LEU:HD11	1:A:347:PRO:HD3	1.65	0.78
1:B:189:ARG:NH2	1:C:189:ARG:NH2	2.31	0.78
1:C:287:HIS:CD2	1:D:102:HIS:HE1	2.02	0.78
1:A:189:ARG:NH2	1:D:189:ARG:NH2	2.32	0.78
1:D:94:GLU:HG2	1:D:102:HIS:O	1.84	0.77
1:B:189:ARG:HH21	1:C:189:ARG:HH21	1.31	0.77
1:A:189:ARG:HH21	1:D:189:ARG:NH2	1.81	0.77
1:D:49:ILE:HG22	1:D:52:ARG:NH2	1.99	0.77
1:D:36:PRO:HG2	1:D:337:MET:HE2	1.67	0.77
1:A:34:VAL:HG12	1:A:61:ALA:HB2	1.66	0.77
1:A:102:HIS:CD2	1:A:107:LEU:H	2.03	0.77
1:B:152:MET:HG2	1:B:305:LEU:HD13	1.65	0.77
1:A:287:HIS:HE1	1:B:157:HIS:HE1	1.28	0.77
1:A:178:VAL:HG21	3:A:353:NAP:O4D	1.84	0.76
1:C:42:HIS:O	1:C:46:GLU:HB2	1.85	0.76
1:A:287:HIS:HD2	1:B:102:HIS:HE1	1.33	0.76
1:C:38:THR:HB	3:C:2353:NAP:O3D	1.86	0.76
1:A:169:THR:HB	1:A:193:ARG:HB3	1.66	0.76
1:C:157:HIS:CE1	1:D:287:HIS:HE1	2.03	0.75
1:C:22:PRO:HD3	1:C:28:ILE:CD1	2.15	0.75
1:B:31:PRO:HB2	1:B:351:LEU:HD12	1.67	0.75
1:C:177:PRO:HD2	3:C:2353:NAP:O1N	1.85	0.75
1:B:169:THR:HB	1:B:193:ARG:HB3	1.69	0.75
1:C:245:ASN:HB2	1:C:247:ASP:OD1	1.87	0.74
1:A:102:HIS:HD2	1:A:107:LEU:H	1.36	0.74
1:B:189:ARG:HH21	1:C:189:ARG:NH2	1.86	0.73
1:B:82:VAL:HG12	1:B:302:MET:HE2	1.70	0.73
1:A:267:TYR:N	3:A:353:NAP:O2D	2.21	0.73
1:C:281:TRP:CE3	1:C:284:GLY:HA2	2.24	0.73
1:D:49:ILE:HG12	1:D:267:TYR:OH	1.87	0.72
1:C:102:HIS:CE1	1:D:287:HIS:HD2	2.08	0.72
1:D:36:PRO:HG2	1:D:337:MET:CE	2.19	0.72
1:A:217:ASN:HD22	1:A:219:LYS:H	1.38	0.72
1:C:287:HIS:HD2	1:D:102:HIS:HE1	1.36	0.71
1:B:148:ILE:HG22	1:B:302:MET:HE1	1.71	0.71
1:A:49:ILE:HG22	1:A:52:ARG:NH2	2.03	0.71
1:D:86:ILE:HG22	1:D:88:PRO:HD3	1.71	0.71
1:A:202:VAL:HG21	1:A:343:ASP:HA	1.73	0.71
1:A:217:ASN:HD22	1:A:218:TYR:N	1.89	0.70
1:B:22:PRO:HD3	1:B:28:ILE:HD11	1.72	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:31:PRO:HB2	1:D:351:LEU:HD12	1.72	0.70
1:B:254:LYS:HG2	1:B:280:GLU:HG2	1.73	0.70
1:D:14:TRP:CZ3	1:D:327:PHE:CE2	2.80	0.69
1:B:32:LEU:HG	1:B:63:GLY:HA2	1.75	0.69
1:A:287:HIS:CD2	1:B:102:HIS:HE1	2.09	0.69
1:B:302:MET:O	1:B:306:ILE:HG12	1.92	0.69
1:A:287:HIS:CE1	1:B:157:HIS:HE1	2.10	0.69
1:B:217:ASN:HD22	1:B:218:TYR:N	1.91	0.69
1:D:9:ILE:HG23	1:D:51:GLU:HG3	1.74	0.69
1:C:80:ARG:HH11	1:C:80:ARG:HG2	1.58	0.69
1:A:94:GLU:HG2	1:A:102:HIS:O	1.93	0.69
1:A:175:ILE:N	3:A:353:NAP:O3B	2.24	0.69
1:A:217:ASN:C	1:A:217:ASN:HD22	1.97	0.69
1:A:45:PHE:O	1:A:46:GLU:HG3	1.93	0.69
1:D:57:LEU:HA	1:D:113:SER:HB2	1.76	0.68
1:A:149:PRO:HA	1:A:302:MET:HE1	1.74	0.68
1:C:157:HIS:HA	1:C:160:GLU:OE1	1.94	0.68
1:D:245:ASN:HB3	1:D:247:ASP:OD1	1.93	0.68
1:B:22:PRO:HD3	1:B:28:ILE:CD1	2.23	0.68
1:B:42:HIS:O	1:B:46:GLU:HB2	1.92	0.68
1:D:178:VAL:HB	3:D:3353:NAP:H52N	1.75	0.68
1:A:82:VAL:HG12	1:A:302:MET:HE2	1.75	0.68
1:D:267:TYR:HB3	3:D:3353:NAP:O2D	1.92	0.68
1:C:287:HIS:CE1	1:D:157:HIS:HE1	2.12	0.68
1:C:179:GLY:O	1:C:183:VAL:HG23	1.95	0.67
1:A:31:PRO:HB2	1:A:351:LEU:HD12	1.77	0.67
1:D:12:VAL:HG12	1:D:13:GLY:N	2.09	0.67
1:D:59:HIS:NE2	1:D:113:SER:CB	2.58	0.67
1:B:267:TYR:HB3	3:B:1353:NAP:O2D	1.94	0.67
1:C:66:VAL:HG21	1:C:123:PHE:HE1	1.60	0.67
1:D:37:CYS:HB2	1:D:60:GLU:OE2	1.96	0.66
1:A:34:VAL:HG12	1:A:61:ALA:CB	2.25	0.66
1:A:281:TRP:CE3	1:A:284:GLY:HA2	2.31	0.66
1:D:324:PHE:HB2	1:D:349:VAL:HG22	1.78	0.66
1:C:218:TYR:HB3	1:C:223:ILE:HD11	1.77	0.66
1:A:174:GLY:O	1:A:179:GLY:HA3	1.95	0.65
1:D:102:HIS:CD2	1:D:107:LEU:H	2.14	0.65
1:C:9:ILE:HD12	1:C:51:GLU:OE2	1.97	0.65
1:D:49:ILE:CG2	1:D:52:ARG:HH22	2.06	0.65
1:D:217:ASN:HD22	1:D:218:TYR:N	1.94	0.65
1:C:344:LEU:HD11	1:C:347:PRO:HD3	1.79	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:256:VAL:HG21	1:D:262:ILE:HD11	1.79	0.65
1:B:49:ILE:HG22	1:B:52:ARG:HH22	1.60	0.65
1:A:36:PRO:HG2	1:A:337:MET:CE	2.27	0.64
1:C:200:ARG:NH1	1:C:344:LEU:O	2.30	0.64
1:A:217:ASN:HD21	1:A:219:LYS:HG3	1.63	0.64
1:B:49:ILE:HG22	1:B:49:ILE:O	1.97	0.64
1:A:157:HIS:HE1	1:B:287:HIS:CE1	2.15	0.64
1:C:231:THR:O	1:C:234:LYS:HG3	1.98	0.64
1:B:188:LEU:HD11	1:B:314:VAL:HG21	1.80	0.64
1:D:200:ARG:HG3	3:D:3353:NAP:P2B	2.38	0.63
1:D:245:ASN:HD22	1:D:246:ALA:N	1.96	0.63
1:B:148:ILE:HG22	1:B:302:MET:CE	2.28	0.63
1:A:209:TYR:CE2	1:A:318:LYS:HG3	2.34	0.63
1:B:252:ALA:O	1:B:256:VAL:HG13	1.98	0.63
1:B:178:VAL:HG21	3:B:1353:NAP:H52N	1.80	0.63
1:A:200:ARG:NH1	1:A:203:CYS:SG	2.72	0.63
1:B:116:LYS:NZ	1:B:125:HIS:HD2	1.98	0.62
1:C:173:LEU:HD12	1:C:241:ILE:HG12	1.81	0.62
1:A:287:HIS:HD2	1:B:102:HIS:CE1	2.17	0.62
1:B:178:VAL:CG2	3:B:1353:NAP:H52N	2.29	0.62
1:D:178:VAL:HG21	3:D:3353:NAP:H6N	1.81	0.62
1:A:31:PRO:HB2	1:A:351:LEU:CD1	2.30	0.62
1:B:1:MET:HG3	1:B:122:GLU:HB2	1.82	0.62
1:C:256:VAL:HG11	1:C:262:ILE:HD13	1.82	0.62
1:D:254:LYS:HG2	1:D:280:GLU:HG2	1.82	0.62
1:B:329:ASN:OD1	1:B:332:LYS:HE3	1.99	0.61
1:A:86:ILE:HB	1:A:297:GLY:HA3	1.81	0.61
1:A:189:ARG:NH2	1:D:189:ARG:HH21	1.98	0.61
1:A:143:GLU:H	1:A:143:GLU:CD	2.03	0.61
1:D:299:ARG:O	1:D:303:GLU:HG3	2.00	0.61
1:A:267:TYR:HB2	3:A:353:NAP:O2D	2.00	0.61
1:D:245:ASN:ND2	1:D:246:ALA:N	2.44	0.61
1:B:202:VAL:HG21	1:B:343:ASP:HA	1.81	0.61
1:A:146:VAL:HG13	1:A:147:MET:HE2	1.83	0.61
1:B:300:LEU:CD2	1:D:97:ARG:HG2	2.31	0.61
1:C:287:HIS:HD2	1:D:102:HIS:CE1	2.19	0.61
1:D:107:LEU:HD23	1:D:294:LEU:HD22	1.82	0.60
1:C:321:THR:HG21	1:C:345:ILE:HD13	1.81	0.60
1:D:32:LEU:HB2	1:D:63:GLY:HA2	1.84	0.60
1:D:58:GLY:C	1:D:119:VAL:HG22	2.22	0.60
1:B:94:GLU:HG2	1:B:102:HIS:O	2.01	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:7:LEU:HD23	1:C:54:ASN:HD22	1.66	0.60
1:D:193:ARG:NH2	1:D:230:LEU:HB3	2.17	0.60
1:B:150:ASP:OD1	3:B:1353:NAP:H4N	2.01	0.59
1:B:29:VAL:HB	1:B:124:PHE:CE1	2.37	0.59
1:B:217:ASN:HD22	1:B:217:ASN:C	2.05	0.59
1:C:49:ILE:HG22	1:C:52:ARG:NH2	2.12	0.59
1:A:1:MET:HG3	1:A:122:GLU:HB2	1.84	0.59
1:C:91:ARG:NH1	1:C:91:ARG:HB3	2.17	0.59
1:C:157:HIS:CE1	1:D:287:HIS:CE1	2.85	0.59
1:D:29:VAL:HB	1:D:124:PHE:CE1	2.37	0.59
1:D:8:SER:OG	1:D:11:LYS:HD2	2.03	0.59
1:C:32:LEU:HG	1:C:63:GLY:HA2	1.85	0.58
1:B:200:ARG:NH1	1:B:344:LEU:O	2.36	0.58
1:C:24:PRO:O	1:C:71:GLU:HB2	2.04	0.58
1:C:9:ILE:HG23	1:C:51:GLU:HG3	1.85	0.58
1:D:49:ILE:HG12	1:D:267:TYR:HH	1.67	0.58
1:D:264:ASN:O	1:D:292:GLY:HA2	2.04	0.58
1:D:344:LEU:HD11	1:D:347:PRO:HD3	1.86	0.57
1:B:327:PHE:O	1:B:330:ILE:HD13	2.04	0.57
1:D:4:PHE:O	1:D:56:ILE:HA	2.05	0.57
1:C:65:VAL:O	1:C:77:PRO:HA	2.05	0.57
1:C:91:ARG:HH11	1:C:91:ARG:HB3	1.69	0.57
1:D:126:VAL:HG11	1:D:133:LEU:HD21	1.87	0.57
1:D:146:VAL:HG13	1:D:147:MET:CE	2.34	0.56
1:D:267:TYR:HB3	3:D:3353:NAP:C2D	2.36	0.56
1:C:286:ALA:O	1:C:287:HIS:HB2	2.06	0.56
1:D:12:VAL:HG12	1:D:13:GLY:H	1.69	0.56
1:D:301:ARG:O	1:D:305:LEU:HG	2.06	0.56
1:C:254:LYS:HG2	1:C:280:GLU:HG2	1.86	0.56
1:A:36:PRO:HG2	1:A:337:MET:HE2	1.86	0.56
1:B:282:GLY:HA3	1:B:286:ALA:HB2	1.87	0.56
1:C:33:ALA:HB3	1:C:62:VAL:HG22	1.87	0.56
1:A:256:VAL:HG21	1:A:262:ILE:CD1	2.34	0.56
1:A:149:PRO:HA	1:A:302:MET:CE	2.36	0.56
1:B:102:HIS:HD2	1:B:107:LEU:N	1.89	0.56
1:C:285:MET:HB2	1:D:107:LEU:HD21	1.86	0.56
1:A:86:ILE:CD1	1:A:110:TRP:HB2	2.36	0.56
1:A:42:HIS:O	1:A:46:GLU:HB2	2.05	0.56
1:B:137:PRO:HB2	1:B:140:ILE:HD13	1.87	0.56
1:C:94:GLU:HB2	1:C:103:SER:HA	1.88	0.56
1:C:49:ILE:O	1:C:49:ILE:HG22	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:94:GLU:HB2	1:D:103:SER:HA	1.87	0.55
1:C:245:ASN:O	1:C:248:ILE:HG22	2.07	0.55
1:A:12:VAL:HG12	1:A:13:GLY:N	2.21	0.55
1:D:325:ARG:O	1:D:329:ASN:ND2	2.37	0.55
1:D:341:PRO:HG2	1:D:344:LEU:HB2	1.87	0.55
1:C:102:HIS:HD2	1:C:107:LEU:H	1.54	0.55
1:C:280:GLU:N	1:C:280:GLU:OE1	2.39	0.55
1:D:6:MET:SD	1:D:44:VAL:HG22	2.46	0.55
1:A:337:MET:CE	1:A:347:PRO:HD2	2.37	0.55
1:D:102:HIS:HD2	1:D:107:LEU:H	1.52	0.55
1:A:116:LYS:NZ	1:A:125:HIS:HD2	2.05	0.55
1:C:39:SER:O	1:C:43:THR:HG23	2.07	0.55
1:C:202:VAL:HG21	1:C:343:ASP:HA	1.89	0.55
1:D:146:VAL:HG13	1:D:147:MET:HE3	1.87	0.55
1:B:24:PRO:O	1:B:71:GLU:HB2	2.06	0.55
1:B:300:LEU:HD22	1:D:97:ARG:HG2	1.88	0.55
1:A:274:LEU:HB2	1:B:276:VAL:HB	1.88	0.55
1:D:161:LEU:O	1:D:291:LYS:HD2	2.07	0.55
1:B:156:PHE:O	1:B:160:GLU:HG3	2.07	0.54
1:A:294:LEU:HA	1:B:285:MET:HE2	1.90	0.54
1:C:34:VAL:HG12	1:C:61:ALA:HB2	1.88	0.54
1:D:20:PRO:HG2	1:D:28:ILE:HD12	1.90	0.54
1:B:264:ASN:O	1:B:292:GLY:HA2	2.08	0.54
1:C:175:ILE:HD11	1:C:198:GLY:HA3	1.88	0.54
1:C:5:ALA:HB2	1:C:56:ILE:HA	1.88	0.54
1:C:37:CYS:HB2	1:C:60:GLU:OE2	2.08	0.54
1:B:131:MET:HG3	1:D:91:ARG:HH11	1.73	0.54
1:C:31:PRO:O	1:C:351:LEU:HD12	2.07	0.54
1:B:139:GLU:N	1:B:139:GLU:OE1	2.37	0.54
1:A:324:PHE:HB3	1:A:329:ASN:ND2	2.22	0.54
1:C:116:LYS:NZ	1:C:125:HIS:HD2	2.06	0.54
1:B:223:ILE:HG22	1:B:227:ILE:HD12	1.90	0.53
1:C:36:PRO:HG2	1:C:337:MET:HE2	1.90	0.53
1:B:66:VAL:HG21	1:B:123:PHE:HE1	1.73	0.53
1:B:6:MET:HB2	1:B:57:LEU:HD21	1.90	0.53
1:B:256:VAL:HG11	1:B:262:ILE:HD13	1.90	0.53
1:B:1:MET:HB3	1:B:17:LYS:O	2.08	0.53
1:B:193:ARG:HB2	1:C:311:TYR:CE2	2.43	0.53
1:B:89:ASP:OD1	1:B:89:ASP:N	2.33	0.53
1:D:321:THR:HG23	1:D:345:ILE:HB	1.91	0.53
1:B:82:VAL:HG12	1:B:302:MET:CE	2.38	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:8:SER:O	1:A:9:ILE:C	2.47	0.53
1:D:344:LEU:HG	1:D:345:ILE:N	2.22	0.53
1:B:188:LEU:HD11	1:B:314:VAL:CG2	2.39	0.53
1:C:171:ALA:HB2	1:C:236:VAL:HG11	1.90	0.53
1:A:49:ILE:O	1:A:49:ILE:HG22	2.09	0.53
1:A:102:HIS:HE1	1:B:287:HIS:HD2	1.56	0.53
1:A:152:MET:HG2	1:A:305:LEU:HD13	1.91	0.53
1:C:8:SER:OG	1:C:11:LYS:HD2	2.09	0.52
1:A:339:ASP:O	1:A:341:PRO:HD3	2.09	0.52
1:A:82:VAL:HG12	1:A:302:MET:CE	2.39	0.52
1:B:281:TRP:CE3	1:B:284:GLY:HA2	2.45	0.52
1:A:267:TYR:CB	3:A:353:NAP:O2D	2.57	0.52
1:C:293:GLY:HA2	1:D:285:MET:HA	1.91	0.52
1:C:80:ARG:HG2	1:C:80:ARG:NH1	2.23	0.52
1:C:102:HIS:CD2	1:C:107:LEU:HD23	2.45	0.52
1:C:7:LEU:CD2	1:C:54:ASN:HD22	2.22	0.52
1:B:322:HIS:CD2	1:B:336:LEU:HD21	2.45	0.52
1:A:157:HIS:HE1	1:B:287:HIS:HE1	1.58	0.52
1:C:172:VAL:HG21	1:C:183:VAL:HG22	1.90	0.52
1:C:293:GLY:HA2	1:D:285:MET:C	2.30	0.52
1:C:250:ALA:O	1:C:254:LYS:HG3	2.08	0.52
1:C:294:LEU:N	1:D:285:MET:HB3	2.25	0.51
1:D:191:ALA:HB3	1:D:194:ILE:HD11	1.91	0.51
1:C:106:MET:HG2	1:C:107:LEU:HG	1.93	0.51
1:A:302:MET:O	1:A:306:ILE:HG13	2.11	0.51
1:D:327:PHE:O	1:D:330:ILE:HG12	2.09	0.51
1:D:218:TYR:O	1:D:219:LYS:HG3	2.10	0.51
1:D:281:TRP:CZ3	1:D:284:GLY:HA2	2.44	0.51
1:B:175:ILE:HG21	1:B:196:ALA:HB1	1.92	0.51
1:A:102:HIS:HE1	1:B:287:HIS:CD2	2.28	0.51
1:C:51:GLU:HG2	1:C:52:ARG:N	2.25	0.51
1:C:49:ILE:CG2	1:C:52:ARG:HH22	2.17	0.51
1:D:59:HIS:CE1	1:D:110:TRP:HE1	2.27	0.51
1:A:217:ASN:HB3	1:A:220:ASP:OD1	2.11	0.51
1:C:102:HIS:CD2	1:C:107:LEU:H	2.29	0.51
1:C:256:VAL:CG2	1:C:288:LYS:HD3	2.41	0.51
1:D:51:GLU:HG2	1:D:52:ARG:N	2.25	0.51
1:D:183:VAL:HG13	1:D:194:ILE:HG21	1.93	0.51
1:B:256:VAL:HG11	1:B:262:ILE:CD1	2.40	0.50
1:B:280:GLU:OE1	1:B:280:GLU:N	2.44	0.50
1:C:339:ASP:O	1:C:341:PRO:HD3	2.10	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:6:MET:HG3	1:A:44:VAL:HG22	1.93	0.50
1:D:1:MET:HB3	1:D:17:LYS:O	2.12	0.50
1:D:97:ARG:HB3	1:D:97:ARG:HH11	1.76	0.50
1:C:217:ASN:C	1:C:219:LYS:H	2.15	0.50
1:D:59:HIS:HE1	1:D:110:TRP:HZ2	1.60	0.50
1:C:267:TYR:H	3:C:2353:NAP:H1D	1.75	0.50
1:C:199:SER:HB3	1:C:217:ASN:ND2	2.27	0.50
1:C:188:LEU:HD11	1:C:314:VAL:CG2	2.42	0.50
1:A:191:ALA:HB3	1:A:194:ILE:HD11	1.93	0.50
1:B:49:ILE:HG12	1:B:267:TYR:OH	2.12	0.50
1:D:302:MET:O	1:D:306:ILE:HD13	2.12	0.50
1:C:83:VAL:O	1:C:149:PRO:HB3	2.12	0.50
1:B:43:THR:HG22	1:B:267:TYR:OH	2.12	0.49
1:D:82:VAL:HG12	1:D:302:MET:CE	2.41	0.49
1:C:287:HIS:HE1	1:D:157:HIS:CE1	2.30	0.49
1:A:82:VAL:HG23	1:A:136:LEU:HB2	1.94	0.49
1:D:256:VAL:HG21	1:D:262:ILE:CD1	2.42	0.49
1:B:9:ILE:HD12	1:B:47:GLY:CA	2.42	0.49
1:C:20:PRO:O	1:C:28:ILE:HD13	2.12	0.49
1:B:281:TRP:CZ3	1:B:284:GLY:HA2	2.47	0.49
1:C:149:PRO:HA	1:C:302:MET:HE2	1.95	0.49
1:B:6:MET:CE	1:B:9:ILE:HD13	2.43	0.49
1:C:224:GLU:OE1	1:C:254:LYS:HD3	2.12	0.49
1:D:321:THR:C	1:D:322:HIS:ND1	2.66	0.49
1:C:252:ALA:HB1	1:C:262:ILE:HD12	1.94	0.49
1:A:122:GLU:HA	1:A:327:PHE:CZ	2.48	0.49
1:C:264:ASN:O	1:C:292:GLY:HA2	2.12	0.49
1:A:299:ARG:O	1:A:303:GLU:HG3	2.13	0.49
1:C:143:GLU:OE1	1:C:350:ILE:HD11	2.12	0.49
1:C:224:GLU:O	1:C:228:MET:HG2	2.11	0.49
1:D:153:THR:HG23	1:D:154:THR:N	2.28	0.49
1:A:86:ILE:HD12	1:A:110:TRP:HB2	1.94	0.48
1:C:173:LEU:HD23	1:C:197:VAL:HG21	1.94	0.48
1:C:287:HIS:CD2	1:D:102:HIS:CE1	2.92	0.48
1:D:341:PRO:HG2	1:D:344:LEU:CB	2.43	0.48
1:D:12:VAL:CG1	1:D:13:GLY:N	2.76	0.48
1:A:177:PRO:HD2	3:A:353:NAP:O1N	2.13	0.48
1:C:293:GLY:HA3	1:D:287:HIS:CE1	2.49	0.48
1:A:337:MET:HE1	1:A:347:PRO:HD2	1.96	0.48
1:B:91:ARG:HH11	1:D:128:ASP:CG	2.16	0.48
1:B:91:ARG:HD2	1:D:128:ASP:OD2	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:52:ARG:HD2	1:D:55:MET:SD	2.53	0.48
1:D:334:PHE:O	1:D:337:MET:HB3	2.14	0.48
1:B:254:LYS:HG2	1:B:280:GLU:CG	2.41	0.48
1:C:183:VAL:HG13	1:C:194:ILE:HG21	1.95	0.48
1:B:39:SER:O	1:B:43:THR:HG23	2.13	0.48
1:C:21:ALA:HA	1:C:28:ILE:HD11	1.96	0.48
1:D:280:GLU:OE1	1:D:280:GLU:N	2.44	0.48
1:A:223:ILE:HG22	1:A:227:ILE:HD12	1.95	0.48
1:C:293:GLY:CA	1:D:285:MET:HA	2.43	0.48
1:D:148:ILE:HB	1:D:149:PRO:HD3	1.96	0.48
1:A:36:PRO:HG2	1:A:337:MET:HE3	1.96	0.47
1:B:175:ILE:N	3:B:1353:NAP:O3B	2.44	0.47
1:B:136:LEU:HD12	1:B:137:PRO:HD2	1.95	0.47
1:C:75:PHE:O	1:C:76:LYS:HD3	2.14	0.47
1:B:82:VAL:HG21	1:B:306:ILE:HD11	1.95	0.47
1:B:49:ILE:O	1:B:49:ILE:CG2	2.63	0.47
1:A:300:LEU:HD21	1:C:97:ARG:HG2	1.96	0.47
1:C:143:GLU:CD	1:C:143:GLU:H	2.17	0.47
1:C:195:ILE:HD12	1:C:195:ILE:N	2.29	0.47
1:D:75:PHE:CE1	1:D:135:HIS:CE1	3.02	0.47
1:B:36:PRO:HG2	1:B:337:MET:HE2	1.96	0.47
1:A:287:HIS:HE1	1:B:157:HIS:CE1	2.19	0.47
1:D:36:PRO:HG2	1:D:337:MET:HE3	1.97	0.47
1:A:157:HIS:HA	1:A:160:GLU:OE1	2.14	0.47
1:D:169:THR:HB	1:D:193:ARG:HB3	1.97	0.47
1:C:116:LYS:HZ2	1:C:125:HIS:HD2	1.63	0.47
1:A:268:PHE:O	1:B:278:ARG:HD2	2.14	0.47
1:D:34:VAL:HG12	1:D:61:ALA:CB	2.44	0.47
1:B:68:VAL:HB	1:B:72:VAL:HB	1.96	0.47
1:D:200:ARG:NH1	1:D:203:CYS:SG	2.88	0.47
1:C:299:ARG:O	1:C:303:GLU:HG3	2.15	0.47
1:D:100:HIS:CD2	1:D:101:GLN:N	2.83	0.47
1:D:43:THR:HG22	1:D:49:ILE:HG12	1.97	0.47
1:A:68:VAL:HB	1:A:72:VAL:HG21	1.97	0.47
1:B:83:VAL:HG22	1:B:133:LEU:CD2	2.45	0.47
1:A:51:GLU:HG2	1:A:52:ARG:N	2.30	0.47
1:D:147:MET:O	1:D:151:MET:HB2	2.15	0.47
1:D:193:ARG:HH22	1:D:230:LEU:HB3	1.81	0.46
1:C:92:THR:OG1	1:C:95:VAL:HG23	2.15	0.46
1:C:287:HIS:CE1	1:D:157:HIS:CE1	2.99	0.46
1:D:335:MET:O	1:D:338:LYS:N	2.49	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:177:PRO:CD	3:C:2353:NAP:O1N	2.60	0.46
1:D:6:MET:SD	1:D:44:VAL:HA	2.55	0.46
1:D:254:LYS:HG2	1:D:280:GLU:CG	2.46	0.46
1:D:58:GLY:O	1:D:119:VAL:HG22	2.15	0.46
1:D:148:ILE:HG22	1:D:302:MET:HE1	1.96	0.46
1:D:240:ILE:HA	1:D:263:ALA:O	2.16	0.46
1:D:200:ARG:HG3	3:D:3353:NAP:O3X	2.14	0.46
1:B:6:MET:SD	1:B:52:ARG:HB2	2.56	0.46
1:B:325:ARG:O	1:B:329:ASN:ND2	2.49	0.46
1:C:9:ILE:H	1:C:9:ILE:HD13	1.81	0.46
1:B:155:GLY:O	1:B:182:ALA:HA	2.16	0.46
1:A:217:ASN:C	1:A:217:ASN:ND2	2.67	0.46
1:A:245:ASN:O	1:A:248:ILE:HG22	2.15	0.46
1:A:157:HIS:CE1	1:B:287:HIS:HE1	2.34	0.46
1:B:273:VAL:O	1:B:274:LEU:HD23	2.16	0.46
1:B:152:MET:CG	1:B:305:LEU:HD13	2.42	0.45
1:C:341:PRO:HG2	1:C:344:LEU:HB3	1.98	0.45
1:B:311:TYR:CE2	1:C:193:ARG:HB2	2.50	0.45
1:A:37:CYS:HB2	1:A:60:GLU:OE2	2.16	0.45
1:D:113:SER:OG	1:D:113:SER:O	2.32	0.45
1:A:311:TYR:CE2	1:D:193:ARG:HB2	2.50	0.45
1:C:82:VAL:HG12	1:C:149:PRO:HG3	1.98	0.45
1:A:86:ILE:HD12	1:A:110:TRP:CA	2.46	0.45
1:C:322:HIS:CD2	1:C:336:LEU:HD21	2.51	0.45
1:A:337:MET:HE2	1:A:347:PRO:CD	2.47	0.45
1:A:96:GLN:HG2	1:C:90:TRP:CZ3	2.52	0.45
1:C:293:GLY:HA2	1:D:285:MET:O	2.17	0.45
1:D:71:GLU:O	1:D:73:LYS:HG3	2.16	0.45
1:C:9:ILE:HD13	1:C:9:ILE:N	2.31	0.45
1:D:83:VAL:HG22	1:D:133:LEU:CD2	2.47	0.45
1:D:24:PRO:O	1:D:71:GLU:HB2	2.17	0.45
1:B:181:MET:CE	1:B:181:MET:HA	2.47	0.45
1:C:49:ILE:HG13	1:C:267:TYR:OH	2.17	0.45
1:C:101:GLN:HG2	1:C:294:LEU:O	2.17	0.45
1:C:154:THR:HG21	3:C:2353:NAP:C5N	2.47	0.45
1:C:66:VAL:HG21	1:C:123:PHE:CE1	2.46	0.45
1:B:321:THR:C	1:B:322:HIS:ND1	2.70	0.45
1:D:106:MET:O	1:D:107:LEU:HB2	2.16	0.44
1:D:254:LYS:HE2	1:D:280:GLU:OE2	2.16	0.44
1:A:10:GLY:N	1:A:44:VAL:O	2.47	0.44
1:B:116:LYS:HZ1	1:B:125:HIS:HD2	1.65	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:206:ALA:O	1:D:209:TYR:HB3	2.17	0.44
1:D:107:LEU:HD22	1:D:294:LEU:HD13	1.99	0.44
1:A:181:MET:CE	1:A:181:MET:HA	2.47	0.44
1:B:34:VAL:HG12	1:B:61:ALA:HB2	1.99	0.44
1:A:102:HIS:HD2	1:A:107:LEU:N	2.10	0.44
1:B:40:ASP:OD2	1:B:57:LEU:HB3	2.18	0.44
1:C:282:GLY:HA3	1:C:286:ALA:HB2	1.99	0.44
1:B:86:ILE:HG22	1:B:88:PRO:HD3	1.99	0.44
1:D:175:ILE:N	3:D:3353:NAP:O3B	2.38	0.44
1:A:327:PHE:HD1	1:A:351:LEU:HD22	1.82	0.44
1:A:5:ALA:HB2	1:A:56:ILE:HA	2.00	0.44
1:D:306:ILE:HG22	1:D:307:ASP:N	2.33	0.44
1:A:197:VAL:HG22	1:A:216:VAL:HB	1.99	0.44
1:C:142:LEU:HA	1:C:142:LEU:HD23	1.81	0.44
1:D:15:ILE:HD13	1:D:54:ASN:HB3	1.99	0.44
1:D:49:ILE:CG1	1:D:267:TYR:OH	2.63	0.44
1:D:107:LEU:CD2	1:D:294:LEU:HD13	2.48	0.44
1:C:149:PRO:HA	1:C:302:MET:CE	2.48	0.44
1:A:22:PRO:HD3	1:A:28:ILE:CG1	2.47	0.44
1:D:14:TRP:HZ3	1:D:327:PHE:CE2	2.32	0.43
1:C:36:PRO:CG	1:C:347:PRO:HG2	2.47	0.43
1:C:327:PHE:HA	1:C:351:LEU:HD22	1.99	0.43
1:D:34:VAL:HG12	1:D:61:ALA:HB2	2.00	0.43
1:A:224:GLU:O	1:A:228:MET:HG2	2.18	0.43
1:C:150:ASP:OD1	3:C:2353:NAP:H4N	2.17	0.43
1:D:335:MET:C	1:D:337:MET:N	2.70	0.43
1:A:101:GLN:HB3	1:A:294:LEU:HD23	2.00	0.43
1:D:231:THR:O	1:D:234:LYS:HG3	2.18	0.43
1:C:4:PHE:CZ	1:C:12:VAL:CG1	3.01	0.43
1:A:76:LYS:HE3	1:A:79:ASP:OD1	2.18	0.43
1:B:256:VAL:CG1	1:B:262:ILE:HD11	2.48	0.43
1:A:337:MET:HE2	1:A:347:PRO:HD2	2.00	0.43
1:A:94:GLU:HB2	1:A:103:SER:HA	2.00	0.43
1:D:256:VAL:CG2	1:D:262:ILE:HD11	2.48	0.43
1:A:327:PHE:CD1	1:A:351:LEU:HD22	2.54	0.43
1:D:82:VAL:HG21	1:D:306:ILE:HD11	2.00	0.43
1:A:190:GLY:HA2	1:D:188:LEU:HD22	2.01	0.43
1:A:170:VAL:HG12	1:A:171:ALA:N	2.32	0.43
1:B:241:ILE:HD11	1:B:252:ALA:CB	2.48	0.43
1:B:32:LEU:CG	1:B:63:GLY:HA2	2.45	0.43
1:B:59:HIS:HB3	1:B:112:PHE:HB2	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:150:ASP:CG	3:A:353:NAP:H5N	2.39	0.43
1:B:321:THR:O	1:B:322:HIS:ND1	2.52	0.43
1:B:91:ARG:NH1	1:D:128:ASP:OD1	2.52	0.43
1:C:287:HIS:CE1	1:D:293:GLY:HA3	2.54	0.43
1:C:174:GLY:O	1:C:179:GLY:HA3	2.19	0.43
1:B:116:LYS:NZ	1:B:125:HIS:CD2	2.85	0.43
1:A:5:ALA:O	1:A:12:VAL:HG13	2.19	0.43
1:B:164:ILE:HD12	1:B:170:VAL:HG22	2.00	0.43
1:D:337:MET:CE	1:D:347:PRO:HD2	2.48	0.42
1:C:268:PHE:O	1:D:278:ARG:NH1	2.52	0.42
1:C:126:VAL:HG11	1:C:133:LEU:HD21	2.01	0.42
1:B:154:THR:OG1	1:B:295:CYS:HB3	2.19	0.42
1:D:178:VAL:CB	3:D:3353:NAP:H52N	2.47	0.42
1:D:41:ILE:HD11	1:D:337:MET:HG3	2.01	0.42
1:D:30:ARG:HA	1:D:31:PRO:HD3	1.89	0.42
1:A:206:ALA:O	1:A:209:TYR:HB3	2.19	0.42
1:A:217:ASN:ND2	1:A:219:LYS:N	2.51	0.42
1:D:9:ILE:HG23	1:D:51:GLU:CG	2.46	0.42
1:B:214:ASP:C	1:B:215:ILE:HD12	2.39	0.42
1:D:209:TYR:CZ	1:D:318:LYS:HG3	2.54	0.42
1:B:20:PRO:O	1:B:28:ILE:HD13	2.19	0.42
1:C:34:VAL:HG12	1:C:61:ALA:CB	2.50	0.42
1:C:188:LEU:HA	1:C:188:LEU:HD23	1.89	0.42
1:D:222:PRO:O	1:D:226:GLN:HG3	2.19	0.42
1:B:5:ALA:HB2	1:B:56:ILE:HA	2.01	0.42
1:C:295:CYS:HA	1:C:296:PRO:HD3	1.91	0.42
1:B:245:ASN:HB2	1:B:247:ASP:OD1	2.19	0.42
1:D:52:ARG:HH11	1:D:52:ARG:HG3	1.84	0.42
1:B:188:LEU:O	1:C:190:GLY:N	2.53	0.42
1:C:228:MET:HE3	1:C:233:GLY:HA2	2.00	0.42
1:C:58:GLY:C	1:C:119:VAL:HG22	2.39	0.42
1:B:218:TYR:CB	1:B:223:ILE:HD11	2.38	0.42
1:B:90:TRP:CZ3	1:D:96:GLN:HG2	2.53	0.42
1:B:30:ARG:HA	1:B:31:PRO:HD3	1.85	0.42
1:D:135:HIS:O	1:D:137:PRO:HD3	2.20	0.42
1:A:336:LEU:HD11	1:A:341:PRO:HG3	2.00	0.42
1:B:176:GLY:O	1:B:180:LEU:HG	2.20	0.42
1:B:44:VAL:HG12	1:B:45:PHE:N	2.34	0.42
1:A:49:ILE:HG12	1:A:267:TYR:OH	2.20	0.42
1:A:285:MET:HB2	1:B:107:LEU:HD21	2.01	0.42
1:D:199:SER:O	1:D:200:ARG:C	2.58	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:148:ILE:N	1:A:149:PRO:CD	2.83	0.42
1:D:204:VAL:HG12	1:D:208:LYS:HE3	2.01	0.42
1:A:252:ALA:O	1:A:256:VAL:HG13	2.20	0.42
1:B:329:ASN:O	1:B:330:ILE:C	2.59	0.42
1:D:302:MET:O	1:D:306:ILE:CD1	2.68	0.42
1:B:82:VAL:CG1	1:B:302:MET:CE	2.97	0.42
1:D:217:ASN:HD22	1:D:217:ASN:C	2.22	0.42
1:A:191:ALA:CB	1:A:194:ILE:HD11	2.50	0.42
1:B:4:PHE:CE2	1:B:12:VAL:HG11	2.54	0.42
1:A:91:ARG:NH1	1:C:128:ASP:OD1	2.53	0.42
1:A:49:ILE:CG2	1:A:52:ARG:HH22	2.13	0.41
1:D:322:HIS:O	1:D:347:PRO:HA	2.20	0.41
1:B:31:PRO:HB2	1:B:351:LEU:CD1	2.42	0.41
1:D:128:ASP:OD1	1:D:131:MET:HB2	2.20	0.41
1:C:20:PRO:O	1:C:28:ILE:CD1	2.68	0.41
1:D:227:ILE:HA	1:D:230:LEU:HD12	2.01	0.41
1:C:132:ASN:O	1:C:133:LEU:HD23	2.20	0.41
1:A:175:ILE:HA	1:A:179:GLY:HA3	2.03	0.41
1:D:75:PHE:CE2	1:D:133:LEU:HB2	2.56	0.41
1:D:1:MET:HG3	1:D:122:GLU:HB2	2.02	0.41
1:A:177:PRO:HG2	3:A:353:NAP:O1N	2.20	0.41
1:D:202:VAL:HG12	1:D:345:ILE:HD11	2.03	0.41
1:A:6:MET:HB2	1:A:57:LEU:HD21	2.03	0.41
1:C:148:ILE:HB	1:C:149:PRO:HD3	2.02	0.41
1:A:25:PHE:CD1	1:C:91:ARG:NH2	2.88	0.41
1:B:128:ASP:OD1	1:D:91:ARG:NH1	2.54	0.41
1:B:4:PHE:CZ	1:B:12:VAL:HG12	2.55	0.41
1:C:259:GLY:HA2	1:C:287:HIS:HB3	2.03	0.41
1:A:266:ASN:OD1	3:A:353:NAP:O3D	2.25	0.41
1:A:146:VAL:HG13	1:A:147:MET:CE	2.49	0.41
1:D:308:LEU:HA	1:D:308:LEU:HD23	1.93	0.41
1:D:6:MET:CE	1:D:52:ARG:HB2	2.51	0.41
1:D:175:ILE:HG21	1:D:196:ALA:HB1	2.03	0.41
1:C:92:THR:HG21	1:C:103:SER:OG	2.21	0.41
1:D:337:MET:HE1	1:D:346:LYS:HD2	2.02	0.41
1:B:344:LEU:HD11	1:B:347:PRO:HD3	2.03	0.41
1:C:264:ASN:OD1	1:C:266:ASN:HB3	2.21	0.41
1:A:53:HIS:O	1:A:54:ASN:HB2	2.21	0.41
1:B:339:ASP:O	1:B:341:PRO:HD3	2.21	0.41
1:D:314:VAL:HG12	1:D:315:ASP:N	2.36	0.41
1:D:337:MET:HE1	1:D:347:PRO:HD2	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:300:LEU:CD2	1:C:97:ARG:HG2	2.50	0.41
1:C:169:THR:HA	1:C:193:ARG:O	2.21	0.41
1:D:164:ILE:HD12	1:D:170:VAL:CG2	2.51	0.41
1:A:97:ARG:HG2	1:C:300:LEU:CD2	2.50	0.41
1:D:6:MET:HE2	1:D:52:ARG:HB2	2.03	0.40
1:B:90:TRP:CH2	1:D:96:GLN:HG2	2.56	0.40
1:C:88:PRO:HG3	1:C:100:HIS:HB2	2.04	0.40
1:D:12:VAL:CG1	1:D:13:GLY:H	2.31	0.40
1:C:84:PRO:HD3	1:C:302:MET:HG3	2.03	0.40
1:B:319:LEU:O	1:B:345:ILE:HB	2.21	0.40
1:B:152:MET:CE	1:B:316:PRO:HB3	2.51	0.40
1:D:32:LEU:HD23	1:D:32:LEU:HA	1.86	0.40
1:B:174:GLY:O	1:B:179:GLY:HA3	2.22	0.40
1:D:295:CYS:HA	1:D:296:PRO:HD3	1.90	0.40
1:A:254:LYS:HG2	1:A:280:GLU:HG2	2.02	0.40
1:B:147:MET:CB	1:B:316:PRO:HB2	2.52	0.40
1:A:133:LEU:HA	1:A:133:LEU:HD23	1.91	0.40
1:A:49:ILE:O	1:A:52:ARG:NH2	2.55	0.40
1:A:256:VAL:O	1:A:288:LYS:HE2	2.20	0.40
1:A:294:LEU:HD12	1:B:285:MET:HE1	2.03	0.40
1:C:86:ILE:HD11	1:C:110:TRP:HB2	2.04	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	350/352 (99%)	316 (90%)	32 (9%)	2 (1%)	30	65
1	B	350/352 (99%)	320 (91%)	27 (8%)	3 (1%)	21	55
1	C	350/352 (99%)	325 (93%)	24 (7%)	1 (0%)	46	79
1	D	350/352 (99%)	310 (89%)	33 (9%)	7 (2%)	9	30

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	1400/1408 (99%)	1271 (91%)	116 (8%)	13 (1%)	21	55

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	9	ILE
1	A	46	GLU
1	D	46	GLU
1	D	326	GLY
1	D	114	ASN
1	D	328	ASP
1	D	329	ASN
1	D	338	LYS
1	B	245	ASN
1	C	218	TYR
1	B	44	VAL
1	D	115	VAL
1	B	347	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	274/274 (100%)	265 (97%)	9 (3%)	45	79
1	B	274/274 (100%)	264 (96%)	10 (4%)	42	76
1	C	274/274 (100%)	266 (97%)	8 (3%)	50	83
1	D	274/274 (100%)	264 (96%)	10 (4%)	42	76
All	All	1096/1096 (100%)	1059 (97%)	37 (3%)	44	78

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	43	THR
1	A	64	GLU

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Mol	Chain	Res	Type
1	A	169	THR
1	A	217	ASN
1	A	245	ASN
1	A	307	ASP
1	A	325	ARG
1	A	327	PHE
1	A	339	ASP
1	B	62	VAL
1	B	89	ASP
1	B	169	THR
1	B	217	ASN
1	B	220	ASP
1	B	256	VAL
1	B	267	TYR
1	B	325	ARG
1	B	339	ASP
1	B	344	LEU
1	C	9	ILE
1	C	175	ILE
1	C	214	ASP
1	C	217	ASN
1	C	220	ASP
1	C	328	ASP
1	C	339	ASP
1	C	344	LEU
1	D	26	ASP
1	D	46	GLU
1	D	89	ASP
1	D	169	THR
1	D	217	ASN
1	D	229	ASN
1	D	245	ASN
1	D	307	ASP
1	D	325	ARG
1	D	339	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	102	HIS
1	A	125	HIS
1	A	157	HIS

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Mol	Chain	Res	Type
1	A	217	ASN
1	A	287	HIS
1	B	102	HIS
1	B	125	HIS
1	B	157	HIS
1	B	217	ASN
1	B	287	HIS
1	C	42	HIS
1	C	54	ASN
1	C	102	HIS
1	C	125	HIS
1	C	157	HIS
1	C	217	ASN
1	C	245	ASN
1	C	287	HIS
1	D	102	HIS
1	D	125	HIS
1	D	157	HIS
1	D	217	ASN
1	D	245	ASN
1	D	287	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 for Mogul analysis.

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	NAP	A	353	-	42,52,52	1.28	5 (11%)	54,80,80	1.87	14 (25%)
3	NAP	B	1353	-	42,52,52	1.30	6 (14%)	54,80,80	1.95	15 (27%)
3	NAP	C	2353	-	42,52,52	1.73	4 (9%)	54,80,80	2.04	14 (25%)
3	NAP	D	3353	-	42,52,52	1.42	7 (16%)	54,80,80	1.60	8 (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	NAP	A	353	-	-	0/27/67/67	0/5/5/5
3	NAP	B	1353	-	-	0/27/67/67	0/5/5/5
3	NAP	C	2353	-	-	0/27/67/67	0/5/5/5
3	NAP	D	3353	-	-	0/27/67/67	0/5/5/5

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1353	NAP	C3N-C7N	-2.50	1.46	1.50
3	A	353	NAP	C5A-C4A	-2.40	1.35	1.40
3	B	1353	NAP	PA-O2A	-2.19	1.45	1.54
3	C	2353	NAP	C5A-C4A	-2.12	1.35	1.40
3	D	3353	NAP	PA-O2A	-2.09	1.46	1.54
3	D	3353	NAP	C5A-C4A	-2.05	1.35	1.40
3	B	1353	NAP	C5A-C4A	-2.05	1.35	1.40
3	A	353	NAP	P2B-O2X	-2.03	1.47	1.54
3	A	353	NAP	P2B-O3X	-2.01	1.47	1.54
3	B	1353	NAP	C6N-N1N	2.05	1.41	1.35
3	D	3353	NAP	C6N-N1N	2.26	1.41	1.35
3	D	3353	NAP	O4D-C1D	2.59	1.44	1.41
3	C	2353	NAP	C4A-N3A	2.74	1.39	1.35
3	D	3353	NAP	C4A-N3A	2.75	1.39	1.35
3	B	1353	NAP	C4A-N3A	2.80	1.39	1.35
3	A	353	NAP	C2N-C3N	3.08	1.43	1.39
3	D	3353	NAP	O4B-C1B	3.14	1.45	1.41
3	A	353	NAP	C4A-N3A	3.19	1.40	1.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1353	NAP	C2N-C3N	3.89	1.44	1.39
3	D	3353	NAP	C2N-C3N	4.17	1.45	1.39
3	C	2353	NAP	C2N-C3N	4.33	1.45	1.39
3	C	2353	NAP	O4D-C1D	6.99	1.50	1.41

All (51) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1353	NAP	N3A-C2A-N1A	-5.78	124.47	128.89
3	C	2353	NAP	N3A-C2A-N1A	-5.60	124.61	128.89
3	D	3353	NAP	N3A-C2A-N1A	-5.49	124.69	128.89
3	A	353	NAP	N3A-C2A-N1A	-5.20	124.91	128.89
3	B	1353	NAP	C4B-O4B-C1B	-4.95	104.28	109.72
3	C	2353	NAP	O7N-C7N-N7N	-4.32	116.51	122.59
3	D	3353	NAP	O7N-C7N-N7N	-4.14	116.76	122.59
3	A	353	NAP	O7N-C7N-N7N	-4.10	116.83	122.59
3	A	353	NAP	C4B-O4B-C1B	-4.04	105.28	109.72
3	B	1353	NAP	O7N-C7N-N7N	-3.93	117.06	122.59
3	C	2353	NAP	O4D-C4D-C3D	-3.88	97.32	105.15
3	A	353	NAP	C3B-C2B-C1B	-3.54	95.87	102.73
3	B	1353	NAP	C1B-N9A-C4A	-3.46	121.72	126.94
3	B	1353	NAP	O4B-C1B-C2B	-3.31	100.61	106.60
3	B	1353	NAP	C3B-C2B-C1B	-3.21	96.52	102.73
3	C	2353	NAP	C3N-C2N-N1N	-3.16	116.73	120.36
3	D	3353	NAP	O4B-C1B-C2B	-3.04	101.11	106.60
3	A	353	NAP	C3N-C2N-N1N	-2.99	116.91	120.36
3	C	2353	NAP	O4B-C1B-C2B	-2.64	101.83	106.60
3	A	353	NAP	C1B-N9A-C4A	-2.62	122.99	126.94
3	D	3353	NAP	C4B-O4B-C1B	-2.56	106.91	109.72
3	B	1353	NAP	C3N-C2N-N1N	-2.50	117.48	120.36
3	C	2353	NAP	C4B-O4B-C1B	-2.49	106.98	109.72
3	B	1353	NAP	C2B-C3B-C4B	-2.47	96.00	101.85
3	B	1353	NAP	O3-PA-O5B	-2.43	96.50	102.94
3	A	353	NAP	O4B-C1B-C2B	-2.38	102.30	106.60
3	D	3353	NAP	C3N-C2N-N1N	-2.35	117.66	120.36
3	C	2353	NAP	C1B-N9A-C4A	-2.31	123.46	126.94
3	A	353	NAP	C4D-O4D-C1D	2.04	111.97	109.72
3	B	1353	NAP	C3N-C7N-N7N	2.06	120.08	117.82
3	D	3353	NAP	C5N-C4N-C3N	2.14	123.03	120.33
3	A	353	NAP	O7N-C7N-C3N	2.16	121.95	119.59
3	D	3353	NAP	C3N-C7N-N7N	2.19	120.22	117.82
3	B	1353	NAP	O2B-P2B-O1X	2.20	112.59	107.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1353	NAP	O3-PN-O5D	2.21	108.80	102.94
3	A	353	NAP	C4A-C5A-N7A	2.21	111.51	109.48
3	B	1353	NAP	O2B-C2B-C3B	2.27	120.34	111.51
3	B	1353	NAP	C5N-C4N-C3N	2.34	123.28	120.33
3	C	2353	NAP	O7N-C7N-C3N	2.43	122.24	119.59
3	C	2353	NAP	C4A-C5A-N7A	2.44	111.72	109.48
3	C	2353	NAP	O3D-C3D-C4D	2.48	118.50	111.05
3	C	2353	NAP	C5N-C4N-C3N	2.90	123.99	120.33
3	A	353	NAP	O3-PN-O5D	3.00	110.88	102.94
3	B	1353	NAP	O7N-C7N-C3N	3.00	122.86	119.59
3	A	353	NAP	C5N-C4N-C3N	3.03	124.15	120.33
3	A	353	NAP	C3N-C7N-N7N	3.11	121.22	117.82
3	C	2353	NAP	C3N-C7N-N7N	3.14	121.25	117.82
3	D	3353	NAP	O7N-C7N-C3N	3.14	123.02	119.59
3	A	353	NAP	O4D-C1D-N1N	3.93	112.44	108.13
3	C	2353	NAP	O3-PN-O5D	4.09	113.78	102.94
3	C	2353	NAP	O4D-C1D-N1N	5.86	114.57	108.13

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 30 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	353	NAP	9	0
3	B	1353	NAP	6	0
3	C	2353	NAP	7	0
3	D	3353	NAP	8	0

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	352/352 (100%)	-0.44	0 100 100	16, 32, 54, 67	0
1	B	352/352 (100%)	-0.43	0 100 100	14, 32, 50, 62	0
1	C	352/352 (100%)	-0.38	3 (0%) 85 79	20, 38, 66, 77	0
1	D	352/352 (100%)	0.01	11 (3%) 52 40	23, 48, 75, 87	0
All	All	1408/1408 (100%)	-0.31	14 (0%) 84 77	14, 36, 66, 87	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	53	HIS	4.5
1	C	352	ALA	4.5
1	D	55	MET	3.8
1	D	352	ALA	3.7
1	D	9	ILE	3.0
1	D	10	GLY	2.9
1	D	16	GLU	2.8
1	D	269	GLY	2.7
1	D	222	PRO	2.5
1	D	13	GLY	2.5
1	D	4	PHE	2.5
1	D	52	ARG	2.1
1	C	218	TYR	2.0
1	C	270	GLU	2.0

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

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

6.3 Carbohydrates [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(\AA^2)	Q<0.9
3	NAP	C	2353	48/48	0.81	0.29	2.47	40,88,99,101	0
3	NAP	A	353	48/48	0.86	0.23	2.40	40,65,82,86	0
3	NAP	D	3353	48/48	0.82	0.27	1.63	40,89,98,99	0
3	NAP	B	1353	48/48	0.89	0.21	1.11	40,57,66,68	0
2	ZN	D	3354	1/1	0.94	0.12	-1.36	78,78,78,78	0
2	ZN	A	354	1/1	0.99	0.10	-1.77	52,52,52,52	0
2	ZN	B	1354	1/1	0.98	0.11	-1.94	49,49,49,49	0
2	ZN	C	2354	1/1	0.98	0.10	-3.15	69,69,69,69	0

6.5 Other polymers [i](#)

There are no such residues in this entry.