



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:15 PM GMT

PDB ID : 4O9U  
Title : Mechanism of transhydrogenase coupling proton translocation and hydride transfer  
Authors : Leung, J.H.; Yamaguchi, M.; Moeller, A.; Schurig-Briccio, L.A.; Gennis, R.B.; Potter, C.S.; Carragher, B.; Stout, C.D.  
Deposited on : 2014-01-02  
Resolution : 6.93 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](mailto: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.

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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

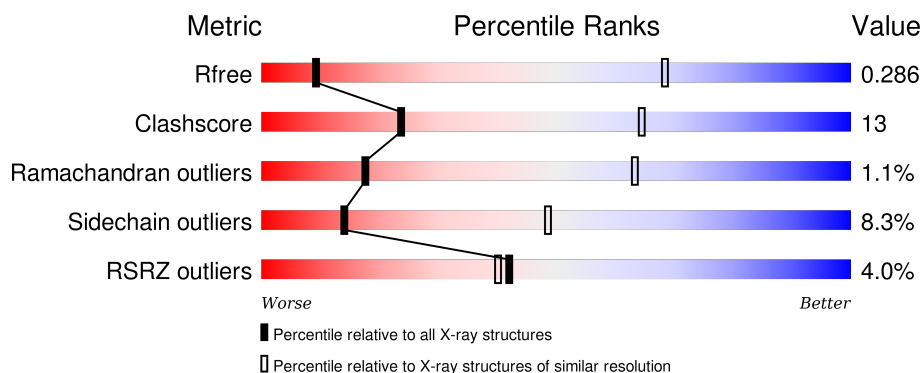
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 6.93 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	1014 (9.50-3.66)
Clashscore	102246	1063 (10.00-3.70)
Ramachandran outliers	100387	1035 (9.50-3.66)
Sidechain outliers	100360	1005 (9.50-3.66)
RSRZ outliers	91569	1013 (9.50-3.66)

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  $\geq 3$ , 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	100	<div> <div>3%</div> <div>72%</div> <div>20%</div> <div>6%</div> </div>
1	C	100	<div> <div>%</div> <div>63%</div> <div>25%</div> <div>10%</div> </div>
2	B	450	<div> <div>2%</div> <div>68%</div> <div>27%</div> <div>.</div> </div>
2	D	450	<div> <div>2%</div> <div>64%</div> <div>30%</div> <div>.</div> </div>
3	E	384	<div> <div>7%</div> <div>68%</div> <div>25%</div> <div>.</div> </div>

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Mol	Chain	Length	Quality of chain
3	F	384	

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
4	NAP	D	500	-	-	-	X

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 13618 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 NAD(P) transhydrogenase subunit alpha 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	94	Total	C	N	O	S	0	0	0
			713	475	114	119	5			
1	C	90	Total	C	N	O	S	0	0	0
			677	453	106	113	5			

- Molecule 2 is a protein called NAD(P) transhydrogenase subunit beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	440	Total	C	N	O	S	0	0	0
			3250	2131	534	563	22			
2	D	437	Total	C	N	O	S	0	0	0
			3235	2122	531	560	22			

- Molecule 3 is a protein called NAD/NADP transhydrogenase alpha subunit 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	372	Total	C	N	O	S	0	0	0
			2794	1772	498	512	12			
3	F	369	Total	C	N	O	S	0	0	0
			2765	1753	494	506	12			

There are 18 discrepancies between the modelled and reference sequences:

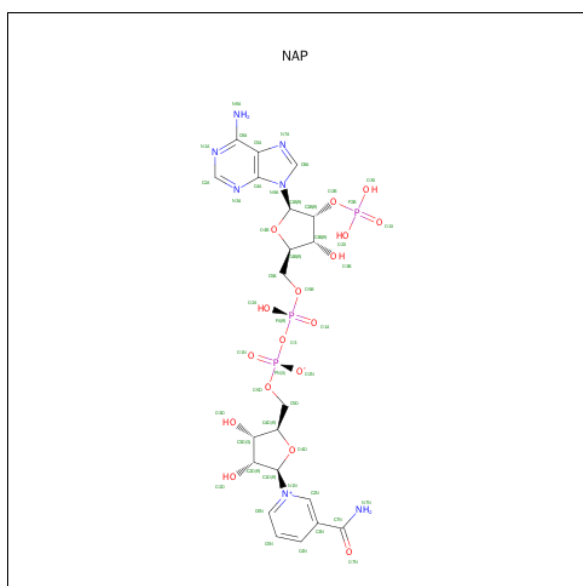
Chain	Residue	Modelled	Actual	Comment	Reference
E	-8	MET	-	EXPRESSION TAG	UNP Q72GR8
E	-7	HIS	-	EXPRESSION TAG	UNP Q72GR8
E	-6	HIS	-	EXPRESSION TAG	UNP Q72GR8
E	-5	HIS	-	EXPRESSION TAG	UNP Q72GR8
E	-4	HIS	-	EXPRESSION TAG	UNP Q72GR8
E	-3	HIS	-	EXPRESSION TAG	UNP Q72GR8
E	-2	HIS	-	EXPRESSION TAG	UNP Q72GR8
E	-1	HIS	-	EXPRESSION TAG	UNP Q72GR8

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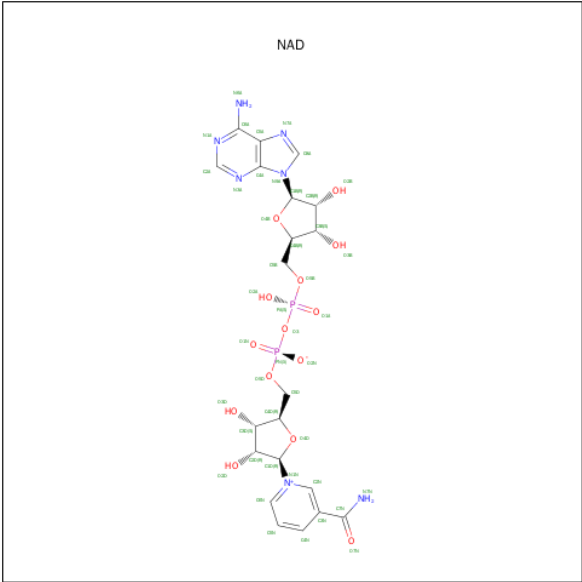
Chain	Residue	Modelled	Actual	Comment	Reference
E	0	HIS	-	EXPRESSION TAG	UNP Q72GR8
F	-8	MET	-	EXPRESSION TAG	UNP Q72GR8
F	-7	HIS	-	EXPRESSION TAG	UNP Q72GR8
F	-6	HIS	-	EXPRESSION TAG	UNP Q72GR8
F	-5	HIS	-	EXPRESSION TAG	UNP Q72GR8
F	-4	HIS	-	EXPRESSION TAG	UNP Q72GR8
F	-3	HIS	-	EXPRESSION TAG	UNP Q72GR8
F	-2	HIS	-	EXPRESSION TAG	UNP Q72GR8
F	-1	HIS	-	EXPRESSION TAG	UNP Q72GR8
F	0	HIS	-	EXPRESSION TAG	UNP Q72GR8

- Molecule 4 is NADP NICOTINAMIDE-ADENINE-DINUCLEOTIDE PHOSPHATE (three-letter code: NAP) (formula:  $C_{21}H_{28}N_7O_{17}P_3$ ).



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

- Molecule 5 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (three-letter code: NAD) (formula:  $C_{21}H_{27}N_7O_{14}P_2$ ).



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

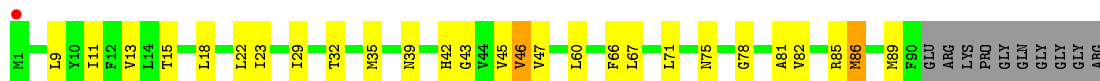
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

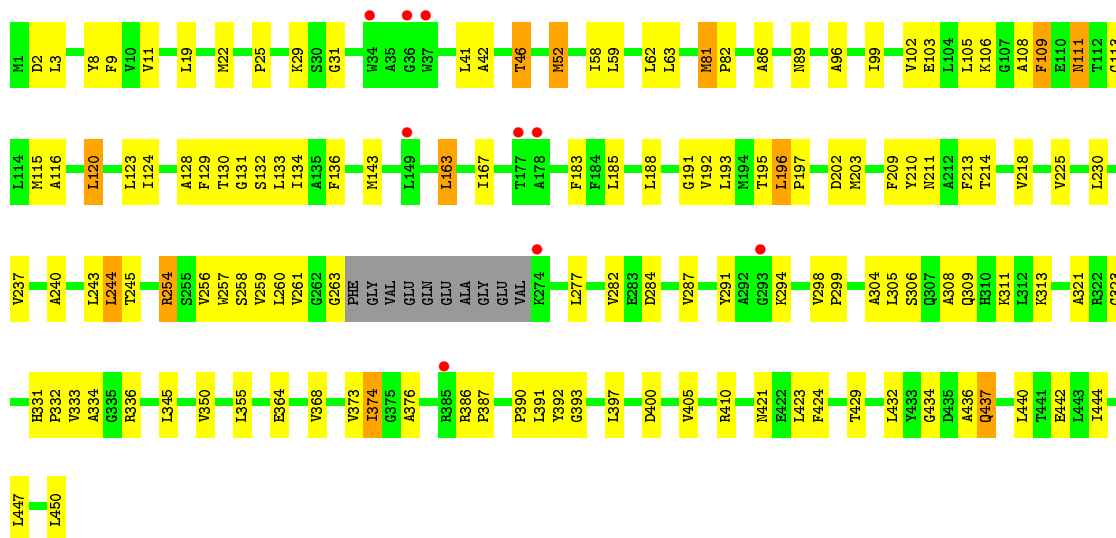
- Molecule 1: NAD(P) transhydrogenase subunit alpha 2



- Molecule 1: NAD(P) transhydrogenase subunit alpha 2

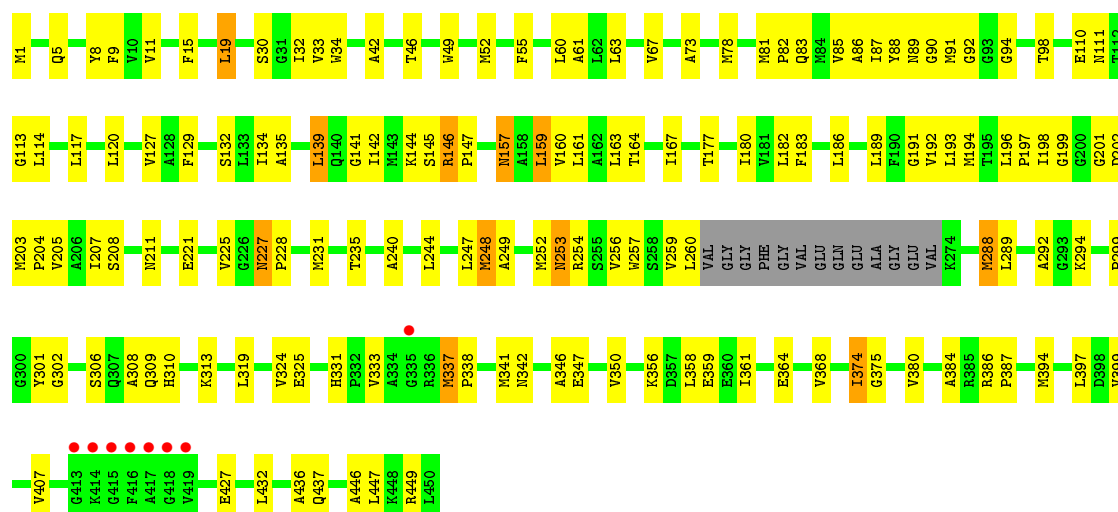


- Molecule 2: NAD(P) transhydrogenase subunit beta

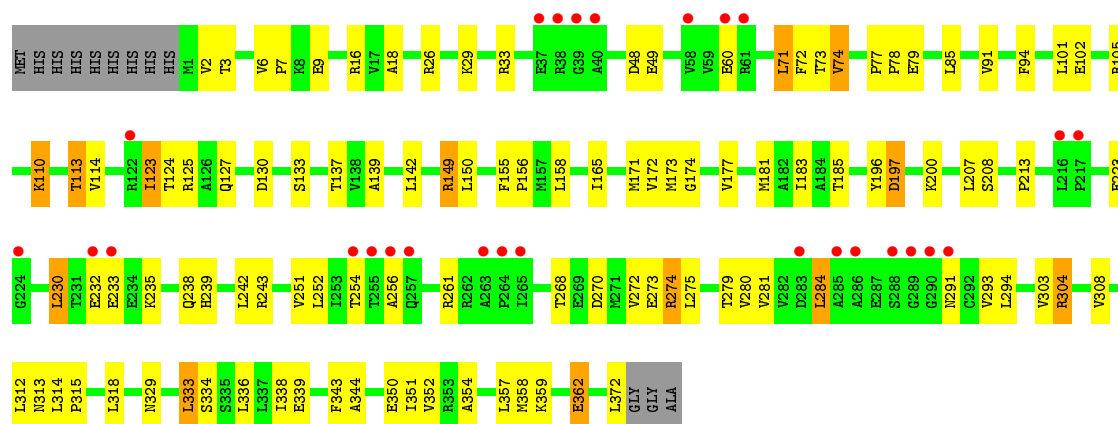


- Molecule 2: NAD(P) transhydrogenase subunit beta

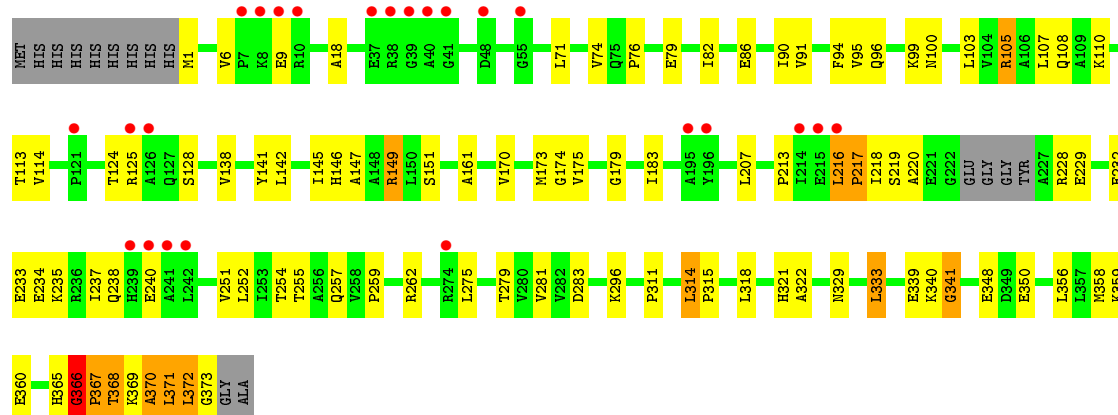




• Molecule 3: NAD/NADP transhydrogenase alpha subunit 1



• Molecule 3: NAD/NADP transhydrogenase alpha subunit 1





## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	175.33Å 160.59Å 139.62Å 90.00° 128.36° 90.00°	Depositor
Resolution (Å)	38.90 – 6.93 38.89 – 6.93	Depositor EDS
% Data completeness (in resolution range)	98.1 (38.90-6.93) 98.8 (38.89-6.93)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.72 (at 6.64Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, $R_{free}$	0.213 , 0.284 0.218 , 0.286	Depositor DCC
$R_{free}$ test set	228 reflections (4.86%)	DCC
Wilson B-factor (Å <sup>2</sup> )	176.7	Xtriage
Anisotropy	0.322	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.18 , 71.7	EDS
Estimated twinning fraction	0.048 for h,-k,-h-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.41$ , $\langle L^2 \rangle = 0.23$	Xtriage
Outliers	1 of 4922 reflections (0.020%)	Xtriage
$F_o, F_c$ correlation	0.84	EDS
Total number of atoms	13618	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	159.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.35% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

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

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z  > 5$	RMSZ	$\# Z  > 5$
1	A	0.35	0/730	0.54	0/990
1	C	0.34	0/693	0.53	0/941
2	B	0.32	0/3314	0.52	0/4503
2	D	0.33	0/3299	0.52	0/4483
3	E	0.32	0/2840	0.54	0/3852
3	F	0.34	1/2809 (0.0%)	0.57	1/3810 (0.0%)
All	All	0.33	1/13685 (0.0%)	0.54	1/18579 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	F	367	PRO	N-CD	5.20	1.55	1.47

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	366	GLY	C-N-CD	5.73	140.43	128.40

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	713	0	739	17	0
1	C	677	0	700	26	0
2	B	3250	0	3399	98	0
2	D	3235	0	3384	114	0
3	E	2794	0	2911	77	0
3	F	2765	0	2881	75	0
4	B	48	0	25	1	0
4	D	48	0	25	5	0
5	E	44	0	26	5	0
5	F	44	0	26	2	0
All	All	13618	0	14116	366	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 13.

All (366) 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)
3:F:365:HIS:ND1	3:F:368:THR:CG2	1.96	1.26
3:F:371:LEU:O	3:F:373:GLY:N	1.66	1.26
3:F:365:HIS:CB	3:F:368:THR:HG21	1.74	1.16
3:F:114:VAL:HG12	3:F:358:MET:HG2	1.30	1.12
2:B:103:GLU:OE2	2:B:225:VAL:HG11	1.50	1.10
2:D:164:THR:HG22	2:D:186:LEU:HB3	1.34	1.10
3:F:365:HIS:HB3	3:F:368:THR:CG2	1.84	1.06
3:E:274:ARG:HG3	3:E:274:ARG:HH11	1.17	1.05
3:F:365:HIS:ND1	3:F:368:THR:HG23	1.66	1.05
3:F:365:HIS:HB3	3:F:368:THR:HG21	1.00	0.98
3:F:367:PRO:O	3:F:370:ALA:N	1.98	0.96
3:F:371:LEU:O	3:F:373:GLY:O	1.84	0.95
3:F:365:HIS:CG	3:F:368:THR:CG2	2.51	0.92
3:F:216:LEU:HB3	3:F:217:PRO:HA	1.52	0.92
2:B:202:ASP:HB3	2:B:259:VAL:HG13	1.51	0.91
3:F:367:PRO:O	3:F:368:THR:C	2.11	0.89
3:E:3:THR:HG22	3:E:33:ARG:HB2	1.56	0.86
2:B:103:GLU:OE2	2:B:225:VAL:CG1	2.23	0.86
3:E:133:SER:HA	3:E:177:VAL:HG11	1.58	0.86
3:F:371:LEU:HD12	3:F:371:LEU:O	1.77	0.84
2:D:252:MET:HA	2:D:333:VAL:HG11	1.60	0.84
3:F:365:HIS:ND1	3:F:368:THR:HG22	1.90	0.83
3:F:371:LEU:HD12	3:F:373:GLY:O	1.78	0.83
3:F:365:HIS:CG	3:F:368:THR:HG21	2.12	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:216:LEU:HB3	3:F:217:PRO:CA	2.08	0.82
2:D:384:ALA:HB1	2:D:394:MET:HB3	1.61	0.82
3:F:91:VAL:HB	3:F:114:VAL:HG23	1.63	0.81
3:F:371:LEU:O	3:F:372:LEU:C	2.18	0.79
2:B:25:PRO:HB3	2:B:291:TYR:OH	1.82	0.79
2:D:201:GLY:C	4:D:500:NAP:H5N	2.04	0.78
3:E:238:GLN:O	3:E:242:LEU:HB2	1.82	0.77
1:C:47:VAL:HG13	2:D:46:THR:HG21	1.66	0.77
2:D:198:ILE:HD12	2:D:259:VAL:HG12	1.66	0.77
2:D:198:ILE:HD11	2:D:260:LEU:HD23	1.66	0.76
1:C:39:ASN:HA	1:C:42:HIS:CE1	2.19	0.76
2:D:30:SER:O	2:D:34:TRP:HD1	1.70	0.75
3:E:274:ARG:HG3	3:E:274:ARG:NH1	1.97	0.74
2:D:225:VAL:HG23	2:D:227:ASN:HB2	1.69	0.74
3:F:365:HIS:CB	3:F:368:THR:CG2	2.51	0.74
2:D:248:MET:SD	2:D:338:PRO:HB3	2.28	0.73
1:C:35:MET:HE1	2:D:338:PRO:HG2	1.69	0.73
3:F:365:HIS:CG	3:F:368:THR:HG23	2.22	0.73
3:F:367:PRO:O	3:F:369:LYS:N	2.22	0.73
3:F:371:LEU:C	3:F:373:GLY:N	2.42	0.73
2:B:392:TYR:O	3:E:125:ARG:HA	1.91	0.71
2:B:261:VAL:HG12	2:B:263:GLY:H	1.53	0.71
2:B:86:ALA:O	2:B:132:SER:HB3	1.90	0.71
3:E:274:ARG:HH11	3:E:274:ARG:CG	2.00	0.69
3:E:272:VAL:HG11	3:E:294:LEU:HD11	1.73	0.69
2:D:201:GLY:O	2:D:204:PRO:HD2	1.93	0.69
2:B:209:PHE:HD1	2:B:245:THR:HG21	1.58	0.68
3:E:110:LYS:HA	3:E:110:LYS:HE3	1.75	0.68
3:F:113:THR:HG23	3:F:359:LYS:HA	1.75	0.68
2:B:103:GLU:OE1	2:B:225:VAL:HG13	1.94	0.68
2:B:191:GLY:O	2:B:195:THR:HG22	1.92	0.67
2:B:261:VAL:HG12	2:B:263:GLY:N	2.09	0.67
2:D:81:MET:N	2:D:82:PRO:HD2	2.09	0.67
2:D:63:LEU:O	2:D:67:VAL:HG22	1.93	0.66
3:E:149:ARG:HH11	3:F:146:HIS:HE1	1.43	0.66
3:E:114:VAL:HG13	3:E:357:LEU:HB3	1.77	0.66
3:F:371:LEU:C	3:F:371:LEU:HD12	2.16	0.65
2:D:82:PRO:HG2	2:D:83:GLN:OE1	1.95	0.65
2:D:204:PRO:HG2	4:D:500:NAP:H4N	1.77	0.65
2:B:131:GLY:HA2	2:B:195:THR:HG21	1.79	0.65
3:E:149:ARG:NH2	3:F:149:ARG:HH21	1.95	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:218:ILE:HG22	3:F:219:SER:H	1.62	0.64
3:E:256:ALA:HB3	3:E:291:ASN:HD21	1.61	0.64
3:E:127:GLN:HA	3:E:130:ASP:HB2	1.80	0.64
1:C:39:ASN:HD21	2:D:88:TYR:HB2	1.62	0.63
2:B:299:PRO:HA	2:B:374:ILE:HG23	1.79	0.63
2:B:103:GLU:OE1	2:B:225:VAL:CG1	2.47	0.63
3:F:371:LEU:O	3:F:373:GLY:CA	2.47	0.63
2:B:103:GLU:CD	2:B:225:VAL:CG1	2.68	0.62
2:D:204:PRO:HG3	2:D:337:MET:HA	1.82	0.62
3:F:275:LEU:HD22	3:F:279:THR:HG21	1.81	0.62
2:D:319:LEU:HB3	2:D:324:VAL:HB	1.80	0.62
3:E:137:THR:O	3:E:181:MET:HG2	2.00	0.62
3:E:171:MET:HB3	3:E:252:LEU:HD13	1.82	0.62
3:F:365:HIS:O	3:F:368:THR:OG1	2.18	0.61
2:D:160:VAL:O	2:D:164:THR:HG23	1.99	0.61
3:E:174:GLY:HA2	5:E:500:NAD:O2A	2.00	0.61
2:B:391:LEU:HA	3:E:124:THR:OG1	2.00	0.61
2:D:253:ASN:HA	2:D:394:MET:HA	1.83	0.60
2:D:193:LEU:HD23	2:D:196:LEU:HD22	1.83	0.60
2:B:254:ARG:HG2	2:B:258:SER:OG	2.01	0.60
2:D:8:TYR:HA	2:D:11:VAL:HB	1.83	0.59
3:E:16:ARG:HH21	3:E:313:ASN:HD21	1.50	0.59
2:D:177:THR:HA	2:D:180:ILE:HD12	1.84	0.59
3:E:171:MET:HB3	3:E:252:LEU:CD1	2.33	0.59
2:B:256:VAL:O	2:B:260:LEU:HG	2.02	0.59
1:C:35:MET:CE	2:D:338:PRO:HG2	2.32	0.59
3:F:311:PRO:HB2	3:F:314:LEU:HD13	1.85	0.59
1:A:42:HIS:CE1	1:A:75:ASN:HD21	2.20	0.59
2:D:198:ILE:CD1	2:D:259:VAL:HG12	2.33	0.59
3:E:238:GLN:O	3:E:242:LEU:CB	2.51	0.58
3:E:256:ALA:HB1	5:E:500:NAD:H52A	1.86	0.58
1:A:42:HIS:HD2	2:B:237:VAL:HG11	1.68	0.58
2:B:89:ASN:HD21	2:B:211:ASN:HA	1.68	0.58
2:B:400:ASP:HB3	2:B:423:LEU:HD13	1.83	0.58
2:B:282:VAL:HG11	2:B:442:GLU:OE1	2.03	0.58
1:C:43:GLY:O	1:C:46:VAL:HG23	2.04	0.58
3:E:139:ALA:HB1	3:E:284:LEU:HD21	1.85	0.58
1:C:11:ILE:O	1:C:15:THR:HG22	2.05	0.57
2:B:306:SER:CB	2:B:436:ALA:HB3	2.35	0.57
1:A:51:VAL:HA	2:B:52:MET:CE	2.34	0.57
3:E:155:PHE:N	3:E:156:PRO:HD2	2.20	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:371:LEU:O	3:F:373:GLY:C	2.42	0.57
2:D:358:LEU:HG	2:D:359:GLU:HG3	1.86	0.57
3:E:243:ARG:HH11	3:E:270:ASP:HB2	1.69	0.57
2:D:87:ILE:O	2:D:91:MET:HG2	2.04	0.57
2:B:332:PRO:HB3	3:F:161:ALA:HA	1.85	0.56
2:D:30:SER:O	2:D:34:TRP:CD1	2.57	0.56
2:B:294:LYS:HG2	2:B:368:VAL:HG22	1.88	0.56
3:F:114:VAL:CG1	3:F:358:MET:HG2	2.21	0.56
2:D:8:TYR:OH	2:D:46:THR:HG22	2.06	0.56
3:F:90:ILE:HG12	3:F:113:THR:HB	1.88	0.56
3:E:350:GLU:O	3:E:354:ALA:HB3	2.05	0.56
3:E:173:MET:HB2	3:E:254:THR:HG22	1.87	0.56
2:D:331:HIS:ND1	2:D:333:VAL:HG22	2.21	0.56
2:D:289:LEU:HB3	2:D:319:LEU:HD21	1.87	0.55
3:F:141:TYR:O	3:F:145:ILE:HG12	2.05	0.55
2:B:116:ALA:O	2:B:120:LEU:HB2	2.06	0.55
2:D:82:PRO:HB3	2:D:337:MET:HB2	1.87	0.55
3:E:273:GLU:HG2	3:E:304:ARG:HH22	1.70	0.55
1:A:3:PHE:N	1:A:4:GLY:HA2	2.21	0.55
3:F:175:VAL:HA	3:F:179:GLY:HA3	1.87	0.55
3:F:76:PRO:HB3	3:F:95:VAL:HA	1.88	0.55
2:D:86:ALA:O	2:D:132:SER:HB3	2.06	0.55
1:A:52:VAL:HG21	2:B:99:ILE:HG23	1.88	0.55
2:D:146:ARG:HB3	2:D:147:PRO:CD	2.37	0.55
2:D:288:MET:O	2:D:292:ALA:HB2	2.07	0.55
3:E:114:VAL:HG12	3:E:358:MET:HB2	1.90	0.54
3:F:1:MET:HA	3:F:341:GLY:O	2.07	0.54
2:B:42:ALA:O	2:B:46:THR:HG22	2.07	0.54
3:E:233:GLU:O	3:E:233:GLU:HG3	2.08	0.54
3:F:173:MET:HB2	3:F:254:THR:HG22	1.87	0.54
2:B:313:LYS:HG3	2:B:350:VAL:HG22	1.89	0.54
2:B:374:ILE:HD11	2:B:436:ALA:HB1	1.88	0.54
3:F:9:GLU:HG3	3:F:74:VAL:HB	1.89	0.54
2:B:311:LYS:O	2:B:444:ILE:HD11	2.08	0.54
3:E:6:VAL:HG23	3:E:18:ALA:HB3	1.90	0.54
2:D:374:ILE:HD11	2:D:436:ALA:HB1	1.90	0.54
2:B:390:PRO:O	3:E:124:THR:OG1	2.24	0.54
2:B:58:ILE:O	2:B:62:LEU:HG	2.08	0.54
1:A:51:VAL:HA	2:B:52:MET:HE2	1.90	0.53
1:C:71:LEU:O	2:D:91:MET:HB3	2.08	0.53
1:C:29:ILE:HD12	1:C:29:ILE:H	1.71	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:81:MET:N	2:D:82:PRO:CD	2.72	0.53
3:E:275:LEU:HD22	3:E:279:THR:HG21	1.88	0.53
2:D:157:ASN:HD21	2:D:193:LEU:HD22	1.73	0.53
2:B:440:LEU:O	2:B:444:ILE:HG13	2.08	0.53
2:B:111:ASN:HD22	2:B:113:GLY:H	1.57	0.53
2:B:89:ASN:ND2	2:B:211:ASN:HA	2.24	0.53
3:F:368:THR:OG1	3:F:369:LYS:N	2.39	0.52
2:B:25:PRO:CB	2:B:291:TYR:OH	2.54	0.52
2:B:103:GLU:CD	2:B:225:VAL:HG11	2.24	0.52
1:A:38:SER:HB2	2:B:237:VAL:HG13	1.92	0.52
2:B:421:ASN:HB3	2:B:424:PHE:CE2	2.43	0.52
3:F:311:PRO:HB2	3:F:314:LEU:CD1	2.40	0.52
3:F:367:PRO:C	3:F:369:LYS:N	2.63	0.52
2:D:306:SER:HB2	4:D:500:NAP:N6A	2.24	0.52
3:E:105:ARG:HH12	3:E:372:LEU:HA	1.75	0.52
3:F:175:VAL:HG11	3:F:183:ILE:HD11	1.90	0.52
1:C:89:MET:HB3	2:D:346:ALA:HB1	1.92	0.52
3:F:371:LEU:C	3:F:373:GLY:H	2.07	0.52
3:E:336:LEU:HD23	3:E:352:VAL:HG23	1.92	0.52
2:B:106:LYS:HD3	2:B:108:ALA:HB2	1.92	0.52
2:D:201:GLY:O	4:D:500:NAP:H5N	2.09	0.51
3:F:283:ASP:O	3:F:311:PRO:HD2	2.10	0.51
2:D:252:MET:C	2:D:254:ARG:H	2.12	0.51
2:D:231:MET:O	2:D:235:THR:HG22	2.11	0.51
1:A:13:VAL:HG21	2:D:9:PHE:HZ	1.75	0.51
2:B:96:ALA:HB2	2:B:218:VAL:HG12	1.93	0.51
2:B:294:LYS:NZ	2:B:364:GLU:HG2	2.26	0.51
1:C:85:ARG:O	1:C:89:MET:HG2	2.11	0.51
2:B:210:TYR:HA	2:B:213:PHE:HB2	1.92	0.51
2:D:253:ASN:N	2:D:253:ASN:OD1	2.42	0.51
1:A:42:HIS:CD2	2:B:237:VAL:HG11	2.45	0.51
2:D:198:ILE:HD12	2:D:259:VAL:CG1	2.39	0.51
2:B:391:LEU:HD23	3:E:124:THR:HG21	1.93	0.51
2:D:204:PRO:HB3	2:D:338:PRO:HG3	1.93	0.50
3:E:230:LEU:HD21	3:E:238:GLN:HG3	1.93	0.50
2:D:78:MET:HB2	2:D:347:GLU:OE1	2.10	0.50
2:D:289:LEU:HD22	2:D:319:LEU:HD11	1.93	0.50
3:E:239:HIS:HE1	3:E:268:THR:HG21	1.77	0.50
2:B:89:ASN:O	2:B:214:THR:HG21	2.11	0.50
2:B:192:VAL:O	2:B:196:LEU:HD13	2.11	0.50
2:B:261:VAL:CG1	2:B:263:GLY:C	2.81	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:240:ALA:O	2:D:244:LEU:HB2	2.11	0.49
3:F:233:GLU:HB2	3:F:237:ILE:HD11	1.95	0.49
2:D:249:ALA:O	2:D:252:MET:HB2	2.11	0.49
3:F:232:GLU:O	3:F:233:GLU:HB2	2.13	0.49
1:A:8:ALA:HA	1:A:11:ILE:HB	1.95	0.49
3:F:220:ALA:H	3:F:229:GLU:HB2	1.77	0.49
3:E:334:SER:HB2	3:E:338:ILE:HD12	1.95	0.49
2:B:304:ALA:HB1	3:E:208:SER:HB2	1.95	0.49
1:A:39:ASN:HD21	2:B:89:ASN:HA	1.77	0.49
2:B:129:PHE:O	2:B:133:LEU:HG	2.13	0.49
2:D:134:ILE:HG13	2:D:192:VAL:HG13	1.95	0.48
3:F:138:VAL:HG11	3:F:322:ALA:HB2	1.94	0.48
3:E:74:VAL:O	3:E:94:PHE:HB2	2.14	0.48
2:B:373:VAL:HG13	2:B:376:ALA:HB3	1.95	0.48
2:B:59:LEU:HA	2:B:62:LEU:HD12	1.95	0.48
3:E:339:GLU:HB2	3:E:344:ALA:HB3	1.94	0.48
3:E:26:ARG:HA	3:E:29:LYS:HB2	1.95	0.48
2:D:446:ALA:HA	2:D:449:ARG:HE	1.78	0.48
2:D:164:THR:HG22	2:D:186:LEU:CB	2.25	0.48
2:D:159:LEU:O	2:D:163:LEU:HG	2.13	0.48
2:D:49:TRP:HB3	2:D:52:MET:HG3	1.95	0.48
2:B:392:TYR:O	3:E:125:ARG:HG2	2.14	0.48
2:D:135:ALA:O	2:D:139:LEU:HD12	2.14	0.48
2:D:141:GLY:HA2	2:D:144:LYS:HB2	1.96	0.48
3:F:252:LEU:HB2	3:F:281:VAL:HG22	1.96	0.48
2:D:248:MET:SD	2:D:338:PRO:CB	3.02	0.47
2:D:308:ALA:HA	2:D:437:GLN:HB3	1.96	0.47
2:D:319:LEU:HD22	2:D:324:VAL:HG21	1.96	0.47
3:E:173:MET:HA	3:E:196:TYR:HB3	1.97	0.47
2:B:261:VAL:HG12	2:B:263:GLY:C	2.34	0.47
2:B:52:MET:HE3	2:B:52:MET:HA	1.97	0.47
1:A:46:VAL:O	1:A:50:MET:HG3	2.14	0.47
2:D:356:LYS:HG3	2:D:361:ILE:HD12	1.97	0.47
2:D:380:VAL:HB	2:D:399:VAL:HG11	1.95	0.47
3:E:293:VAL:HG23	3:E:294:LEU:HG	1.96	0.47
2:B:209:PHE:CD1	2:B:245:THR:HG21	2.45	0.47
2:B:193:LEU:HA	2:B:196:LEU:HD22	1.97	0.47
2:D:194:MET:O	2:D:197:PRO:HD2	2.15	0.47
2:B:185:LEU:HD23	2:B:188:LEU:HD12	1.97	0.46
2:B:257:TRP:CE3	2:B:260:LEU:HD12	2.50	0.46
1:C:23:ILE:HD11	2:D:19:LEU:HD22	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:11:VAL:HG21	2:B:42:ALA:HB2	1.97	0.46
3:E:274:ARG:NH1	3:E:274:ARG:CG	2.67	0.46
3:E:26:ARG:HA	3:E:29:LYS:HD2	1.98	0.46
2:D:309:GLN:HE21	2:D:310:HIS:CD2	2.33	0.46
2:D:94:GLY:O	2:D:98:THR:HG23	2.16	0.46
1:C:66:PHE:CD1	2:D:61:ALA:HB2	2.51	0.46
2:B:214:THR:O	2:B:218:VAL:HG13	2.16	0.46
3:F:76:PRO:HG3	3:F:100:ASN:ND2	2.31	0.46
2:B:9:PHE:CZ	1:C:13:VAL:HG21	2.51	0.46
3:F:105:ARG:HA	3:F:108:GLN:HG2	1.98	0.46
3:E:314:LEU:N	3:E:315:PRO:CD	2.79	0.46
2:B:331:HIS:HB3	2:B:334:ALA:HB2	1.98	0.46
3:F:90:ILE:HA	3:F:113:THR:O	2.16	0.46
3:E:149:ARG:HG3	3:E:150:LEU:HG	1.98	0.46
2:D:147:PRO:HD3	2:D:199:GLY:HA2	1.98	0.45
2:D:294:LYS:HB3	2:D:368:VAL:HA	1.97	0.45
1:C:78:GLY:O	1:C:82:VAL:HG22	2.15	0.45
3:E:230:LEU:HD23	3:E:235:LYS:HG3	1.97	0.45
2:B:336:ARG:NH2	4:B:500:NAP:O2A	2.45	0.45
3:F:105:ARG:HG3	3:F:372:LEU:HD12	1.97	0.45
2:D:90:GLY:O	2:D:129:PHE:HD1	1.98	0.45
2:D:127:VAL:HG13	2:D:191:GLY:HA2	1.99	0.45
3:E:251:VAL:HG22	3:E:280:VAL:HB	1.97	0.45
2:D:85:VAL:O	2:D:89:ASN:HB2	2.17	0.45
2:D:302:GLY:HA3	2:D:375:GLY:HA3	1.98	0.45
3:E:2:VAL:HG22	3:E:3:THR:N	2.32	0.45
1:C:71:LEU:HB3	2:D:91:MET:O	2.17	0.45
1:C:75:ASN:HD22	2:D:88:TYR:HB3	1.82	0.45
2:D:221:GLU:O	2:D:225:VAL:HG22	2.17	0.45
2:B:331:HIS:CE1	2:B:333:VAL:HG22	2.52	0.45
3:E:77:PRO:HA	3:E:78:PRO:HD3	1.76	0.45
3:F:207:LEU:HD23	3:F:213:PRO:HD3	1.98	0.45
3:F:257:GLN:H	5:F:500:NAD:H52A	1.82	0.45
2:D:208:SER:HA	2:D:211:ASN:HB3	1.98	0.45
2:D:364:GLU:O	2:D:368:VAL:HG23	2.17	0.45
3:F:238:GLN:HE22	5:F:500:NAD:H62A	1.65	0.45
3:E:207:LEU:HD12	3:E:213:PRO:HD3	1.99	0.45
2:B:81:MET:N	2:B:82:PRO:CD	2.80	0.45
1:C:82:VAL:HG13	2:D:73:ALA:HB1	1.98	0.44
3:E:281:VAL:HB	3:E:308:VAL:HG22	1.99	0.44
2:B:22:MET:HG3	2:B:31:GLY:HA3	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:371:LEU:CD1	3:F:373:GLY:O	2.57	0.44
2:D:42:ALA:O	2:D:46:THR:HG23	2.17	0.44
3:E:303:VAL:HG22	3:E:304:ARG:HE	1.81	0.44
1:C:29:ILE:HA	2:D:342:ASN:ND2	2.33	0.44
1:C:18:LEU:O	1:C:22:LEU:HG	2.17	0.44
3:F:365:HIS:O	3:F:366:GLY:O	2.36	0.44
1:A:52:VAL:HG13	2:B:102:VAL:HG11	2.00	0.44
3:F:339:GLU:C	3:F:341:GLY:H	2.21	0.44
1:A:23:ILE:HB	2:B:243:LEU:HD13	1.99	0.44
3:E:230:LEU:HB2	3:E:235:LYS:NZ	2.32	0.44
2:D:227:ASN:HA	2:D:228:PRO:HD3	1.81	0.44
2:D:144:LYS:HG2	2:D:145:SER:N	2.33	0.44
2:D:11:VAL:HG13	2:D:15:PHE:CE2	2.53	0.43
2:B:284:ASP:O	2:B:287:VAL:HB	2.18	0.43
2:B:130:THR:O	2:B:134:ILE:HG12	2.18	0.43
2:B:196:LEU:N	2:B:197:PRO:CD	2.82	0.43
2:D:113:GLY:O	2:D:117:LEU:HG	2.18	0.43
3:F:314:LEU:N	3:F:315:PRO:CD	2.81	0.43
3:F:74:VAL:O	3:F:94:PHE:HB2	2.19	0.43
2:B:109:PHE:CE1	2:B:115:MET:HG3	2.53	0.43
1:C:86:MET:CE	2:D:81:MET:HG2	2.49	0.43
2:D:358:LEU:HG	2:D:359:GLU:N	2.33	0.43
2:B:410:ARG:HA	2:B:434:GLY:H	1.84	0.42
2:D:301:TYR:HB3	4:D:500:NAP:H52A	2.01	0.42
2:B:128:ALA:HB1	2:B:214:THR:HG23	2.00	0.42
3:E:113:THR:HG23	3:E:359:LYS:HA	2.00	0.42
2:B:128:ALA:HB2	2:B:213:PHE:HB3	2.00	0.42
3:E:155:PHE:N	3:E:156:PRO:CD	2.81	0.42
3:E:139:ALA:CB	3:E:284:LEU:HD21	2.50	0.42
2:D:299:PRO:HA	2:D:374:ILE:HG23	2.00	0.42
3:E:329:ASN:O	3:E:333:LEU:HB2	2.19	0.42
2:D:201:GLY:C	2:D:203:MET:H	2.23	0.42
2:D:337:MET:HB2	2:D:338:PRO:CD	2.49	0.42
2:B:196:LEU:H	2:B:197:PRO:HD3	1.84	0.42
2:B:29:LYS:H	2:B:29:LYS:HG2	1.68	0.42
3:E:235:LYS:HE3	5:E:500:NAD:H2A	2.02	0.42
3:E:125:ARG:HB3	3:E:351:ILE:HD11	2.01	0.42
2:B:240:ALA:O	2:B:244:LEU:HB2	2.19	0.42
2:D:299:PRO:HB2	2:D:341:MET:SD	2.59	0.42
2:D:94:GLY:N	2:D:129:PHE:HB2	2.35	0.42
1:A:13:VAL:HG21	2:D:9:PHE:CZ	2.53	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:111:ASN:HB3	2:D:114:LEU:HG	2.01	0.42
1:C:42:HIS:HE1	2:D:211:ASN:OD1	2.02	0.42
2:D:30:SER:O	2:D:33:VAL:HB	2.20	0.42
3:F:147:ALA:O	3:F:151:SER:HB2	2.19	0.42
3:E:71:LEU:C	3:E:72:PHE:HD1	2.23	0.41
2:D:331:HIS:HD2	2:D:397:LEU:HB2	1.84	0.41
3:E:91:VAL:HB	3:E:114:VAL:HA	2.01	0.41
2:D:167:ILE:HG22	2:D:183:PHE:HB2	2.01	0.41
2:B:345:LEU:HD22	2:B:350:VAL:HG21	2.02	0.41
3:E:9:GLU:HG3	3:E:74:VAL:HB	2.01	0.41
2:D:201:GLY:O	2:D:203:MET:N	2.53	0.41
3:E:165:ILE:HG12	3:F:321:HIS:CD2	2.55	0.41
2:B:386:ARG:HA	2:B:387:PRO:HD3	1.89	0.41
3:F:6:VAL:HG13	3:F:18:ALA:HB3	2.03	0.41
2:B:321:ALA:C	2:B:323:GLY:H	2.23	0.41
2:B:392:TYR:CG	2:B:393:GLY:N	2.88	0.41
1:A:3:PHE:N	1:A:4:GLY:CA	2.83	0.41
3:E:172:VAL:HG21	3:E:183:ILE:HG13	2.02	0.41
3:E:7:PRO:HG3	3:E:73:THR:HG22	2.02	0.41
3:F:329:ASN:O	3:F:333:LEU:HB2	2.21	0.41
1:C:42:HIS:HD2	1:C:75:ASN:OD1	2.04	0.41
2:D:163:LEU:O	2:D:167:ILE:HG13	2.21	0.41
3:F:79:GLU:OE1	3:F:103:LEU:HD23	2.20	0.41
1:A:48:GLY:HA3	2:B:230:LEU:HD11	2.03	0.41
3:E:142:LEU:HD11	3:E:318:LEU:HB2	2.03	0.41
3:E:197:ASP:OD1	5:E:500:NAD:O1A	2.38	0.41
3:E:358:MET:HA	3:E:362:GLU:O	2.21	0.41
1:C:71:LEU:HA	2:D:91:MET:HE3	2.03	0.41
3:F:174:GLY:O	3:F:255:THR:OG1	2.31	0.41
2:D:407:VAL:O	2:D:432:LEU:N	2.50	0.41
2:D:205:VAL:HG12	2:D:205:VAL:O	2.19	0.41
2:D:386:ARG:HA	2:D:387:PRO:HD3	1.90	0.41
3:E:174:GLY:CA	5:E:500:NAD:O2A	2.67	0.41
2:B:8:TYR:OH	2:B:46:THR:HB	2.20	0.41
2:B:109:PHE:HE1	2:B:115:MET:HG3	1.86	0.41
2:B:308:ALA:HA	2:B:437:GLN:HB3	2.02	0.41
3:E:124:THR:HA	3:E:127:GLN:CD	2.41	0.40
2:B:124:ILE:HG23	2:B:213:PHE:CD2	2.55	0.40
3:F:82:ILE:HG23	3:F:107:LEU:HD23	2.03	0.40
1:C:45:VAL:HG11	2:D:92:GLY:HA2	2.02	0.40
3:E:149:ARG:CZ	3:F:149:ARG:HH21	2.33	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:298:VAL:HB	2:B:373:VAL:HG22	2.03	0.40
2:D:313:LYS:HG3	2:D:350:VAL:HG22	2.03	0.40
2:B:405:VAL:HB	2:B:429:THR:HA	2.03	0.40
2:B:25:PRO:HG2	2:B:450:LEU:HD22	2.03	0.40
2:B:133:LEU:HA	2:B:136:PHE:HB3	2.03	0.40
2:B:123:LEU:HD12	2:B:183:PHE:HE1	1.85	0.40
3:F:170:VAL:HG22	3:F:251:VAL:HB	2.03	0.40
2:B:163:LEU:O	2:B:167:ILE:HG22	2.22	0.40
2:D:86:ALA:HB2	2:D:207:ILE:HG12	2.02	0.40
3:E:251:VAL:HA	3:E:280:VAL:O	2.21	0.40
2:D:207:ILE:O	2:D:211:ASN:N	2.54	0.40
2:D:294:LYS:HE2	2:D:364:GLU:HG2	2.04	0.40
1:C:81:ALA:HB3	2:D:73:ALA:HB3	2.03	0.40
3:F:110:LYS:HD3	3:F:110:LYS:O	2.21	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	92/100 (92%)	82 (89%)	9 (10%)	1 (1%)	17	63
1	C	88/100 (88%)	77 (88%)	11 (12%)	0	100	100
2	B	436/450 (97%)	413 (95%)	22 (5%)	1 (0%)	52	86
2	D	433/450 (96%)	403 (93%)	29 (7%)	1 (0%)	52	86
3	E	370/384 (96%)	333 (90%)	34 (9%)	3 (1%)	24	69
3	F	365/384 (95%)	321 (88%)	31 (8%)	13 (4%)	4	38
All	All	1784/1868 (96%)	1629 (91%)	136 (8%)	19 (1%)	17	63

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	F	370	ALA
3	F	372	LEU
3	E	123	ILE
3	E	343	PHE
3	F	216	LEU
3	F	366	GLY
2	B	309	GLN
2	D	202	ASP
3	F	217	PRO
3	F	235	LYS
3	F	368	THR
1	A	45	VAL
3	F	356	LEU
3	F	234	GLU
3	F	340	LYS
3	E	74	VAL
3	F	341	GLY
3	F	96	GLN
3	F	259	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	73/75 (97%)	65 (89%)	8 (11%)	8	34
1	C	69/75 (92%)	63 (91%)	6 (9%)	13	45
2	B	327/335 (98%)	301 (92%)	26 (8%)	15	50
2	D	326/335 (97%)	298 (91%)	28 (9%)	13	47
3	E	287/296 (97%)	261 (91%)	26 (9%)	12	43
3	F	284/296 (96%)	264 (93%)	20 (7%)	19	56
All	All	1366/1412 (97%)	1252 (92%)	114 (8%)	14	48

All (114) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	34	LEU
1	A	44	VAL
1	A	45	VAL
1	A	46	VAL
1	A	60	LEU
1	A	64	ILE
1	A	87	LEU
2	B	2	ASP
2	B	3	LEU
2	B	19	LEU
2	B	41	LEU
2	B	46	THR
2	B	52	MET
2	B	63	LEU
2	B	81	MET
2	B	105	LEU
2	B	109	PHE
2	B	111	ASN
2	B	120	LEU
2	B	143	MET
2	B	163	LEU
2	B	196	LEU
2	B	203	MET
2	B	244	LEU
2	B	254	ARG
2	B	277	LEU
2	B	305	LEU
2	B	355	LEU
2	B	374	ILE
2	B	397	LEU
2	B	432	LEU
2	B	437	GLN
2	B	447	LEU
1	C	9	LEU
1	C	32	THR
1	C	46	VAL
1	C	60	LEU
1	C	67	LEU
1	C	86	MET
2	D	1	MET
2	D	5	GLN
2	D	19	LEU

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Mol	Chain	Res	Type
2	D	32	ILE
2	D	55	PHE
2	D	60	LEU
2	D	110	GLU
2	D	120	LEU
2	D	139	LEU
2	D	142	ILE
2	D	146	ARG
2	D	157	ASN
2	D	159	LEU
2	D	161	LEU
2	D	182	LEU
2	D	189	LEU
2	D	227	ASN
2	D	247	LEU
2	D	248	MET
2	D	253	ASN
2	D	256	VAL
2	D	257	TRP
2	D	288	MET
2	D	325	GLU
2	D	337	MET
2	D	374	ILE
2	D	427	GLU
2	D	447	LEU
3	E	48	ASP
3	E	49	GLU
3	E	60	GLU
3	E	71	LEU
3	E	79	GLU
3	E	85	LEU
3	E	101	LEU
3	E	102	GLU
3	E	110	LYS
3	E	113	THR
3	E	123	ILE
3	E	149	ARG
3	E	158	LEU
3	E	185	THR
3	E	197	ASP
3	E	200	LYS
3	E	223	GLU

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Mol	Chain	Res	Type
3	E	230	LEU
3	E	232	GLU
3	E	261	ARG
3	E	274	ARG
3	E	284	LEU
3	E	304	ARG
3	E	312	LEU
3	E	333	LEU
3	E	362	GLU
3	F	71	LEU
3	F	86	GLU
3	F	99	LYS
3	F	105	ARG
3	F	124	THR
3	F	125	ARG
3	F	128	SER
3	F	142	LEU
3	F	149	ARG
3	F	228	ARG
3	F	240	GLU
3	F	262	ARG
3	F	296	LYS
3	F	314	LEU
3	F	318	LEU
3	F	333	LEU
3	F	348	GLU
3	F	350	GLU
3	F	360	GLU
3	F	371	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	39	ASN
1	A	75	ASN
2	B	89	ASN
2	B	111	ASN
2	B	140	GLN
2	B	310	HIS
1	C	39	ASN
1	C	42	HIS
1	C	55	HIS

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Mol	Chain	Res	Type
1	C	75	ASN
2	D	5	GLN
2	D	153	GLN
2	D	157	ASN
2	D	310	HIS
3	E	98	HIS
3	E	291	ASN
3	E	329	ASN
3	F	146	HIS
3	F	238	GLN
3	F	245	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](#)

4 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
4	NAP	B	500	-	42,52,52	0.89	2 (4%)	54,80,80	1.65	5 (9%)
4	NAP	D	500	-	42,52,52	0.89	2 (4%)	54,80,80	1.54	5 (9%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
5	NAD	E	500	-	38,48,48	0.98	2 (5%)	47,73,73	1.59	3 (6%)
5	NAD	F	500	-	38,48,48	0.97	3 (7%)	47,73,73	1.66	5 (10%)

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
4	NAP	B	500	-	-	0/27/67/67	0/5/5/5
4	NAP	D	500	-	-	0/27/67/67	0/5/5/5
5	NAD	E	500	-	-	0/22/62/62	0/5/5/5
5	NAD	F	500	-	-	0/22/62/62	0/5/5/5

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	500	NAP	O4D-C1D	2.03	1.43	1.41
5	F	500	NAD	O4B-C1B	2.05	1.43	1.41
5	E	500	NAD	O4B-C1B	2.30	1.44	1.41
4	B	500	NAP	O4D-C1D	2.36	1.44	1.41
5	F	500	NAD	O4D-C1D	2.43	1.44	1.41
4	B	500	NAP	C5A-C4A	3.22	1.47	1.40
4	D	500	NAP	C5A-C4A	3.25	1.47	1.40
5	F	500	NAD	C5A-C4A	3.32	1.48	1.40
5	E	500	NAD	C5A-C4A	3.56	1.48	1.40

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	500	NAP	N3A-C2A-N1A	-7.29	123.31	128.89
5	F	500	NAD	N3A-C2A-N1A	-7.27	123.33	128.89
5	E	500	NAD	N3A-C2A-N1A	-6.90	123.61	128.89
4	B	500	NAP	N3A-C2A-N1A	-6.89	123.62	128.89
5	E	500	NAD	PN-O3-PA	-4.43	120.28	132.73
4	B	500	NAP	PN-O3-PA	-4.06	121.34	132.73
5	F	500	NAD	PN-O3-PA	-4.05	121.35	132.73
4	D	500	NAP	PN-O3-PA	-3.83	121.99	132.73
5	E	500	NAD	C4A-C5A-N7A	-3.49	106.27	109.48
4	B	500	NAP	C4A-C5A-N7A	-3.14	106.59	109.48
5	F	500	NAD	C4A-C5A-N7A	-3.13	106.60	109.48
4	D	500	NAP	C4A-C5A-N7A	-2.96	106.75	109.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	F	500	NAD	O5D-C5D-C4D	2.20	117.22	109.12
4	D	500	NAP	O4B-C1B-N9A	2.21	112.73	108.10
4	B	500	NAP	O4B-C1B-N9A	2.44	113.20	108.10
4	D	500	NAP	O4D-C1D-N1N	2.83	111.24	108.13
5	F	500	NAD	O4D-C1D-N1N	3.45	111.92	108.13
4	B	500	NAP	O4D-C1D-N1N	5.88	114.60	108.13

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	500	NAP	1	0
4	D	500	NAP	5	0
5	E	500	NAD	5	0
5	F	500	NAD	2	0

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	94/100 (94%)	-0.13	3 (3%)	51	47	107, 187, 276, 328	0
1	C	90/100 (90%)	-0.44	1 (1%)	82	78	84, 136, 201, 285	0
2	B	440/450 (97%)	-0.35	9 (2%)	68	63	65, 148, 218, 287	0
2	D	437/450 (97%)	-0.25	8 (1%)	71	66	73, 152, 251, 298	0
3	E	372/384 (96%)	-0.02	27 (7%)	18	21	50, 145, 221, 249	0
3	F	369/384 (96%)	0.06	24 (6%)	22	23	84, 173, 238, 275	0
All	All	1802/1868 (96%)	-0.17	72 (3%)	42	40	50, 154, 236, 328	0

All (72) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	F	40	ALA	5.3
3	F	39	GLY	5.3
3	E	263	ALA	5.1
1	A	58	THR	4.8
3	E	60	GLU	4.4
3	E	264	PRO	4.4
2	D	335	GLY	4.2
1	A	59	GLY	4.2
2	D	414	LYS	4.1
2	D	418	GLY	4.0
2	D	415	GLY	4.0
3	F	196	TYR	3.9
3	F	10	ARG	3.8
3	F	8	LYS	3.6
2	D	417	ALA	3.4
3	E	291	ASN	3.4
2	D	419	VAL	3.4
3	E	38	ARG	3.3
3	F	38	ARG	3.2

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Mol	Chain	Res	Type	RSRZ
3	E	288	SER	3.2
2	B	274	LYS	3.1
3	F	9	GLU	3.1
3	E	224	GLY	3.1
3	E	40	ALA	3.1
3	E	39	GLY	3.0
3	F	214	ILE	3.0
3	F	239	HIS	2.9
3	F	241	ALA	2.9
3	E	37	GLU	2.8
3	E	290	GLY	2.8
3	E	254	THR	2.8
3	E	217	PRO	2.7
3	F	55	GLY	2.7
3	E	283	ASP	2.7
3	E	257	GLN	2.6
3	F	125	ARG	2.6
3	F	215	GLU	2.6
3	F	240	GLU	2.6
3	F	121	PRO	2.5
1	A	60	LEU	2.5
2	B	149	LEU	2.5
3	E	232	GLU	2.5
3	E	285	ALA	2.5
3	F	48	ASP	2.4
3	E	233	GLU	2.4
2	B	385	ARG	2.4
3	F	242	LEU	2.4
2	B	37	TRP	2.4
3	E	61	ARG	2.4
3	F	7	PRO	2.4
3	F	37	GLU	2.3
2	B	178	ALA	2.3
3	F	126	ALA	2.3
3	E	58	VAL	2.3
3	E	286	ALA	2.3
3	F	195	ALA	2.2
2	D	416	PHE	2.2
3	E	256	ALA	2.2
2	D	413	GLY	2.2
1	C	1	MET	2.2
3	E	216	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
2	B	36	GLY	2.2
2	B	34	TRP	2.2
2	B	177	THR	2.2
3	E	255	THR	2.1
3	E	265	ILE	2.1
3	F	216	LEU	2.1
3	E	122	ARG	2.1
3	F	274	ARG	2.1
2	B	293	GLY	2.1
3	F	41	GLY	2.0
3	E	289	GLY	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
4	NAP	D	500	48/48	0.89	0.47	0.20	190,212,228,229	0
5	NAD	E	500	44/44	0.76	0.37	-0.06	137,203,307,312	0
5	NAD	F	500	44/44	0.86	0.31	-0.22	165,193,250,257	0
4	NAP	B	500	48/48	0.94	0.15	-0.81	89,94,103,107	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.