



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 26, 2026 – 05:05 AM EDT

PDB ID : 9ETJ / pdb_00009etj
Title : Mouse CNPase catalytic domain with nanobody 10E
Authors : Markusson, S.; Raasakka, A.; Opazo, F.; Kursula, P.
Deposited on : 2024-03-26
Resolution : 2.55 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

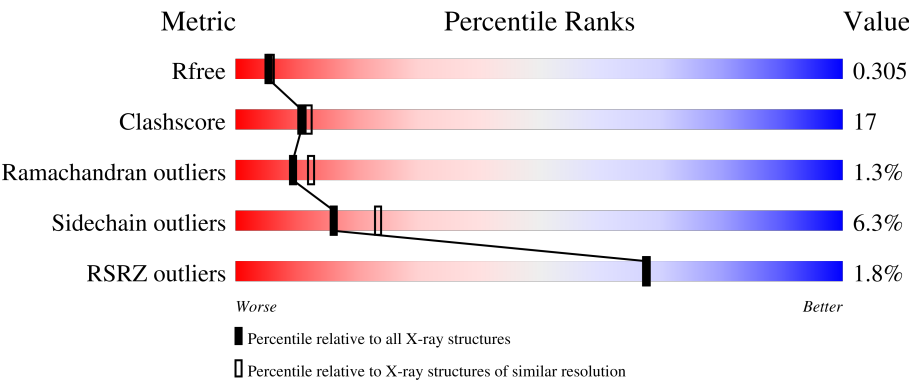
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.55 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}	180053	1091 (2.54-2.54)
Clashscore	190562	1120 (2.54-2.54)
Ramachandran outliers	187476	1106 (2.54-2.54)
Sidechain outliers	187428	1106 (2.54-2.54)
RSRZ outliers	180081	1091 (2.54-2.54)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	220	3% 64% 27% • 7%
1	B	220	2% 66% 26% • 7%
1	E	220	2% 61% 31% • •
1	G	220	% 65% 27% • •
2	C	131	% 60% 32% 8% •

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Mol	Chain	Length	Quality of chain
2	D	131	2% 56% 39%
2	F	131	% 68% 26% 5%
2	H	131	62% 35%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 20778 atoms, of which 10284 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 2',3'-cyclic-nucleotide 3'-phosphodiesterase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	204	Total	C	H	N	O	S	0	0	0
			3198	1026	1610	268	289	5			
1	B	205	Total	C	H	N	O	S	0	0	0
			3215	1032	1621	269	288	5			
1	E	211	Total	C	H	N	O	S	0	0	0
			3306	1060	1664	276	301	5			
1	G	211	Total	C	H	N	O	S	0	0	0
			3306	1060	1664	276	301	5			

- Molecule 2 is a protein called Chains: C,D,F,H.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	C	130	Total	C	H	N	O	S	0	0	0
			1932	630	929	170	199	4			
2	D	130	Total	C	H	N	O	S	0	0	0
			1932	630	929	170	199	4			
2	F	130	Total	C	H	N	O	S	0	0	0
			1932	630	929	170	199	4			
2	H	131	Total	C	H	N	O	S	0	0	0
			1950	635	938	171	202	4			

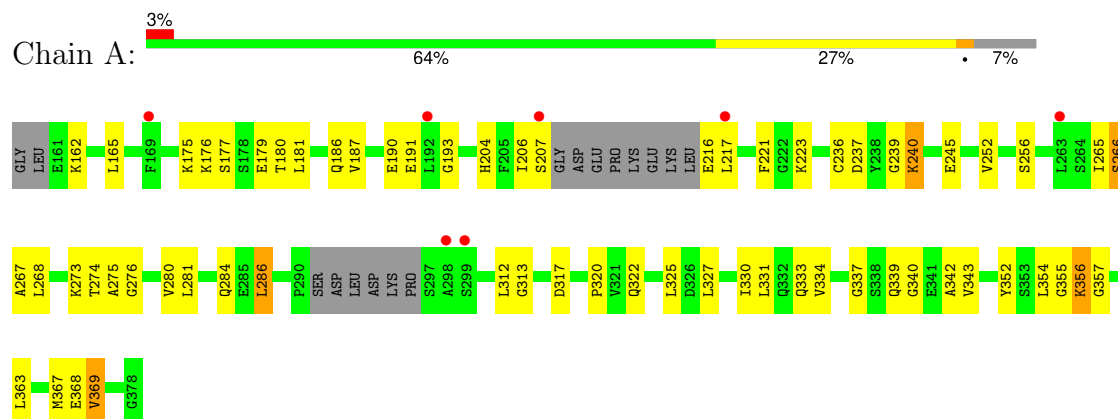
- Molecule 3 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	C	2	Total	Ca	0	0
			2	2		
3	D	2	Total	Ca	0	0
			2	2		
3	F	1	Total	Ca	0	0
			1	1		
3	H	2	Total	Ca	0	0
			2	2		

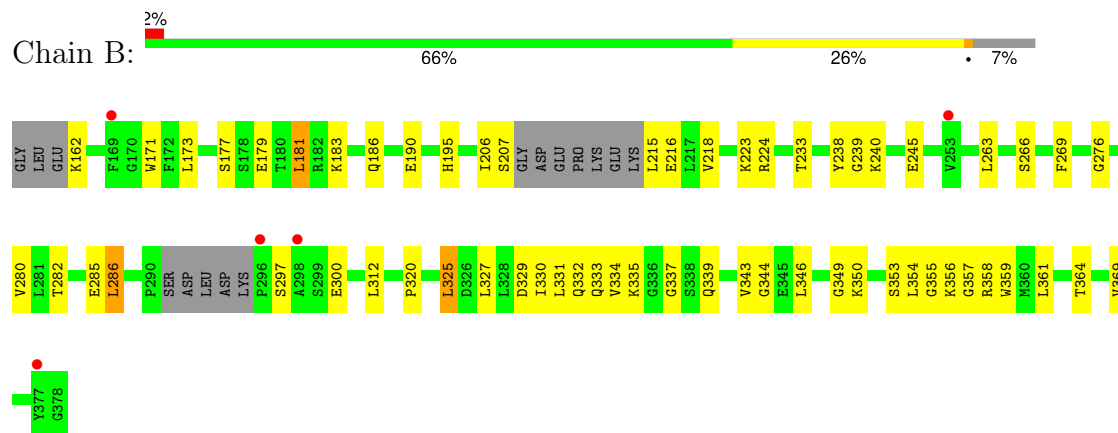
3 Residue-property plots [i](#)

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

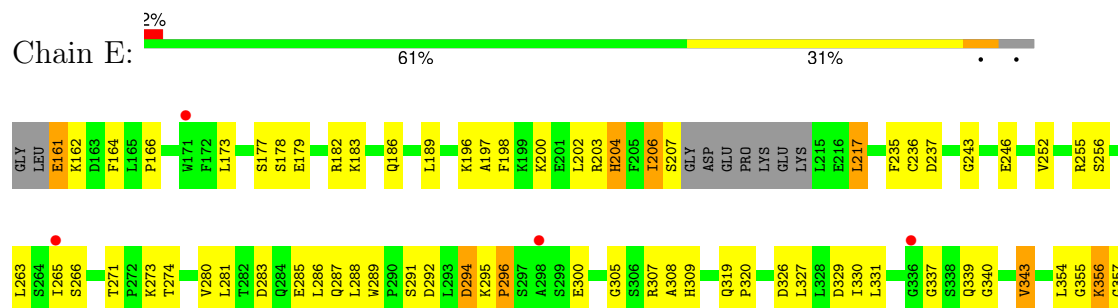
- Molecule 1: 2',3'-cyclic-nucleotide 3'-phosphodiesterase



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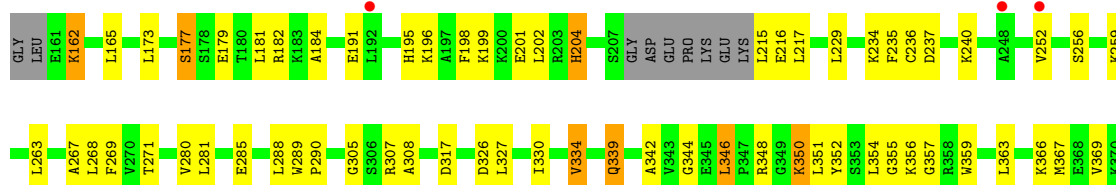


- Molecule 1: 2',3'-cyclic-nucleotide 3'-phosphodiesterase





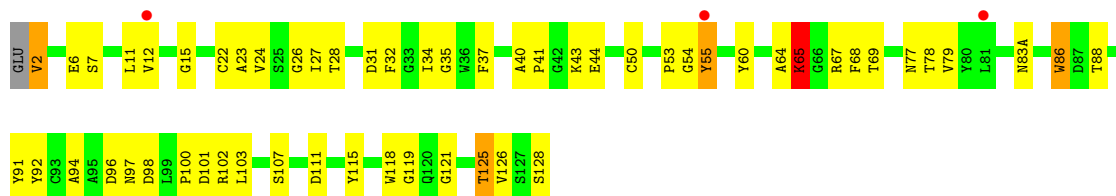
- Molecule 1: 2',3'-cyclic-nucleotide 3'-phosphodiesterase



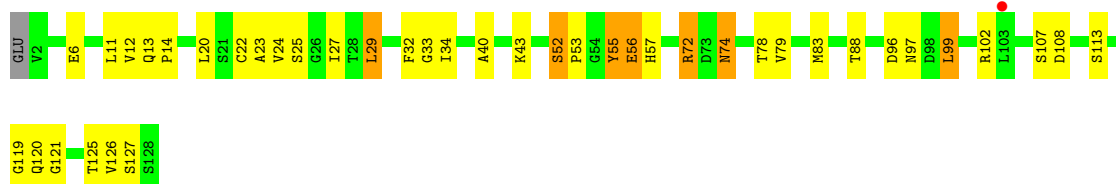
- Molecule 2: Chains: C,D,F,H



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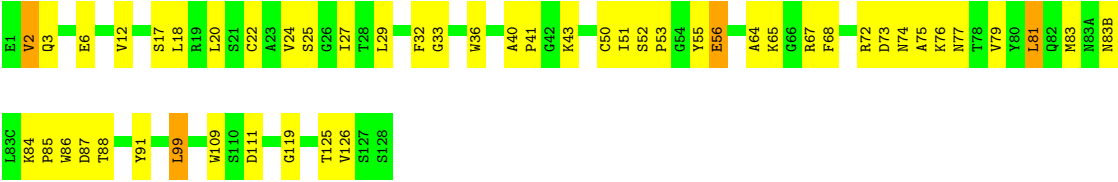


- Molecule 2: Chains: C,D,F,H

Chain H:

62%

35%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	101.73Å 118.25Å 122.45Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.95 – 2.55 47.95 – 2.55	Depositor EDS
% Data completeness (in resolution range)	99.6 (47.95-2.55) 86.8 (47.95-2.55)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.43 (at 2.54Å)	Xtriage
Refinement program	PHENIX 1.21rc1_5127	Depositor
R, R_{free}	0.252 , 0.304 0.252 , 0.305	Depositor DCC
R_{free} test set	2003 reflections (4.11%)	wwPDB-VP
Wilson B-factor (Å ²)	66.2	Xtriage
Anisotropy	0.387	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 64.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.000 for -h,l,k	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	20778	wwPDB-VP
Average B, all atoms (Å ²)	113.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 18.32% 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, 0.333 respectively for untwinned datasets, and 0.375, 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: CA

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	1.00	0/1623	0.97	0/2182
1	B	0.82	0/1630	0.83	0/2192
1	E	0.89	0/1679	0.92	0/2260
1	G	0.79	0/1679	0.84	0/2260
2	C	1.09	2/1031 (0.2%)	1.03	1/1404 (0.1%)
2	D	1.03	2/1031 (0.2%)	0.96	0/1404
2	F	1.00	0/1031	0.96	0/1404
2	H	0.92	0/1040	0.91	0/1416
All	All	0.93	4/10744 (0.0%)	0.92	1/14522 (0.0%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	49	ALA	CA-C	5.97	1.59	1.52
2	D	97	ASN	CG-OD1	5.56	1.34	1.23
2	D	35	GLY	CA-C	-5.18	1.48	1.52
2	C	52	SER	CB-OG	-5.09	1.31	1.42

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	102	ARG	CG-CD-NE	-6.08	98.62	112.00

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1588	1610	1608	40	0
1	B	1594	1621	1621	48	0
1	E	1642	1664	1664	58	0
1	G	1642	1664	1664	49	0
2	C	1003	929	929	47	0
2	D	1003	929	929	40	0
2	F	1003	929	929	29	0
2	H	1012	938	938	45	0
3	C	2	0	0	0	0
3	D	2	0	0	0	0
3	F	1	0	0	0	0
3	H	2	0	0	0	0
All	All	10494	10284	10282	345	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 17.

All (345) 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)
2:H:40:ALA:HB3	2:H:43:LYS:HD3	1.47	0.97
2:D:23:ALA:HA	2:D:78:THR:HG22	1.52	0.92
2:F:27:ILE:HD11	2:F:32:PHE:CE1	2.04	0.92
2:F:11:LEU:HD23	2:F:125:THR:OG1	1.77	0.85
2:F:40:ALA:HB3	2:F:43:LYS:HD3	1.58	0.85
2:C:20:LEU:HD13	2:C:83:MET:HE3	1.59	0.84
2:C:96:ASP:OD2	2:C:102:ARG:NH1	2.11	0.83
1:B:300:GLU:O	1:B:332:GLN:NE2	2.16	0.78
1:B:355:GLY:O	1:B:357:GLY:N	2.17	0.78
1:A:355:GLY:O	1:A:357:GLY:N	2.18	0.76
1:A:177:SER:HB3	1:A:369:VAL:HG12	1.69	0.73
2:F:6:GLU:OE2	2:F:119:GLY:HA3	1.89	0.72
2:C:81:LEU:HD13	2:C:83:MET:HE2	1.72	0.71
2:D:128:SER:O	1:E:295:LYS:HB3	1.90	0.71
2:C:2:VAL:HG23	2:C:27:ILE:HD13	1.73	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:173:LEU:HD23	1:G:371:ALA:HB2	1.72	0.70
2:D:40:ALA:HB3	2:D:43:LYS:HD2	1.74	0.69
1:E:330:ILE:HG23	1:E:354:LEU:CD1	2.24	0.67
2:C:85:PRO:HA	2:C:126:VAL:HB	1.77	0.67
2:H:53:PRO:HA	2:H:72:ARG:HH22	1.60	0.67
2:D:6:GLU:O	2:D:7:SER:OG	2.10	0.67
1:G:334:VAL:HG22	1:G:339:GLN:HB2	1.76	0.66
2:C:23:ALA:HA	2:C:78:THR:HG22	1.76	0.66
1:B:269:PHE:CE1	1:B:327:LEU:HD13	2.30	0.65
1:A:162:LYS:HG3	1:A:245:GLU:OE2	1.96	0.65
2:D:102:ARG:NH1	2:D:107:SER:OG	2.29	0.64
1:A:165:LEU:HD22	1:A:239:GLY:HA2	1.80	0.64
2:D:6:GLU:OE2	2:D:119:GLY:HA3	1.98	0.64
1:G:355:GLY:O	1:G:357:GLY:N	2.30	0.64
2:C:9:GLY:HA2	2:C:18:LEU:HD21	1.80	0.63
1:G:215:LEU:HD13	1:G:216:GLU:N	2.12	0.63
1:B:179:GLU:O	1:B:183:LYS:HG2	1.99	0.63
1:G:281:LEU:HD11	1:G:308:ALA:CB	2.28	0.63
1:B:349:GLY:HA2	1:B:364:THR:HG23	1.81	0.63
2:H:17:SER:C	2:H:18:LEU:HD12	2.24	0.63
2:H:53:PRO:HA	2:H:72:ARG:NH2	2.15	0.62
1:A:334:VAL:HG22	1:A:339:GLN:HG3	1.81	0.62
2:D:24:VAL:CG1	2:D:27:ILE:HG22	2.29	0.61
2:H:67:ARG:NH2	2:H:87:ASP:OD2	2.22	0.61
2:F:34:ILE:HG21	2:F:79:VAL:HG11	1.82	0.61
2:D:88:THR:HG23	2:D:125:THR:HA	1.83	0.61
2:F:72:ARG:HH12	2:F:74:ASN:CG	2.09	0.61
1:G:281:LEU:HD12	1:G:281:LEU:N	2.15	0.61
1:E:243:GLY:HA2	1:E:246:GLU:OE1	2.01	0.60
1:E:281:LEU:HD11	1:E:308:ALA:CB	2.30	0.60
1:B:263:LEU:HD23	1:B:285:GLU:OE2	2.01	0.60
2:C:4:LEU:CD2	2:C:24:VAL:HG22	2.32	0.60
1:A:187:VAL:O	1:A:191:GLU:HG3	2.01	0.60
2:C:20:LEU:HD22	2:C:83:MET:CE	2.32	0.59
2:H:22:CYS:HB3	2:H:79:VAL:HG12	1.85	0.59
1:B:162:LYS:HG3	1:B:245:GLU:HG2	1.83	0.59
1:G:363:LEU:HD13	1:G:367:MET:HG3	1.85	0.59
2:H:85:PRO:HA	2:H:126:VAL:HB	1.85	0.59
1:G:173:LEU:HD11	1:G:181:LEU:CD2	2.32	0.58
1:E:294:ASP:C	1:E:296:PRO:HD3	2.27	0.58
2:D:37:PHE:O	2:D:91:TYR:HA	2.04	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:196:LYS:HD3	1:G:199:LYS:HD2	1.86	0.58
1:E:173:LEU:CD2	1:E:369:VAL:HG11	2.34	0.58
2:H:83:MET:HE1	2:H:91:TYR:CE2	2.39	0.58
2:C:64:ALA:HB1	2:C:68:PHE:CD2	2.39	0.57
1:B:171:TRP:CB	1:B:263:LEU:HD11	2.34	0.57
2:F:72:ARG:HH12	2:F:74:ASN:CB	2.17	0.57
1:A:333:GLN:HB3	1:A:354:LEU:HD23	1.85	0.57
1:A:281:LEU:HB2	1:A:286:LEU:HD12	1.86	0.57
2:C:52:SER:HB3	2:C:53:PRO:HD2	1.86	0.57
2:C:68:PHE:HA	2:C:82:GLN:O	2.05	0.57
1:G:173:LEU:CD1	1:G:181:LEU:HD23	2.35	0.57
1:G:216:GLU:OE1	1:G:216:GLU:HA	2.05	0.56
2:D:22:CYS:HB3	2:D:79:VAL:HG12	1.87	0.56
2:F:23:ALA:HA	2:F:78:THR:HG22	1.88	0.56
1:B:216:GLU:HG3	1:B:218:VAL:HG22	1.88	0.56
2:C:55:TYR:HD1	2:C:55:TYR:O	1.88	0.56
2:H:83:MET:HE1	2:H:91:TYR:CZ	2.41	0.56
2:C:64:ALA:HB1	2:C:68:PHE:CG	2.41	0.55
2:C:81:LEU:CD1	2:C:83:MET:HE2	2.34	0.55
1:A:281:LEU:CB	1:A:286:LEU:HD12	2.37	0.55
1:A:276:GLY:HA2	1:A:312:LEU:HG	1.88	0.55
1:E:274:THR:OG1	1:E:326:ASP:OD2	2.20	0.55
2:C:76:LYS:O	2:C:77:ASN:OD1	2.24	0.55
1:B:269:PHE:CD1	1:B:327:LEU:HD13	2.41	0.55
1:A:186:GLN:O	1:A:190:GLU:HG2	2.07	0.55
2:H:88:THR:HG23	2:H:125:THR:HA	1.88	0.55
2:H:55:TYR:HD2	2:H:56:GLU:OE2	1.90	0.54
2:F:96:ASP:OD2	2:F:99:LEU:HD23	2.07	0.54
2:H:27:ILE:HD11	2:H:32:PHE:CZ	2.43	0.54
1:B:334:VAL:HG23	1:B:339:GLN:HG3	1.89	0.54
2:D:24:VAL:O	2:D:77:ASN:ND2	2.40	0.54
1:E:179:GLU:OE1	1:E:183:LYS:CD	2.55	0.54
1:A:236:CYS:O	1:A:237:ASP:C	2.50	0.54
1:B:266:SER:OG	1:B:280:VAL:HG13	2.08	0.54
1:E:255:ARG:C	1:E:255:ARG:HD2	2.33	0.54
2:H:36:TRP:NE1	2:H:81:LEU:HB2	2.22	0.54
1:E:161:GLU:OE1	1:E:161:GLU:N	2.41	0.54
1:G:179:GLU:OE1	1:G:182:ARG:NH1	2.41	0.54
1:B:334:VAL:HG23	1:B:339:GLN:CG	2.37	0.53
2:C:45:ARG:HH12	2:H:86:TRP:HB2	1.73	0.53
1:E:291:SER:O	1:E:294:ASP:OD2	2.27	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:179:GLU:OE1	1:E:183:LYS:HD2	2.09	0.53
2:F:12:VAL:O	2:F:126:VAL:HA	2.08	0.53
1:A:342:ALA:HA	1:A:352:TYR:CD1	2.44	0.53
1:B:350:LYS:O	1:B:361:LEU:HD12	2.07	0.53
2:C:26:GLY:O	2:C:27:ILE:HD12	2.09	0.53
2:C:64:ALA:CB	2:C:68:PHE:CD2	2.91	0.53
2:D:67:ARG:HB3	2:D:83(A):ASN:O	2.09	0.53
2:F:52:SER:HB2	2:F:99:LEU:HB3	1.91	0.53
2:D:27:ILE:HD11	2:D:32:PHE:CZ	2.44	0.53
2:C:51:ILE:HG13	2:C:58:ILE:HG12	1.89	0.53
1:B:329:ASP:OD1	1:B:358:ARG:NH1	2.42	0.52
2:D:28:THR:CG2	2:D:31:ASP:OD2	2.57	0.52
1:G:173:LEU:HD22	1:G:369:VAL:HG12	1.91	0.52
1:G:173:LEU:HD13	1:G:369:VAL:HG11	1.90	0.52
1:E:355:GLY:O	1:E:357:GLY:N	2.39	0.52
1:B:177:SER:HB3	1:B:369:VAL:HG13	1.91	0.52
1:B:162:LYS:CG	1:B:245:GLU:HG2	2.39	0.52
2:H:6:GLU:OE2	2:H:119:GLY:HA3	2.09	0.52
1:E:273:LYS:HE2	1:E:326:ASP:OD1	2.10	0.52
2:H:83(B):ASN:O	2:H:84:LYS:NZ	2.43	0.52
1:A:333:GLN:CB	1:A:354:LEU:HD23	2.40	0.52
1:B:337:GLY:C	1:B:339:GLN:H	2.18	0.51
2:D:27:ILE:HG13	2:D:32:PHE:CD1	2.46	0.51
2:F:102:ARG:NH1	2:F:107:SER:OG	2.42	0.51
2:D:40:ALA:HB1	2:D:41:PRO:HD2	1.91	0.51
2:C:60:TYR:OH	2:C:69:THR:HA	2.11	0.51
2:H:20:LEU:HB3	2:H:36:TRP:CH2	2.46	0.51
1:A:342:ALA:HB2	1:A:352:TYR:HE1	1.76	0.51
1:E:182:ARG:O	1:E:186:GLN:OE1	2.28	0.51
2:D:64:ALA:CB	2:D:68:PHE:CD2	2.94	0.50
1:G:217:LEU:N	1:G:217:LEU:HD12	2.25	0.50
1:E:266:SER:OG	1:E:280:VAL:HG23	2.10	0.50
1:B:331:LEU:HA	1:B:334:VAL:HG12	1.92	0.50
1:E:173:LEU:HD21	1:E:369:VAL:HG11	1.94	0.50
1:G:252:VAL:HG13	1:G:288:LEU:HD23	1.94	0.50
2:H:64:ALA:HB1	2:H:68:PHE:HB2	1.94	0.50
2:H:72:ARG:HD2	2:H:74:ASN:CG	2.37	0.50
2:F:24:VAL:CG1	2:F:27:ILE:HG22	2.42	0.50
2:H:27:ILE:HD11	2:H:32:PHE:CE1	2.47	0.49
1:A:179:GLU:OE1	1:A:180:THR:N	2.45	0.49
1:A:267:ALA:C	1:A:268:LEU:HD12	2.37	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:196:LYS:O	1:E:200:LYS:HG2	2.12	0.49
2:C:45:ARG:NH1	2:H:86:TRP:HB2	2.26	0.49
2:D:96:ASP:HB2	2:D:115:TYR:CD2	2.48	0.49
2:D:102:ARG:HB2	2:D:107:SER:HB3	1.95	0.49
1:E:206:ILE:HG12	1:E:207:SER:N	2.28	0.49
2:C:76:LYS:O	2:C:78:THR:HG23	2.13	0.49
1:E:309:HIS:HB3	1:E:327:LEU:HD22	1.93	0.49
2:D:64:ALA:O	2:D:65:LYS:C	2.56	0.48
2:F:27:ILE:HD11	2:F:32:PHE:CZ	2.47	0.48
2:H:73:ASP:OD2	2:H:75:ALA:HB3	2.12	0.48
1:B:218:VAL:O	1:B:224:ARG:NH2	2.47	0.48
2:D:94:ALA:HB2	2:D:118:TRP:CE3	2.48	0.48
1:G:330:ILE:HG23	1:G:354:LEU:HD22	1.95	0.48
2:H:24:VAL:CG1	2:H:27:ILE:HG22	2.44	0.48
1:B:282:THR:O	1:B:286:LEU:HD12	2.13	0.48
1:G:201:GLU:HB3	1:G:204:HIS:NE2	2.28	0.48
2:H:25:SER:C	2:H:27:ILE:H	2.21	0.48
1:A:265:ILE:HB	1:A:367:MET:HB3	1.96	0.48
2:C:2:VAL:HG21	2:C:117:TYR:CE1	2.48	0.48
2:D:2:VAL:HA	2:D:26:GLY:HA3	1.95	0.48
1:E:295:LYS:N	1:E:296:PRO:HD3	2.28	0.48
2:F:13:GLN:HG2	2:F:127:SER:OG	2.14	0.48
1:G:281:LEU:HD11	1:G:308:ALA:HB1	1.96	0.48
2:H:27:ILE:CG1	2:H:32:PHE:CE1	2.96	0.48
1:A:162:LYS:HA	1:A:165:LEU:HD13	1.95	0.47
1:B:215:LEU:C	1:B:215:LEU:HD13	2.39	0.47
1:B:177:SER:HB3	1:B:369:VAL:CG1	2.44	0.47
1:B:297:SER:O	1:B:300:GLU:HG2	2.14	0.47
1:E:266:SER:OG	1:E:280:VAL:CG2	2.63	0.47
2:H:27:ILE:HG13	2:H:32:PHE:CD1	2.50	0.47
1:B:320:PRO:HB2	2:D:103:LEU:O	2.14	0.47
1:A:266:SER:OG	1:A:280:VAL:CG1	2.63	0.47
2:C:20:LEU:HD22	2:C:83:MET:HE1	1.97	0.47
1:B:276:GLY:HA2	1:B:312:LEU:HG	1.96	0.47
2:D:92:TYR:CE1	2:D:121:GLY:HA3	2.49	0.47
2:F:20:LEU:HD13	2:F:83:MET:CE	2.44	0.47
2:C:55:TYR:O	2:C:55:TYR:CD1	2.67	0.47
1:A:176:LYS:O	1:A:179:GLU:HG3	2.15	0.47
1:E:327:LEU:O	1:E:331:LEU:HG	2.14	0.47
1:A:193:GLY:HA3	1:A:217:LEU:HD13	1.97	0.47
1:A:320:PRO:HG2	2:C:103:LEU:HB3	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:162:LYS:HG3	1:B:245:GLU:CG	2.45	0.47
1:E:337:GLY:C	1:E:339:GLN:H	2.23	0.47
1:G:162:LYS:HA	1:G:165:LEU:HD13	1.97	0.47
1:G:289:TRP:CE2	1:G:305:GLY:HA2	2.50	0.47
1:B:162:LYS:HE2	1:B:239:GLY:C	2.40	0.46
1:G:281:LEU:N	1:G:281:LEU:CD1	2.78	0.46
1:A:317:ASP:N	1:A:317:ASP:OD1	2.41	0.46
2:D:86:TRP:HH2	1:E:300:GLU:HB3	1.80	0.46
2:F:72:ARG:NH2	2:F:74:ASN:OD1	2.49	0.46
1:E:255:ARG:HD2	1:E:255:ARG:O	2.15	0.46
2:F:33:GLY:CA	2:F:53:PRO:HD3	2.45	0.46
2:H:50:CYS:HB2	2:H:109:TRP:CE3	2.50	0.46
1:G:173:LEU:CD1	1:G:369:VAL:HG11	2.45	0.46
1:B:333:GLN:NE2	1:B:358:ARG:HH22	2.13	0.46
2:C:87:ASP:O	2:C:88:THR:C	2.58	0.46
1:B:186:GLN:O	1:B:190:GLU:HG3	2.15	0.46
2:D:53:PRO:O	2:D:55:TYR:N	2.49	0.46
2:D:60:TYR:OH	2:D:69:THR:HA	2.16	0.46
1:E:179:GLU:HA	1:E:182:ARG:HB3	1.97	0.46
2:H:83(B):ASN:O	2:H:84:LYS:HD2	2.16	0.46
1:G:269:PHE:CE1	1:G:327:LEU:HD13	2.51	0.46
1:A:181:LEU:HD13	1:A:367:MET:CE	2.45	0.46
1:A:281:LEU:N	1:A:281:LEU:HD12	2.31	0.46
2:D:37:PHE:O	2:D:92:TYR:N	2.45	0.46
2:F:120:GLN:NE2	2:F:121:GLY:O	2.48	0.46
1:G:173:LEU:HD11	1:G:181:LEU:HD23	1.97	0.46
2:F:22:CYS:HB3	2:F:79:VAL:HG12	1.98	0.46
1:A:221:PHE:HB3	1:A:313:GLY:HA3	1.98	0.45
1:E:295:LYS:N	1:E:296:PRO:CD	2.78	0.45
2:C:73:ASP:OD2	2:C:76:LYS:HG3	2.16	0.45
2:H:51:ILE:HG23	2:H:51:ILE:O	2.17	0.45
1:E:263:LEU:HD23	1:E:285:GLU:OE2	2.17	0.45
1:E:366:LYS:N	1:E:366:LYS:HD3	2.31	0.45
1:B:238:TYR:OH	2:D:98:ASP:OD1	2.23	0.45
2:D:28:THR:HG21	2:D:31:ASP:OD2	2.15	0.45
2:F:32:PHE:CZ	2:F:97:ASN:HB2	2.52	0.45
2:C:38:ARG:HD3	2:C:91:TYR:CZ	2.51	0.45
1:E:271:THR:HG21	1:E:326:ASP:HB3	1.98	0.45
1:B:195:HIS:CE1	1:B:344:GLY:HA3	2.52	0.45
2:D:40:ALA:O	2:D:41:PRO:C	2.60	0.45
1:E:236:CYS:O	1:E:237:ASP:C	2.60	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:330:ILE:HG23	1:E:354:LEU:HD11	1.98	0.45
1:A:343:VAL:HG22	1:A:343:VAL:O	2.17	0.45
1:G:173:LEU:CD1	1:G:181:LEU:CD2	2.94	0.45
1:A:221:PHE:CE2	1:A:275:ALA:HB2	2.52	0.45
1:A:273:LYS:HE3	1:A:322:GLN:NE2	2.32	0.45
1:A:342:ALA:HA	1:A:352:TYR:HD1	1.81	0.45
1:E:203:ARG:HG3	1:E:204:HIS:N	2.31	0.45
2:C:111:ASP:OD1	2:C:111:ASP:C	2.60	0.44
2:H:2:VAL:HG13	2:H:2:VAL:O	2.17	0.44
2:C:26:GLY:C	2:C:27:ILE:HD12	2.43	0.44
2:D:111:ASP:OD1	2:D:111:ASP:C	2.59	0.44
2:H:40:ALA:HB1	2:H:41:PRO:HD2	1.98	0.44
2:C:38:ARG:NH1	2:C:46:GLU:OE2	2.51	0.44
2:C:45:ARG:NH1	2:H:86:TRP:CG	2.86	0.44
2:H:29:LEU:HB3	2:H:74:ASN:ND2	2.33	0.44
1:B:216:GLU:HG3	1:B:218:VAL:CG2	2.48	0.44
1:E:283:ASP:HA	1:E:286:LEU:HD12	2.00	0.44
2:H:3:GLN:HB2	2:H:25:SER:OG	2.17	0.44
1:A:237:ASP:OD2	1:A:240:LYS:CD	2.66	0.44
2:H:111:ASP:C	2:H:111:ASP:OD1	2.58	0.44
1:E:189:LEU:HB3	1:E:217:LEU:HB3	2.00	0.44
2:F:88:THR:HG23	2:F:125:THR:HA	2.00	0.44
1:G:177:SER:HB3	1:G:369:VAL:HG13	1.99	0.44
1:B:181:LEU:HD23	1:B:181:LEU:O	2.17	0.43
1:E:376:TYR:CD1	1:E:376:TYR:C	2.96	0.43
1:G:259:LYS:HA	1:G:259:LYS:HE2	2.00	0.43
1:A:268:LEU:HD13	1:A:363:LEU:HD21	1.99	0.43
1:B:346:LEU:HD22	1:B:361:LEU:HD22	2.01	0.43
2:D:64:ALA:HB1	2:D:68:PHE:CG	2.54	0.43
2:F:24:VAL:CG1	2:F:27:ILE:CG2	2.97	0.43
2:F:32:PHE:CE2	2:F:97:ASN:HB2	2.52	0.43
1:G:191:GLU:OE1	1:G:346:LEU:HD23	2.18	0.43
2:H:76:LYS:O	2:H:77:ASN:C	2.62	0.43
1:E:307:ARG:C	1:E:327:LEU:HD23	2.44	0.43
1:E:307:ARG:O	1:E:327:LEU:HD23	2.19	0.43
2:D:100:PRO:O	2:D:101:ASP:C	2.62	0.43
1:E:289:TRP:NE1	1:E:305:GLY:HA2	2.33	0.43
1:G:173:LEU:HD12	1:G:181:LEU:HD23	2.01	0.43
1:G:267:ALA:C	1:G:268:LEU:HD12	2.44	0.43
2:D:34:ILE:O	2:D:50:CYS:HA	2.19	0.43
1:B:171:TRP:HB2	1:B:263:LEU:HD11	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:206:ILE:HG13	1:B:207:SER:N	2.34	0.43
1:B:330:ILE:HG23	1:B:354:LEU:CD1	2.49	0.43
2:F:55:TYR:C	2:F:57:HIS:H	2.27	0.43
1:G:268:LEU:HD12	1:G:268:LEU:N	2.33	0.43
1:B:334:VAL:HG13	1:B:335:LYS:N	2.32	0.43
2:D:68:PHE:N	2:D:68:PHE:CD1	2.85	0.43
1:E:198:PHE:CE2	1:E:202:LEU:HD21	2.53	0.43
1:E:291:SER:HA	1:E:294:ASP:OD2	2.19	0.43
2:H:33:GLY:HA3	2:H:99:LEU:HD23	2.01	0.43
1:E:197:ALA:HB1	1:E:343:VAL:CG1	2.49	0.42
1:G:237:ASP:OD2	1:G:240:LYS:HG3	2.19	0.42
2:H:27:ILE:HG13	2:H:32:PHE:CE1	2.54	0.42
2:D:12:VAL:O	2:D:126:VAL:HA	2.19	0.42
1:A:252:VAL:O	1:A:256:SER:OG	2.27	0.42
2:C:84:LYS:O	2:C:126:VAL:HG21	2.19	0.42
2:C:100:PRO:O	2:C:101:ASP:C	2.62	0.42
2:H:40:ALA:HB1	2:H:41:PRO:CD	2.49	0.42
1:A:179:GLU:OE1	1:A:179:GLU:C	2.62	0.42
1:E:329:ASP:OD1	1:E:358:ARG:NH1	2.53	0.42
1:G:237:ASP:OD2	1:G:240:LYS:CG	2.68	0.42
1:G:271:THR:HG21	1:G:326:ASP:HB3	2.02	0.42
1:A:327:LEU:O	1:A:331:LEU:HG	2.20	0.42
2:C:38:ARG:NH2	2:C:46:GLU:OE2	2.53	0.42
1:E:177:SER:O	1:E:178:SER:C	2.62	0.42
1:E:252:VAL:HG11	1:E:287:GLN:HB3	2.02	0.42
2:C:86:TRP:CE3	2:C:86:TRP:N	2.88	0.42
1:E:319:GLN:HA	1:E:320:PRO:HD3	1.87	0.42
1:B:325:LEU:N	1:B:325:LEU:CD1	2.83	0.42
1:G:195:HIS:NE2	1:G:344:GLY:HA3	2.35	0.42
1:A:206:ILE:HG13	1:A:207:SER:N	2.35	0.42
1:B:325:LEU:N	1:B:325:LEU:HD12	2.35	0.42
2:F:25:SER:C	2:F:27:ILE:H	2.28	0.42
2:C:67:ARG:HH22	2:C:87:ASP:CG	2.28	0.41
2:C:102:ARG:O	2:C:103:LEU:C	2.61	0.41
2:C:12:VAL:O	2:C:126:VAL:HA	2.20	0.41
2:C:17:SER:HB2	2:D:15:GLY:O	2.20	0.41
2:C:84:LYS:C	2:C:126:VAL:HG11	2.45	0.41
1:E:173:LEU:HD23	1:E:369:VAL:HG11	2.03	0.41
2:F:29:LEU:CD2	2:F:72:ARG:HD2	2.50	0.41
1:A:281:LEU:HB2	1:A:286:LEU:CD1	2.50	0.41
1:E:179:GLU:OE1	1:E:183:LYS:HG3	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:366:LYS:HD3	1:E:366:LYS:H	1.84	0.41
1:G:263:LEU:HD23	1:G:285:GLU:OE2	2.20	0.41
1:G:342:ALA:HB1	1:G:350:LYS:HG2	2.01	0.41
1:G:366:LYS:HD3	1:G:366:LYS:N	2.35	0.41
1:G:374:THR:OG1	1:G:375:GLY:N	2.52	0.41
2:H:29:LEU:HD12	2:H:29:LEU:HA	1.85	0.41
1:E:164:PHE:N	1:E:164:PHE:CD1	2.88	0.41
1:B:353:SER:HB2	1:B:359:TRP:CZ3	2.55	0.41
2:H:55:TYR:CD2	2:H:56:GLU:OE2	2.71	0.41
2:C:96:ASP:HA	2:C:116:ASP:OD2	2.20	0.41
1:B:329:ASP:OD1	1:B:329:ASP:C	2.64	0.41
1:E:166:PRO:O	1:E:236:CYS:HB2	2.20	0.41
1:E:256:SER:OG	1:E:288:LEU:HD21	2.21	0.41
1:G:215:LEU:HD13	1:G:215:LEU:C	2.45	0.41
1:A:337:GLY:C	1:A:339:GLN:H	2.29	0.41
1:B:162:LYS:HG3	1:B:245:GLU:CD	2.45	0.41
1:B:173:LEU:HD23	1:B:369:VAL:HG11	2.03	0.41
1:B:349:GLY:C	1:B:361:LEU:HD11	2.46	0.41
1:E:319:GLN:HE22	2:F:108:ASP:CG	2.29	0.41
2:H:17:SER:O	2:H:18:LEU:HD12	2.20	0.41
2:D:41:PRO:O	2:D:43:LYS:N	2.53	0.41
1:E:235:PHE:CE2	1:E:237:ASP:HB2	2.55	0.41
1:B:349:GLY:O	1:B:361:LEU:HD11	2.20	0.40
2:F:13:GLN:HB3	2:F:14:PRO:CD	2.51	0.40
1:G:317:ASP:N	1:G:317:ASP:OD1	2.53	0.40
1:A:330:ILE:O	1:A:333:GLN:HB2	2.21	0.40
2:C:37:PHE:CE2	2:C:112:PRO:HB3	2.57	0.40
1:B:327:LEU:O	1:B:331:LEU:HG	2.21	0.40
1:E:265:ILE:HB	1:E:367:MET:HB2	2.02	0.40
1:G:165:LEU:HD23	1:G:236:CYS:HB3	2.02	0.40
1:G:173:LEU:HD11	1:G:181:LEU:HD22	2.01	0.40
1:G:184:ALA:HB1	1:G:348:ARG:HH12	1.86	0.40
1:G:198:PHE:HA	1:G:359:TRP:CH2	2.56	0.40
2:H:67:ARG:H	2:H:67:ARG:HG2	1.70	0.40
2:C:38:ARG:HD3	2:C:91:TYR:OH	2.22	0.40
1:E:166:PRO:HB3	1:E:377:TYR:CZ	2.57	0.40
2:C:103:LEU:N	2:C:103:LEU:CD1	2.85	0.40
2:D:128:SER:O	1:E:295:LYS:CB	2.66	0.40
1:G:234:LYS:HG3	1:G:235:PHE:O	2.21	0.40
1:G:289:TRP:CD2	1:G:290:PRO:HD2	2.56	0.40
1:G:342:ALA:HA	1:G:352:TYR:HD1	1.87	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:64:ALA:CB	2:H:68:PHE:CD2	3.04	0.40
2:H:68:PHE:N	2:H:68:PHE:CD1	2.90	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	198/220 (90%)	188 (95%)	8 (4%)	2 (1%)	12	17
1	B	199/220 (90%)	187 (94%)	11 (6%)	1 (0%)	24	34
1	E	207/220 (94%)	191 (92%)	11 (5%)	5 (2%)	4	4
1	G	207/220 (94%)	197 (95%)	9 (4%)	1 (0%)	24	34
2	C	128/131 (98%)	119 (93%)	7 (6%)	2 (2%)	7	9
2	D	128/131 (98%)	119 (93%)	7 (6%)	2 (2%)	7	9
2	F	128/131 (98%)	116 (91%)	10 (8%)	2 (2%)	7	9
2	H	129/131 (98%)	118 (92%)	9 (7%)	2 (2%)	7	9
All	All	1324/1404 (94%)	1235 (93%)	72 (5%)	17 (1%)	9	12

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	356	LYS
2	D	65	LYS
1	G	356	LYS
2	H	2	VAL
2	H	65	LYS
1	B	356	LYS
2	D	54	GLY
1	E	292	ASP

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Mol	Chain	Res	Type
1	E	296	PRO
2	F	55	TYR
2	C	55	TYR
1	E	294	ASP
2	F	56	GLU
2	C	127	SER
1	E	356	LYS
1	A	340	GLY
1	E	340	GLY

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	166/180 (92%)	153 (92%)	13 (8%)	11	16
1	B	167/180 (93%)	160 (96%)	7 (4%)	26	40
1	E	173/180 (96%)	165 (95%)	8 (5%)	24	36
1	G	173/180 (96%)	160 (92%)	13 (8%)	12	17
2	C	105/106 (99%)	96 (91%)	9 (9%)	10	13
2	D	105/106 (99%)	98 (93%)	7 (7%)	15	21
2	F	105/106 (99%)	98 (93%)	7 (7%)	15	21
2	H	106/106 (100%)	101 (95%)	5 (5%)	23	36
All	All	1100/1144 (96%)	1031 (94%)	69 (6%)	16	23

All (69) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	175	LYS
1	A	204	HIS
1	A	216	GLU
1	A	223	LYS
1	A	240	LYS
1	A	266	SER

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Mol	Chain	Res	Type
1	A	274	THR
1	A	284	GLN
1	A	286	LEU
1	A	325	LEU
1	A	356	LYS
1	A	368	GLU
1	A	369	VAL
2	C	27	ILE
2	C	67	ARG
2	C	69	THR
2	C	81	LEU
2	C	84	LYS
2	C	86	TRP
2	C	99	LEU
2	C	103	LEU
2	C	120	GLN
1	B	181	LEU
1	B	223	LYS
1	B	233	THR
1	B	240	LYS
1	B	286	LEU
1	B	325	LEU
1	B	343	VAL
2	D	2	VAL
2	D	11	LEU
2	D	44	GLU
2	D	55	TYR
2	D	65	LYS
2	D	86	TRP
2	D	125	THR
1	E	161	GLU
1	E	162	LYS
1	E	204	HIS
1	E	206	ILE
1	E	217	LEU
1	E	343	VAL
1	E	356	LYS
1	E	364	THR
2	F	29	LEU
2	F	52	SER
2	F	56	GLU
2	F	72	ARG

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Mol	Chain	Res	Type
2	F	74	ASN
2	F	99	LEU
2	F	113	SER
1	G	162	LYS
1	G	177	SER
1	G	202	LEU
1	G	204	HIS
1	G	229	LEU
1	G	256	SER
1	G	280	VAL
1	G	307	ARG
1	G	334	VAL
1	G	339	GLN
1	G	346	LEU
1	G	350	LYS
1	G	351	LEU
2	H	12	VAL
2	H	52	SER
2	H	56	GLU
2	H	81	LEU
2	H	99	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	284	GLN
1	A	287	GLN
1	A	322	GLN
2	C	82	GLN
1	B	194	ASN
2	D	74	ASN
1	E	278	GLN
2	F	39	GLN
2	F	120	GLN
1	G	284	GLN

5.3.3 RNA ⓘ

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 7 ligands modelled in this entry, 7 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	204/220 (92%)	0.47	7 (3%) 48 49	73, 103, 148, 185	0
1	B	205/220 (93%)	0.32	5 (2%) 59 61	79, 121, 185, 214	0
1	E	211/220 (95%)	0.27	4 (1%) 66 66	74, 112, 150, 181	0
1	G	211/220 (95%)	0.20	3 (1%) 73 74	77, 121, 178, 216	0
2	C	130/131 (99%)	0.25	1 (0%) 82 84	69, 94, 136, 177	0
2	D	130/131 (99%)	0.24	3 (2%) 61 62	69, 96, 123, 136	0
2	F	130/131 (99%)	0.25	1 (0%) 82 84	70, 110, 147, 169	0
2	H	131/131 (100%)	0.21	0 100 100	70, 109, 153, 166	0
All	All	1352/1404 (96%)	0.28	24 (1%) 67 68	69, 108, 164, 216	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	248	ALA	3.4
1	E	336	GLY	3.2
1	B	377	TYR	2.9
1	A	298	ALA	2.8
1	A	207	SER	2.7
1	A	217	LEU	2.7
1	A	299	SER	2.6
1	B	296	PRO	2.6
1	B	298	ALA	2.5
2	F	103	LEU	2.5
1	B	253	VAL	2.4
2	D	55	TYR	2.4
2	D	81	LEU	2.3
1	A	169	PHE	2.3
1	E	265	ILE	2.3
1	E	298	ALA	2.2

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Mol	Chain	Res	Type	RSRZ
1	G	252	VAL	2.1
1	B	169	PHE	2.1
1	G	192	LEU	2.1
2	D	12	VAL	2.0
2	C	127	SER	2.0
1	A	192	LEU	2.0
1	A	263	LEU	2.0
1	E	171	TRP	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 oligosaccharides 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. 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	B-factors($\text{\AA}^2$)	Q<0.9
3	CA	H	201	1/1	0.82	0.13	156,156,156,156	0
3	CA	D	201	1/1	0.85	0.15	80,80,80,80	0
3	CA	D	202	1/1	0.87	0.12	88,88,88,88	0
3	CA	C	202	1/1	0.88	0.11	73,73,73,73	0
3	CA	F	201	1/1	0.89	0.10	76,76,76,76	0
3	CA	H	202	1/1	0.90	0.12	95,95,95,95	0
3	CA	C	201	1/1	0.94	0.07	88,88,88,88	0

6.5 Other polymers [i](#)

There are no such residues in this entry.