



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:37 AM UTC

PDB ID : 1LO5 / pdb_00001lo5
Title : Crystal structure of the D227A variant of Staphylococcal enterotoxin A in complex with human MHC class II
Authors : Petersson, K.; Thunnissen, M.; Forsberg, G.; Walse, B.
Deposited on : 2002-05-06
Resolution : 3.20 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

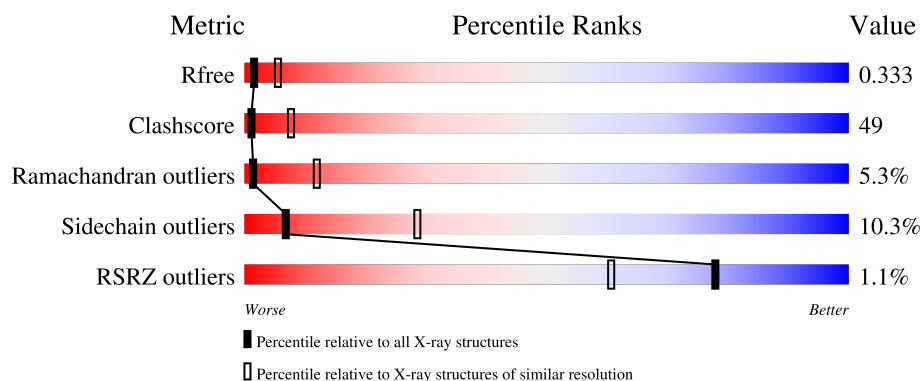
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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	1466 (3.20-3.20)
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (3.20-3.20)

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	182	31% 52% 15% .
2	B	190	3% 27% 56% 14% ..
3	C	13	38% 62%
4	D	233	% 30% 53% 15% .

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 5055 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 HLA class II histocompatibility antigen, DR alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	179	Total	C	N	O	S	0	0	0
			1470	952	239	274	5			

- Molecule 2 is a protein called HLA class II histocompatibility antigen, DR-1 beta chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	188	Total	C	N	O	S	0	0	0
			1545	973	277	289	6			

- Molecule 3 is a protein called Hemagglutinin peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	13	Total	C	N	O	0	0	0
			106	69	18	19			

- Molecule 4 is a protein called enterotoxin A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	233	Total	C	N	O	S	0	0	0
			1911	1210	323	374	4			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	227	ALA	ASP	engineered mutation	UNP P0A0L2

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	8	Total	O	0	0
			8	8		

Continued on next page...

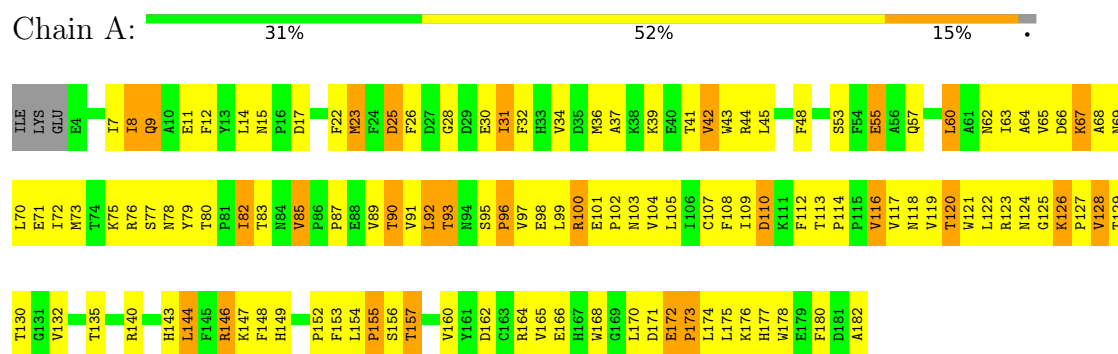
Continued from previous page...

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	8	Total	O	0	0
			8	8		
5	D	7	Total	O	0	0
			7	7		

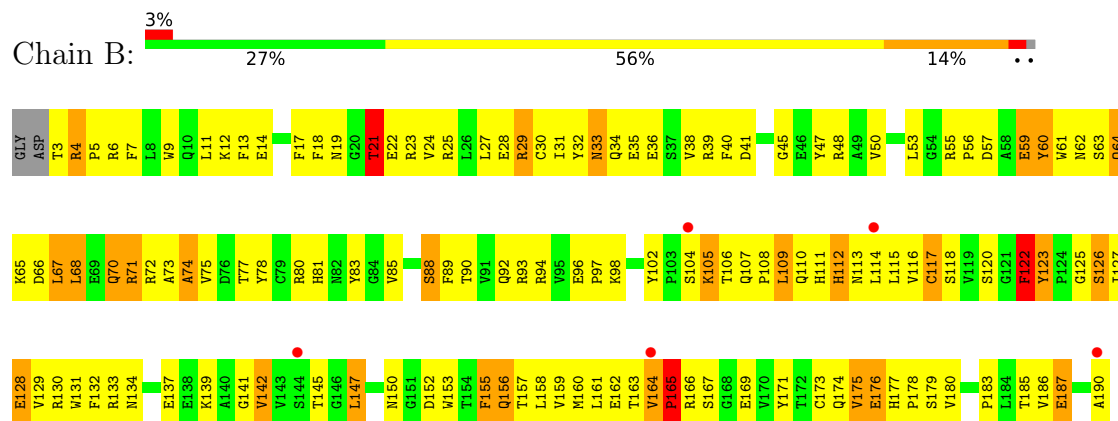
3 Residue-property plots

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: HLA class II histocompatibility antigen, DR alpha chain



- Molecule 2: HLA class II histocompatibility antigen, DR-1 beta chain

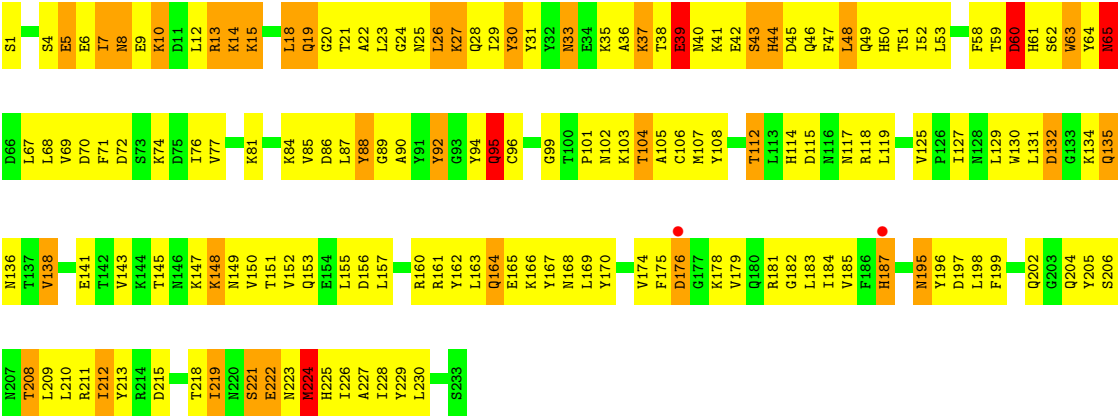


- Molecule 3: Hemagglutinin peptide



- Molecule 4: enterotoxin A





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	52.01Å 75.79Å 198.08Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 3.20 30.00 – 3.20	Depositor EDS
% Data completeness (in resolution range)	93.9 (30.00-3.20) 90.4 (30.00-3.20)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.05 (at 3.23Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.245 , 0.339 0.243 , 0.333	Depositor DCC
R_{free} test set	1005 reflections (8.17%)	wwPDB-VP
Wilson B-factor (Å ²)	48.3	Xtriage
Anisotropy	0.490	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 70.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.41$, $\langle L^2 \rangle = 0.23$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.86	EDS
Total number of atoms	5055	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.07% 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

5.1 Standard geometry

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	0.59	0/1515	1.11	14/2065 (0.7%)
2	B	0.53	0/1585	1.05	9/2152 (0.4%)
3	C	0.60	0/107	1.01	0/141
4	D	0.51	0/1950	1.19	26/2630 (1.0%)
All	All	0.54	0/5157	1.12	49/6988 (0.7%)

There are no bond length outliers.

All (49) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	157	THR	N-CA-C	-10.50	99.17	114.39
1	A	31	ILE	N-CA-C	-9.07	103.02	111.45
4	D	15	LYS	N-CA-C	-8.94	101.81	112.89
4	D	30	TYR	N-CA-C	8.37	121.98	111.69
2	B	122	PHE	N-CA-C	8.36	128.60	110.80
4	D	221	SER	N-CA-C	8.18	119.88	110.97
2	B	139	LYS	N-CA-C	-7.95	102.63	112.88
4	D	7	ILE	N-CA-C	-7.91	101.51	112.50
4	D	74	LYS	N-CA-C	-7.83	102.75	111.28
4	D	187	HIS	N-CA-C	6.96	119.62	108.41
4	D	44	HIS	N-CA-C	-6.90	102.05	110.88
4	D	26	LEU	N-CA-C	-6.85	103.59	112.23
4	D	5	GLU	N-CA-C	-6.79	104.69	114.12
2	B	164	VAL	CA-C-N	6.48	127.94	119.84
2	B	164	VAL	C-N-CA	6.48	127.94	119.84
4	D	39	GLU	N-CA-C	6.40	118.56	110.24
4	D	49	GLN	N-CA-C	6.40	120.92	113.18
1	A	126	LYS	CA-C-N	6.39	126.34	120.21
1	A	126	LYS	C-N-CA	6.39	126.34	120.21
4	D	148	LYS	N-CA-C	-6.04	103.44	111.24
4	D	27	LYS	N-CA-C	-5.98	105.48	112.89
4	D	222	GLU	N-CA-C	5.86	118.11	110.43
4	D	92	TYR	N-CA-C	5.78	117.04	108.60

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	224	MET	N-CA-C	5.75	117.33	110.19
1	A	85	VAL	CA-C-N	5.71	126.26	120.38
1	A	85	VAL	C-N-CA	5.71	126.26	120.38
4	D	138	VAL	CA-C-N	5.68	126.94	119.84
4	D	138	VAL	C-N-CA	5.68	126.94	119.84
4	D	8	ASN	N-CA-C	-5.67	106.20	113.01
1	A	144	LEU	N-CA-C	-5.55	101.42	109.31
4	D	48	LEU	N-CA-C	-5.55	101.25	109.86
1	A	110	ASP	N-CA-C	5.53	117.20	108.96
4	D	104	THR	N-CA-C	5.51	118.45	109.24
1	A	165	VAL	N-CA-C	5.47	115.98	107.28
4	D	13	ARG	N-CA-C	-5.42	103.74	110.41
4	D	38	THR	N-CA-C	5.35	117.77	107.44
1	A	8	ILE	CB-CA-C	-5.34	103.21	110.42
4	D	215	ASP	N-CA-C	-5.31	106.75	113.18
2	B	88	SER	N-CA-C	5.28	116.72	111.07
1	A	67	LYS	N-CA-C	-5.24	104.82	111.11
2	B	117	CYS	N-CA-C	-5.13	102.10	110.20
2	B	175	VAL	N-CA-C	5.11	116.03	107.18
1	A	9	GLN	N-CA-C	-5.10	98.26	107.75
1	A	128	VAL	N-CA-C	5.09	116.33	108.44
4	D	18	LEU	N-CA-C	5.09	116.61	108.41
2	B	14	GLU	N-CA-C	5.08	117.76	109.59
4	D	96	CYS	N-CA-C	5.06	117.81	110.42
1	A	42	VAL	N-CA-C	5.06	114.03	106.85
2	B	165	PRO	N-CA-C	5.04	122.84	112.47

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	1470	0	1406	139	0
2	B	1545	0	1478	189	0
3	C	106	0	119	14	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	D	1911	0	1863	177	0
5	A	8	0	0	0	0
5	B	8	0	0	1	0
5	D	7	0	0	0	0
All	All	5055	0	4866	482	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 49.

All (482) 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:B:110:GLN:H	2:B:165:PRO:HG2	1.06	1.09
2:B:129:VAL:HG22	2:B:175:VAL:HG22	1.32	1.07
4:D:52:ILE:HD11	4:D:77:VAL:HG13	1.40	1.04
1:A:82:ILE:HD13	2:B:33:ASN:HB3	1.40	0.98
2:B:110:GLN:N	2:B:165:PRO:HG2	1.79	0.95
4:D:44:HIS:HA	4:D:81:LYS:HG3	1.50	0.93
4:D:221:SER:HA	4:D:224:MET:HE2	1.49	0.93
1:A:39:LYS:HG2	1:A:60:LEU:HD11	1.53	0.90
2:B:176:GLU:HG3	2:B:183:PRO:HB3	1.54	0.88
4:D:13:ARG:HH12	4:D:202:GLN:NE2	1.71	0.88
2:B:187:GLU:CD	2:B:187:GLU:H	1.82	0.87
4:D:127:ILE:HD11	4:D:143:VAL:HG23	1.55	0.87
4:D:51:THR:C	4:D:52:ILE:HD12	2.01	0.85
4:D:95:GLN:HA	4:D:95:GLN:HE21	1.40	0.84
4:D:37:LYS:O	4:D:37:LYS:HD3	1.77	0.84
2:B:116:VAL:HG13	2:B:160:MET:HG2	1.60	0.84
1:A:160:VAL:HB	1:A:177:HIS:CE1	2.12	0.83
1:A:108:PHE:HZ	1:A:146:ARG:HD2	1.44	0.83
4:D:127:ILE:HD11	4:D:143:VAL:CG2	2.07	0.83
4:D:183:LEU:HD21	4:D:195:ASN:ND2	1.93	0.83
2:B:3:THR:HG22	2:B:4:ARG:HD3	1.61	0.83
4:D:13:ARG:HH12	4:D:202:GLN:HE22	1.28	0.81
2:B:25:ARG:NH2	2:B:27:LEU:HD22	1.97	0.80
2:B:141:GLY:O	2:B:161:LEU:HA	1.82	0.80
1:A:36:MET:HB3	4:D:48:LEU:HD13	1.64	0.79
2:B:122:PHE:HZ	2:B:127:ILE:HG12	1.48	0.79
2:B:132:PHE:CE1	2:B:137:GLU:HB2	2.18	0.79
1:A:87:PRO:HB3	1:A:112:PHE:HB3	1.65	0.78
1:A:26:PHE:CD2	2:B:90:THR:HB	2.17	0.78

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:97:VAL:HA	1:A:103:ASN:ND2	1.99	0.78
4:D:52:ILE:HD13	4:D:71:PHE:CE1	2.19	0.78
4:D:183:LEU:HD21	4:D:195:ASN:HD21	1.49	0.78
4:D:63:TRP:N	4:D:63:TRP:CD1	2.53	0.77
4:D:25:ASN:O	4:D:29:ILE:HG13	1.85	0.77
4:D:221:SER:O	4:D:224:MET:HG2	1.84	0.77
1:A:108:PHE:CZ	1:A:146:ARG:HD2	2.19	0.76
4:D:156:ASP:OD2	4:D:160:ARG:NH2	2.19	0.76
2:B:59:GLU:OE2	2:B:59:GLU:HA	1.84	0.76
1:A:26:PHE:CE2	2:B:90:THR:HB	2.21	0.76
4:D:30:TYR:CE1	4:D:160:ARG:HD3	2.22	0.75
4:D:94:TYR:CD2	4:D:95:GLN:HG2	2.22	0.75
2:B:110:GLN:H	2:B:165:PRO:CG	1.94	0.74
2:B:21:THR:HG23	2:B:80:ARG:NE	2.03	0.74
4:D:162:TYR:HD2	4:D:163:LEU:HD23	1.53	0.73
1:A:69:ASN:ND2	3:C:316:LEU:HG	2.03	0.73
2:B:129:VAL:HG22	2:B:175:VAL:CG2	2.16	0.73
2:B:48:ARG:HG2	2:B:48:ARG:HH11	1.52	0.73
1:A:57:GLN:NE2	4:D:95:GLN:HB3	2.03	0.72
4:D:59:THR:O	4:D:60:ASP:HB2	1.89	0.72
4:D:221:SER:HA	4:D:224:MET:CE	2.20	0.72
2:B:98:LYS:HD2	2:B:98:LYS:C	2.15	0.72
2:B:110:GLN:HG3	2:B:165:PRO:O	1.91	0.70
2:B:85:VAL:HG13	3:C:306:PRO:HB2	1.72	0.70
4:D:94:TYR:O	4:D:95:GLN:HB2	1.91	0.70
4:D:52:ILE:HD11	4:D:77:VAL:CG1	2.19	0.70
2:B:141:GLY:O	2:B:161:LEU:HD12	1.92	0.70
1:A:160:VAL:HB	1:A:177:HIS:HE1	1.57	0.69
2:B:57:ASP:O	2:B:60:TYR:HB3	1.92	0.69
2:B:141:GLY:C	2:B:161:LEU:HD12	2.17	0.69
1:A:82:ILE:HD13	2:B:33:ASN:CB	2.18	0.69
2:B:71:ARG:HH11	2:B:71:ARG:HG2	1.58	0.69
2:B:177:HIS:CD2	2:B:178:PRO:HD2	2.28	0.69
2:B:116:VAL:HG21	2:B:160:MET:HE2	1.75	0.69
2:B:25:ARG:NH1	2:B:41:ASP:OD2	2.24	0.68
2:B:81:HIS:O	2:B:85:VAL:HG23	1.93	0.68
4:D:112:THR:HG23	4:D:151:THR:HG22	1.74	0.68
2:B:133:ARG:HG3	2:B:171:TYR:CE1	2.29	0.68
2:B:31:ILE:HG13	2:B:36:GLU:HA	1.76	0.68
4:D:31:TYR:HE1	4:D:165:GLU:HG2	1.59	0.68
1:A:45:LEU:HD12	1:A:48:PHE:CZ	2.29	0.68

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:36:MET:SD	1:A:63:ILE:HD12	2.34	0.67
1:A:97:VAL:HG11	1:A:178:TRP:CZ2	2.29	0.67
2:B:21:THR:HG23	2:B:80:ARG:CD	2.25	0.67
2:B:96:GLU:HA	2:B:179:SER:OG	1.93	0.67
2:B:150:ASN:HD21	2:B:156:GLN:HG2	1.60	0.67
1:A:128:VAL:HG12	1:A:130:THR:H	1.60	0.66
4:D:174:VAL:C	4:D:176:ASP:H	2.03	0.66
4:D:129:LEU:HD23	4:D:130:TRP:N	2.10	0.66
4:D:205:TYR:O	4:D:209:LEU:HG	1.96	0.66
4:D:4:SER:O	4:D:8:ASN:HB2	1.95	0.65
2:B:142:VAL:HG12	2:B:159:VAL:HG12	1.78	0.65
1:A:98:GLU:HB2	1:A:101:GLU:HB3	1.78	0.65
4:D:35:LYS:HD2	4:D:35:LYS:O	1.96	0.65
2:B:29:ARG:NH1	2:B:36:GLU:OE2	2.30	0.65
1:A:98:GLU:HB2	1:A:101:GLU:CB	2.28	0.64
1:A:87:PRO:CB	1:A:112:PHE:HB3	2.27	0.64
1:A:164:ARG:HH11	1:A:164:ARG:HG2	1.61	0.64
4:D:25:ASN:HA	4:D:28:GLN:HB2	1.78	0.64
4:D:150:VAL:HG11	4:D:155:LEU:HD11	1.80	0.64
1:A:97:VAL:HG11	1:A:178:TRP:HZ2	1.62	0.64
1:A:17:ASP:CG	2:B:6:ARG:NH1	2.56	0.64
1:A:87:PRO:HB3	1:A:112:PHE:CB	2.28	0.64
1:A:77:SER:O	1:A:80:THR:HG23	1.98	0.63
2:B:60:TYR:CE2	3:C:317:ALA:HA	2.34	0.63
4:D:62:SER:HG	4:D:63:TRP:CD1	2.17	0.63
1:A:82:ILE:HD12	1:A:83:THR:H	1.63	0.63
2:B:122:PHE:HB2	2:B:155:PHE:H	1.63	0.63
4:D:112:THR:HG23	4:D:151:THR:CG2	2.28	0.63
1:A:109:ILE:HD12	1:A:109:ILE:H	1.63	0.63
2:B:129:VAL:HA	2:B:174:GLN:O	1.99	0.63
4:D:52:ILE:CD1	4:D:77:VAL:HG13	2.24	0.63
4:D:119:LEU:HD21	4:D:148:LYS:HA	1.80	0.63
2:B:118:SER:HB2	2:B:158:LEU:HD11	1.80	0.63
2:B:29:ARG:HG2	2:B:29:ARG:HH11	1.64	0.62
2:B:127:ILE:HG13	2:B:128:GLU:N	2.14	0.62
4:D:30:TYR:CZ	4:D:160:ARG:HD3	2.35	0.62
2:B:142:VAL:HG12	2:B:159:VAL:CG1	2.30	0.62
2:B:185:THR:O	2:B:186:VAL:HG23	2.00	0.61
1:A:122:LEU:HD22	1:A:125:GLY:O	2.00	0.61
4:D:94:TYR:O	4:D:95:GLN:CB	2.49	0.61
4:D:63:TRP:N	4:D:63:TRP:HD1	1.98	0.61

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:109:LEU:HD11	2:B:190:ALA:OXT	2.00	0.61
2:B:152:ASP:O	2:B:153:TRP:HB2	2.01	0.61
4:D:41:LYS:O	4:D:85:VAL:HG22	2.01	0.60
4:D:37:LYS:O	4:D:37:LYS:CD	2.49	0.60
2:B:115:LEU:O	2:B:160:MET:HA	2.02	0.60
4:D:51:THR:O	4:D:52:ILE:HD12	2.01	0.60
1:A:105:LEU:HG	1:A:153:PHE:CE1	2.36	0.60
1:A:85:VAL:HB	1:A:113:THR:HG22	1.83	0.60
2:B:102:TYR:CD1	2:B:102:TYR:C	2.80	0.60
2:B:62:ASN:HA	2:B:68:LEU:HD23	1.83	0.59
2:B:142:VAL:HA	2:B:160:MET:O	2.01	0.59
4:D:6:GLU:HA	4:D:9:GLU:HG3	1.83	0.59
4:D:12:LEU:HA	4:D:197:ASP:OD2	2.02	0.59
1:A:45:LEU:HD13	2:B:93:ARG:HH12	1.67	0.59
1:A:103:ASN:O	1:A:104:VAL:HG23	2.01	0.59
2:B:114:LEU:HD22	2:B:162:GLU:HG2	1.84	0.59
2:B:117:CYS:HB2	2:B:131:TRP:CZ2	2.38	0.59
1:A:9:GLN:HB3	2:B:13:PHE:HB2	1.84	0.59
2:B:122:PHE:CB	2:B:155:PHE:H	2.16	0.59
1:A:26:PHE:HB2	1:A:31:ILE:HD11	1.85	0.59
4:D:135:GLN:C	4:D:136:ASN:HD22	2.10	0.58
2:B:116:VAL:CG2	2:B:160:MET:HE2	2.32	0.58
4:D:28:GLN:O	4:D:33:ASN:HB2	2.04	0.58
2:B:21:THR:HG23	2:B:80:ARG:HE	1.66	0.58
2:B:129:VAL:CG2	2:B:175:VAL:HG13	2.33	0.58
4:D:135:GLN:O	4:D:136:ASN:ND2	2.35	0.58
4:D:141:GLU:CD	4:D:141:GLU:H	2.12	0.58
1:A:87:PRO:CA	1:A:112:PHE:HB3	2.34	0.58
1:A:57:GLN:HE21	4:D:95:GLN:NE2	2.02	0.58
4:D:101:PRO:C	4:D:103:LYS:H	2.11	0.58
1:A:122:LEU:HA	1:A:126:LYS:O	2.03	0.57
4:D:90:ALA:O	4:D:210:LEU:HD13	2.04	0.57
1:A:118:ASN:HB2	1:A:166:GLU:HB2	1.86	0.57
2:B:28:GLU:HB3	2:B:40:PHE:HB3	1.86	0.57
4:D:42:GLU:O	4:D:43:SER:CB	2.53	0.57
4:D:182:GLY:O	4:D:198:LEU:HB2	2.04	0.57
4:D:59:THR:O	4:D:59:THR:HG23	2.04	0.57
4:D:18:LEU:HD12	4:D:23:LEU:HD13	1.85	0.57
4:D:117:ASN:HD21	4:D:147:LYS:HB3	1.69	0.57
2:B:142:VAL:HG22	2:B:161:LEU:HD13	1.86	0.57
4:D:131:LEU:HD22	4:D:167:TYR:HE2	1.69	0.57

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:153:GLN:NE2	4:D:213:TYR:HB3	2.19	0.57
4:D:72:ASP:HB2	4:D:76:ILE:HD12	1.85	0.57
2:B:47:TYR:CD2	2:B:61:TRP:CE3	2.93	0.57
1:A:39:LYS:HE3	1:A:60:LEU:HD11	1.87	0.56
2:B:122:PHE:CB	2:B:155:PHE:HB2	2.36	0.56
4:D:13:ARG:NH1	4:D:202:GLN:HE22	2.00	0.56
1:A:62:ASN:CG	3:C:313:THR:HG23	2.30	0.56
1:A:174:LEU:HD12	1:A:175:LEU:N	2.20	0.56
2:B:114:LEU:CD2	2:B:162:GLU:HG2	2.35	0.56
2:B:129:VAL:CG2	2:B:175:VAL:HG22	2.21	0.56
4:D:182:GLY:C	4:D:198:LEU:HD12	2.30	0.56
4:D:212:ILE:HD11	4:D:213:TYR:CD2	2.41	0.56
4:D:52:ILE:HD13	4:D:71:PHE:HE1	1.64	0.56
1:A:119:VAL:HG21	1:A:149:HIS:CE1	2.41	0.56
4:D:130:TRP:HE3	4:D:228:ILE:O	1.89	0.56
1:A:39:LYS:CG	1:A:60:LEU:HD11	2.31	0.56
1:A:164:ARG:HG2	1:A:164:ARG:NH1	2.19	0.56
2:B:122:PHE:O	2:B:123:TYR:HB2	2.06	0.55
1:A:97:VAL:HA	1:A:103:ASN:HD21	1.70	0.55
1:A:103:ASN:HB3	1:A:153:PHE:CE1	2.41	0.55
4:D:65:ASN:OD1	4:D:103:LYS:HB3	2.06	0.55
2:B:98:LYS:C	2:B:98:LYS:CD	2.80	0.55
2:B:111:HIS:O	2:B:112:HIS:O	2.24	0.55
4:D:13:ARG:HG3	4:D:13:ARG:HH11	1.70	0.55
2:B:3:THR:HG22	2:B:4:ARG:CD	2.36	0.55
2:B:122:PHE:HB2	2:B:155:PHE:N	2.21	0.55
4:D:39:GLU:HA	4:D:86:ASP:OD1	2.06	0.55
1:A:128:VAL:HG12	1:A:129:THR:N	2.22	0.55
2:B:18:PHE:HD1	2:B:23:ARG:HG2	1.72	0.55
1:A:65:VAL:O	1:A:68:ALA:HB3	2.07	0.55
4:D:161:ARG:O	4:D:165:GLU:HG3	2.07	0.54
2:B:29:ARG:HH11	2:B:29:ARG:CG	2.20	0.54
2:B:73:ALA:C	2:B:75:VAL:H	2.14	0.54
2:B:133:ARG:HG3	2:B:171:TYR:HE1	1.73	0.54
1:A:17:ASP:OD1	2:B:6:ARG:NH1	2.41	0.54
2:B:129:VAL:HG23	2:B:175:VAL:HG13	1.89	0.54
4:D:64:TYR:HA	4:D:103:LYS:O	2.08	0.54
4:D:127:ILE:HD11	4:D:143:VAL:HG21	1.88	0.54
2:B:96:GLU:HG3	2:B:180:VAL:CG1	2.37	0.54
4:D:127:ILE:HG12	4:D:226:ILE:HB	1.89	0.54
1:A:79:TYR:CD1	1:A:79:TYR:N	2.76	0.53

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:105:LEU:HG	1:A:153:PHE:CZ	2.43	0.53
2:B:21:THR:O	2:B:21:THR:CG2	2.56	0.53
2:B:111:HIS:O	2:B:112:HIS:C	2.51	0.53
2:B:122:PHE:CZ	2:B:127:ILE:HG12	2.37	0.53
4:D:182:GLY:HA3	4:D:198:LEU:HD12	1.90	0.53
2:B:27:LEU:HD12	2:B:40:PHE:O	2.08	0.53
4:D:169:LEU:HD23	4:D:170:TYR:CE2	2.44	0.53
2:B:88:SER:O	2:B:92:GLN:HB2	2.09	0.53
1:A:62:ASN:ND2	3:C:313:THR:HG23	2.24	0.53
4:D:46:GLN:NE2	4:D:77:VAL:HG11	2.24	0.53
4:D:42:GLU:HG2	4:D:43:SER:H	1.73	0.53
4:D:61:HIS:CE1	4:D:63:TRP:CD1	2.97	0.53
4:D:156:ASP:OD2	4:D:160:ARG:HD2	2.08	0.53
4:D:179:VAL:HG12	4:D:181:ARG:O	2.09	0.53
1:A:30:GLU:OE1	1:A:44:ARG:HD3	2.09	0.52
1:A:132:VAL:HG23	1:A:132:VAL:O	2.09	0.52
2:B:129:VAL:HG13	2:B:173:CYS:SG	2.49	0.52
2:B:73:ALA:O	2:B:75:VAL:N	2.43	0.52
2:B:93:ARG:HG2	2:B:123:TYR:CD1	2.45	0.52
2:B:145:THR:HG23	2:B:158:LEU:O	2.10	0.52
4:D:42:GLU:HG2	4:D:43:SER:N	2.25	0.52
4:D:52:ILE:HD13	4:D:71:PHE:CD1	2.44	0.52
1:A:34:VAL:HG12	1:A:36:MET:HE2	1.90	0.52
4:D:18:LEU:HD12	4:D:23:LEU:CD1	2.40	0.52
2:B:102:TYR:C	2:B:102:TYR:HD1	2.18	0.52
4:D:24:GLY:O	4:D:28:GLN:HB2	2.09	0.52
4:D:141:GLU:CD	4:D:141:GLU:N	2.68	0.52
4:D:131:LEU:O	4:D:132:ASP:HB2	2.10	0.52
4:D:131:LEU:HD22	4:D:167:TYR:CE2	2.45	0.52
2:B:18:PHE:CD1	2:B:23:ARG:NE	2.78	0.52
4:D:222:GLU:C	4:D:224:MET:H	2.17	0.52
2:B:70:GLN:O	2:B:72:ARG:N	2.42	0.52
4:D:108:TYR:OH	4:D:211:ARG:NH1	2.43	0.52
4:D:101:PRO:O	4:D:103:LYS:N	2.41	0.51
4:D:150:VAL:CG1	4:D:155:LEU:HD11	2.40	0.51
1:A:116:VAL:HG11	1:A:168:TRP:CH2	2.44	0.51
2:B:3:THR:HG22	2:B:4:ARG:H	1.76	0.51
4:D:174:VAL:C	4:D:176:ASP:N	2.68	0.51
2:B:4:ARG:HB3	2:B:5:PRO:CD	2.41	0.51
2:B:25:ARG:HH21	2:B:27:LEU:HD22	1.73	0.51
2:B:30:CYS:C	2:B:31:ILE:HD12	2.36	0.51

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:26:LEU:HD13	4:D:170:TYR:CE1	2.46	0.51
4:D:31:TYR:CE1	4:D:165:GLU:HG2	2.43	0.51
1:A:117:VAL:HG13	1:A:117:VAL:O	2.10	0.51
4:D:138:VAL:HG22	4:D:162:TYR:OH	2.11	0.51
2:B:71:ARG:HG2	2:B:71:ARG:NH1	2.24	0.51
1:A:108:PHE:CE1	1:A:146:ARG:HG3	2.46	0.51
1:A:154:LEU:HD12	1:A:155:PRO:HD2	1.91	0.51
4:D:9:GLU:O	4:D:10:LYS:C	2.52	0.51
2:B:122:PHE:CG	2:B:155:PHE:HB2	2.46	0.51
2:B:105:LYS:O	2:B:106:THR:C	2.54	0.51
1:A:107:CYS:HB2	1:A:121:TRP:CZ2	2.46	0.51
2:B:109:LEU:HD21	2:B:190:ALA:OXT	2.11	0.50
4:D:35:LYS:HB2	4:D:89:GLY:O	2.10	0.50
2:B:29:ARG:NH1	2:B:29:ARG:HG2	2.25	0.50
4:D:94:TYR:CE2	4:D:95:GLN:HG2	2.46	0.50
1:A:119:VAL:HA	1:A:164:ARG:O	2.12	0.50
2:B:23:ARG:NH1	5:B:191:HOH:O	2.44	0.50
2:B:73:ALA:C	2:B:75:VAL:N	2.70	0.50
4:D:1:SER:HA	4:D:187:HIS:HE1	1.75	0.50
4:D:208:THR:HA	4:D:211:ARG:HG3	1.92	0.50
1:A:129:THR:HG22	1:A:129:THR:O	2.12	0.50
2:B:110:GLN:CA	2:B:165:PRO:HG2	2.40	0.50
4:D:15:LYS:HG3	4:D:199:PHE:CE1	2.47	0.50
1:A:14:LEU:HD12	2:B:7:PHE:O	2.12	0.50
1:A:22:PHE:CD1	1:A:22:PHE:C	2.89	0.50
2:B:21:THR:HG23	2:B:80:ARG:HD3	1.93	0.50
2:B:94:ARG:HG3	2:B:94:ARG:HH11	1.77	0.50
2:B:48:ARG:HH11	2:B:48:ARG:CG	2.25	0.49
2:B:130:ARG:HD3	2:B:137:GLU:OE2	2.12	0.49
2:B:11:LEU:CD2	3:C:313:THR:HG22	2.43	0.49
1:A:7:ILE:O	1:A:8:ILE:HD13	2.11	0.49
4:D:40:ASN:CG	4:D:40:ASN:O	2.55	0.49
2:B:4:ARG:HD3	2:B:4:ARG:H	1.78	0.49
2:B:125:GLY:O	2:B:126:SER:C	2.53	0.49
4:D:59:THR:O	4:D:60:ASP:CB	2.58	0.49
2:B:116:VAL:HG11	2:B:158:LEU:HD23	1.93	0.49
2:B:155:PHE:N	2:B:155:PHE:CD1	2.80	0.49
4:D:166:LYS:HB3	4:D:167:TYR:CD1	2.48	0.49
4:D:174:VAL:O	4:D:176:ASP:N	2.37	0.49
1:A:168:TRP:CH2	2:B:6:ARG:CZ	2.96	0.49
2:B:60:TYR:O	2:B:63:SER:HB3	2.12	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:187:GLU:CD	2:B:187:GLU:N	2.55	0.49
2:B:30:CYS:HB2	2:B:38:VAL:HG12	1.95	0.49
1:A:66:ASP:O	1:A:67:LYS:C	2.53	0.49
2:B:105:LYS:O	2:B:107:GLN:HG3	2.13	0.49
3:C:312:ASN:ND2	3:C:312:ASN:H	2.11	0.49
2:B:108:PRO:O	2:B:109:LEU:O	2.31	0.48
4:D:20:GLY:O	4:D:22:ALA:N	2.46	0.48
2:B:17:PHE:CZ	2:B:83:TYR:HB2	2.48	0.48
2:B:166:ARG:HB2	2:B:169:GLU:CD	2.38	0.48
4:D:40:ASN:OD1	4:D:84:LYS:HD2	2.14	0.48
4:D:87:LEU:HB3	4:D:107:MET:HE1	1.95	0.48
2:B:80:ARG:HG3	2:B:80:ARG:HH11	1.79	0.48
2:B:89:PHE:CD1	2:B:90:THR:HG23	2.49	0.48
4:D:114:HIS:O	4:D:115:ASP:C	2.56	0.48
1:A:90:THR:O	1:A:90:THR:OG1	2.15	0.48
2:B:142:VAL:CG1	2:B:159:VAL:HG12	2.41	0.48
1:A:162:ASP:OD1	1:A:177:HIS:HA	2.14	0.48
1:A:162:ASP:CG	1:A:177:HIS:HD1	2.21	0.48
2:B:97:PRO:HD3	2:B:177:HIS:ND1	2.28	0.48
4:D:12:LEU:O	4:D:14:LYS:HE3	2.14	0.48
1:A:128:VAL:HG12	1:A:129:THR:H	1.79	0.48
4:D:212:ILE:HD11	4:D:213:TYR:CE2	2.49	0.48
1:A:168:TRP:CH2	2:B:6:ARG:NH2	2.82	0.48
1:A:180:PHE:HE1	1:A:182:ALA:HB2	1.78	0.48
4:D:187:HIS:HB3	4:D:225:HIS:CE1	2.49	0.48
2:B:122:PHE:HZ	2:B:127:ILE:CG1	2.23	0.47
4:D:67:LEU:HD12	4:D:105:ALA:O	2.14	0.47
4:D:179:VAL:CG1	4:D:181:ARG:O	2.62	0.47
2:B:122:PHE:HB2	2:B:155:PHE:CA	2.44	0.47
4:D:84:LYS:HB3	4:D:114:HIS:HB3	1.97	0.47
2:B:116:VAL:HG13	2:B:160:MET:CG	2.40	0.47
1:A:17:ASP:OD1	2:B:6:ARG:HD2	2.15	0.47
2:B:48:ARG:HG2	2:B:48:ARG:NH1	2.26	0.47
1:A:9:GLN:HB2	2:B:78:TYR:OH	2.15	0.47
2:B:13:PHE:CE2	2:B:28:GLU:HG3	2.49	0.47
1:A:9:GLN:OE1	3:C:311:GLN:N	2.46	0.47
4:D:53:LEU:HB2	4:D:68:LEU:CD1	2.44	0.47
4:D:69:VAL:HG11	4:D:71:PHE:CZ	2.49	0.47
1:A:76:ARG:NH1	2:B:53:LEU:O	2.47	0.47
2:B:31:ILE:HG13	2:B:36:GLU:CA	2.44	0.47
2:B:127:ILE:HD12	2:B:176:GLU:O	2.15	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:118:ARG:HG3	4:D:118:ARG:HH11	1.79	0.47
1:A:11:GLU:HG3	2:B:11:LEU:HB3	1.97	0.46
2:B:28:GLU:O	2:B:39:ARG:HB2	2.14	0.46
2:B:110:GLN:O	2:B:164:VAL:HG13	2.15	0.46
2:B:118:SER:O	2:B:118:SER:OG	2.31	0.46
4:D:20:GLY:C	4:D:22:ALA:H	2.24	0.46
1:A:48:PHE:CD1	1:A:48:PHE:N	2.82	0.46
4:D:167:TYR:O	4:D:168:ASN:C	2.58	0.46
2:B:4:ARG:CD	2:B:4:ARG:H	2.28	0.46
1:A:43:TRP:HZ2	1:A:53:SER:HA	1.80	0.46
1:A:109:ILE:HD12	1:A:109:ILE:N	2.29	0.46
4:D:44:HIS:HA	4:D:81:LYS:CG	2.34	0.46
1:A:162:ASP:HA	1:A:176:LYS:O	2.15	0.46
2:B:18:PHE:HD1	2:B:23:ARG:NE	2.14	0.46
4:D:135:GLN:C	4:D:136:ASN:ND2	2.73	0.46
4:D:222:GLU:O	4:D:223:ASN:HB2	2.15	0.46
1:A:57:GLN:HA	1:A:57:GLN:OE1	2.16	0.46
1:A:135:THR:OG1	1:A:148:PHE:HB2	2.15	0.46
2:B:55:ARG:O	2:B:56:PRO:C	2.55	0.46
4:D:117:ASN:ND2	4:D:147:LYS:HB3	2.30	0.46
4:D:134:LYS:O	4:D:136:ASN:ND2	2.48	0.46
1:A:15:ASN:ND2	1:A:70:LEU:HD23	2.31	0.46
1:A:147:LYS:HG2	1:A:148:PHE:N	2.31	0.46
4:D:153:GLN:HA	4:D:213:TYR:CE1	2.50	0.46
1:A:123:ARG:O	1:A:124:ASN:HB2	2.15	0.46
2:B:122:PHE:HB2	2:B:155:PHE:O	2.15	0.46
2:B:96:GLU:HG3	2:B:180:VAL:HG13	1.97	0.45
1:A:140:ARG:HG2	1:A:140:ARG:HH11	1.80	0.45
2:B:90:THR:O	2:B:93:ARG:HB3	2.16	0.45
4:D:53:LEU:HB2	4:D:68:LEU:HD12	1.96	0.45
2:B:59:GLU:OE2	2:B:59:GLU:CA	2.61	0.45
4:D:135:GLN:HE21	4:D:135:GLN:HB2	1.59	0.45
1:A:34:VAL:CG1	1:A:36:MET:HE2	2.46	0.45
4:D:27:LYS:HB2	4:D:170:TYR:HB2	1.97	0.45
1:A:172:GLU:O	1:A:173:PRO:O	2.34	0.45
2:B:70:GLN:C	2:B:72:ARG:H	2.25	0.45
4:D:25:ASN:HA	4:D:28:GLN:CB	2.44	0.45
4:D:52:ILE:HD11	4:D:77:VAL:HG22	1.99	0.45
4:D:101:PRO:C	4:D:103:LYS:N	2.75	0.45
4:D:210:LEU:HD23	4:D:210:LEU:HA	1.80	0.45
4:D:182:GLY:CA	4:D:198:LEU:HD12	2.47	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:37:ALA:HA	4:D:50:HIS:NE2	2.32	0.45
1:A:64:ALA:HA	4:D:47:PHE:CE1	2.52	0.45
1:A:117:VAL:O	1:A:117:VAL:CG1	2.65	0.45
1:A:120:THR:O	1:A:164:ARG:HB3	2.17	0.45
2:B:32:TYR:O	2:B:33:ASN:HB2	2.17	0.45
2:B:112:HIS:O	2:B:112:HIS:ND1	2.50	0.45
4:D:36:ALA:O	4:D:88:TYR:HA	2.17	0.45
4:D:127:ILE:CD1	4:D:143:VAL:HG23	2.37	0.45
4:D:226:ILE:HG22	4:D:227:ALA:N	2.32	0.45
2:B:104:SER:O	2:B:105:LYS:CB	2.65	0.45
2:B:104:SER:O	2:B:105:LYS:HB2	2.17	0.45
4:D:30:TYR:CD1	4:D:160:ARG:HB3	2.52	0.44
2:B:74:ALA:O	2:B:78:TYR:HB3	2.17	0.44
4:D:1:SER:HA	4:D:187:HIS:CE1	2.51	0.44
4:D:19:GLN:C	4:D:204:GLN:HE21	2.25	0.44
2:B:21:THR:O	2:B:21:THR:HG22	2.18	0.44
2:B:122:PHE:HB2	2:B:155:PHE:C	2.42	0.44
4:D:31:TYR:CD1	4:D:164:GLN:HB3	2.52	0.44
4:D:132:ASP:C	4:D:134:LYS:H	2.25	0.44
1:A:89:VAL:HG23	1:A:174:LEU:HD23	1.99	0.44
2:B:19:ASN:ND2	2:B:22:GLU:CD	2.75	0.44
2:B:67:LEU:C	2:B:67:LEU:HD23	2.42	0.44
2:B:130:ARG:HG3	2:B:130:ARG:HH11	1.82	0.44
2:B:156:GLN:HE21	2:B:156:GLN:HB2	1.59	0.44
1:A:82:ILE:HD12	1:A:83:THR:N	2.29	0.44
4:D:149:ASN:HB3	4:D:218:THR:HG22	2.00	0.44
1:A:87:PRO:HB3	1:A:112:PHE:CG	2.53	0.44
1:A:100:ARG:HB2	1:A:100:ARG:NH1	2.33	0.44
2:B:13:PHE:CE2	3:C:311:GLN:HG2	2.53	0.44
1:A:146:ARG:HH11	1:A:146:ARG:HB3	1.83	0.44
1:A:42:VAL:O	1:A:42:VAL:HG12	2.18	0.43
4:D:27:LYS:O	4:D:27:LYS:HG2	2.17	0.43
4:D:99:GLY:HA3	4:D:104:THR:OG1	2.18	0.43
2:B:57:ASP:O	2:B:60:TYR:CB	2.64	0.43
2:B:127:ILE:HD11	2:B:175:VAL:HG12	2.00	0.43
2:B:98:LYS:HE3	2:B:120:SER:OG	2.18	0.43
4:D:156:ASP:CG	4:D:160:ARG:HD2	2.43	0.43
1:A:104:VAL:HG22	1:A:152:PRO:HA	1.99	0.43
1:A:172:GLU:O	1:A:173:PRO:C	2.59	0.43
2:B:9:TRP:CH2	3:C:316:LEU:HD11	2.54	0.43
2:B:70:GLN:C	2:B:72:ARG:N	2.75	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:133:ARG:O	2:B:134:ASN:HB2	2.18	0.43
1:A:89:VAL:C	1:A:90:THR:CG2	2.91	0.43
1:A:160:VAL:HG23	1:A:160:VAL:O	2.18	0.43
4:D:156:ASP:O	4:D:157:LEU:C	2.62	0.43
1:A:69:ASN:CG	3:C:316:LEU:HG	2.44	0.43
1:A:144:LEU:HG	2:B:34:GLN:NE2	2.33	0.43
2:B:64:GLN:C	2:B:66:ASP:N	2.74	0.43
4:D:13:ARG:NH1	4:D:13:ARG:HG3	2.32	0.43
4:D:30:TYR:CE1	4:D:157:LEU:HD23	2.54	0.43
4:D:95:GLN:HA	4:D:95:GLN:NE2	2.22	0.43
1:A:67:LYS:HE2	4:D:46:GLN:O	2.19	0.43
1:A:113:THR:OG1	1:A:114:PRO:HA	2.19	0.43
1:A:119:VAL:HB	1:A:149:HIS:NE2	2.33	0.43
1:A:93:THR:HG23	1:A:95:SER:N	2.33	0.43
2:B:31:ILE:HD12	2:B:31:ILE:N	2.33	0.43
1:A:99:LEU:HD21	1:A:180:PHE:CZ	2.54	0.42
2:B:61:TRP:C	2:B:63:SER:N	2.77	0.42
4:D:229:TYR:O	4:D:230:LEU:HD23	2.19	0.42
1:A:57:GLN:HE22	4:D:95:GLN:HB3	1.81	0.42
1:A:67:LYS:O	1:A:71:GLU:HG2	2.19	0.42
4:D:58:PHE:HB2	4:D:65:ASN:HA	2.00	0.42
1:A:91:VAL:HG23	1:A:176:LYS:HB3	2.01	0.42
2:B:19:ASN:HD22	2:B:22:GLU:HB2	1.83	0.42
1:A:39:LYS:HG2	1:A:60:LEU:CD1	2.35	0.42
1:A:62:ASN:OD1	3:C:310:LYS:NZ	2.53	0.42
2:B:13:PHE:CD2	3:C:311:GLN:HG2	2.53	0.42
4:D:125:VAL:HG12	4:D:226:ILE:HG12	2.00	0.42
2:B:122:PHE:CZ	2:B:127:ILE:CG1	3.01	0.42
4:D:125:VAL:HB	4:D:143:VAL:HB	2.01	0.42
1:A:132:VAL:O	1:A:132:VAL:CG2	2.67	0.42
2:B:36:GLU:O	2:B:50:VAL:CG2	2.68	0.42
2:B:129:VAL:O	2:B:129:VAL:HG12	2.19	0.42
4:D:149:ASN:HB3	4:D:218:THR:CG2	2.50	0.42
1:A:107:CYS:HB3	1:A:149:HIS:HB2	2.02	0.42
4:D:9:GLU:O	4:D:10:LYS:O	2.38	0.42
4:D:43:SER:OG	4:D:45:ASP:HB2	2.19	0.42
1:A:28:GLY:O	1:A:146:ARG:NH2	2.52	0.42
1:A:128:VAL:CG1	1:A:130:THR:HG23	2.49	0.42
4:D:84:LYS:HD3	4:D:84:LYS:HA	1.87	0.42
1:A:100:ARG:C	1:A:154:LEU:HD11	2.45	0.42
4:D:23:LEU:HG	4:D:170:TYR:O	2.20	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:72:ASP:HB2	4:D:76:ILE:CD1	2.49	0.42
4:D:229:TYR:C	4:D:230:LEU:HD23	2.44	0.42
2:B:122:PHE:O	2:B:123:TYR:CB	2.68	0.41
4:D:7:ILE:HG21	4:D:229:TYR:CD2	2.55	0.41
1:A:108:PHE:CE2	1:A:110:ASP:HB2	2.56	0.41
4:D:37:LYS:O	4:D:37:LYS:CG	2.68	0.41
2:B:11:LEU:CD2	3:C:313:THR:CG2	2.99	0.41
1:A:7:ILE:HA	1:A:25:ASP:O	2.21	0.41
2:B:77:THR:O	2:B:81:HIS:HB3	2.21	0.41
2:B:157:THR:O	2:B:158:LEU:HD12	2.20	0.41
4:D:46:GLN:HE21	4:D:77:VAL:HG11	1.84	0.41
1:A:32:PHE:CD1	1:A:32:PHE:C	2.98	0.41
1:A:73:MET:HA	1:A:76:ARG:HG2	2.02	0.41
1:A:92:LEU:HD12	1:A:92:LEU:HA	1.92	0.41
2:B:60:TYR:HB3	2:B:61:TRP:H	1.66	0.41
2:B:122:PHE:CE2	2:B:127:ILE:HB	2.56	0.41
4:D:92:TYR:CE2	4:D:106:CYS:HB2	2.56	0.41
4:D:206:SER:HA	4:D:209:LEU:HB2	2.02	0.41
1:A:91:VAL:HG11	1:A:178:TRP:HB2	2.02	0.41
1:A:180:PHE:CD1	1:A:180:PHE:C	2.99	0.41
2:B:24:VAL:HG23	2:B:80:ARG:NH2	2.36	0.41
4:D:219:ILE:HD13	4:D:219:ILE:H	1.86	0.41
1:A:89:VAL:C	1:A:90:THR:HG22	2.45	0.41
2:B:147:LEU:HD12	2:B:147:LEU:HA	1.79	0.41
1:A:12:PHE:CD1	1:A:12:PHE:C	2.99	0.41
1:A:23:MET:C	1:A:23:MET:SD	3.04	0.41
1:A:126:LYS:HA	1:A:127:PRO:HD3	1.80	0.41
2:B:107:GLN:HB3	2:B:108:PRO:CD	2.51	0.41
4:D:52:ILE:HD12	4:D:52:ILE:N	2.33	0.41
1:A:76:ARG:C	1:A:78:ASN:H	2.28	0.40
4:D:88:TYR:HD1	4:D:88:TYR:O	2.03	0.40
4:D:152:VAL:HG11	4:D:196:TYR:CE2	2.56	0.40
1:A:121:TRP:O	1:A:128:VAL:HG23	2.22	0.40
1:A:143:HIS:HD2	2:B:12:LYS:NZ	2.18	0.40
2:B:71:ARG:NH1	2:B:71:ARG:CG	2.84	0.40
2:B:116:VAL:CG1	2:B:160:MET:HG2	2.42	0.40
1:A:75:LYS:O	1:A:78:ASN:N	2.47	0.40
2:B:163:THR:HB	2:B:164:VAL:H	1.64	0.40
4:D:92:TYR:O	4:D:105:ALA:HA	2.22	0.40
1:A:95:SER:O	1:A:96:PRO:C	2.63	0.40
2:B:45:GLY:O	2:B:68:LEU:HD11	2.22	0.40

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:183:LEU:HG	4:D:184:ILE:N	2.36	0.40
1:A:87:PRO:HA	1:A:112:PHE:HB3	2.01	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	177/182 (97%)	151 (85%)	19 (11%)	7 (4%)	2	17
2	B	186/190 (98%)	147 (79%)	27 (14%)	12 (6%)	1	8
3	C	11/13 (85%)	11 (100%)	0	0	100	100
4	D	231/233 (99%)	189 (82%)	29 (13%)	13 (6%)	1	11
All	All	605/618 (98%)	498 (82%)	75 (12%)	32 (5%)	1	12

All (32) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	60	TYR
2	B	109	LEU
2	B	112	HIS
2	B	122	PHE
2	B	123	TYR
4	D	21	THR
4	D	43	SER
4	D	60	ASP
4	D	176	ASP
1	A	55	GLU
1	A	155	PRO
2	B	21	THR
2	B	71	ARG

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
2	B	74	ALA
2	B	105	LYS
4	D	10	LYS
4	D	95	GLN
4	D	102	ASN
2	B	126	SER
2	B	165	PRO
4	D	65	ASN
4	D	132	ASP
4	D	175	PHE
1	A	102	PRO
1	A	173	PRO
2	B	33	ASN
4	D	37	LYS
4	D	88	TYR
4	D	178	LYS
1	A	96	PRO
1	A	72	ILE
1	A	116	VAL

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	163/166 (98%)	146 (90%)	17 (10%)	7	28
2	B	170/171 (99%)	151 (89%)	19 (11%)	6	25
3	C	12/12 (100%)	11 (92%)	1 (8%)	10	38
4	D	210/210 (100%)	190 (90%)	20 (10%)	8	32
All	All	555/559 (99%)	498 (90%)	57 (10%)	7	28

All (57) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	23	MET
1	A	25	ASP

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	41	THR
1	A	55	GLU
1	A	60	LEU
1	A	82	ILE
1	A	90	THR
1	A	92	LEU
1	A	93	THR
1	A	100	ARG
1	A	120	THR
1	A	146	ARG
1	A	156	SER
1	A	157	THR
1	A	170	LEU
1	A	171	ASP
1	A	172	GLU
2	B	4	ARG
2	B	21	THR
2	B	29	ARG
2	B	35	GLU
2	B	59	GLU
2	B	64	GLN
2	B	65	LYS
2	B	67	LEU
2	B	68	LEU
2	B	70	GLN
2	B	113	ASN
2	B	128	GLU
2	B	142	VAL
2	B	147	LEU
2	B	155	PHE
2	B	156	GLN
2	B	167	SER
2	B	176	GLU
2	B	187	GLU
3	C	307	LYS
4	D	5	GLU
4	D	14	LYS
4	D	19	GLN
4	D	33	ASN
4	D	39	GLU
4	D	60	ASP
4	D	63	TRP

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
4	D	65	ASN
4	D	70	ASP
4	D	95	GLN
4	D	112	THR
4	D	135	GLN
4	D	145	THR
4	D	164	GLN
4	D	185	VAL
4	D	195	ASN
4	D	208	THR
4	D	212	ILE
4	D	219	ILE
4	D	224	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	15	ASN
1	A	18	GLN
1	A	84	ASN
1	A	143	HIS
2	B	10	GLN
2	B	19	ASN
2	B	34	GLN
2	B	64	GLN
2	B	111	HIS
2	B	113	ASN
2	B	150	ASN
2	B	156	GLN
3	C	312	ASN
4	D	46	GLN
4	D	49	GLN
4	D	95	GLN
4	D	135	GLN
4	D	136	ASN
4	D	158	GLN
4	D	187	HIS
4	D	195	ASN
4	D	202	GLN
4	D	204	GLN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

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

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	179/182 (98%)	-0.35	0 100 100	2, 26, 95, 145	0
2	B	188/190 (98%)	-0.16	5 (2%) 56 36	2, 38, 126, 157	0
3	C	13/13 (100%)	-0.72	0 100 100	3, 9, 24, 43	0
4	D	233/233 (100%)	-0.11	2 (0%) 81 64	2, 39, 100, 126	0
All	All	613/618 (99%)	-0.21	7 (1%) 78 61	2, 35, 108, 157	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	164	VAL	3.5
4	D	176	ASP	2.6
4	D	187	HIS	2.5
2	B	114	LEU	2.3
2	B	144	SER	2.3
2	B	190	ALA	2.2
2	B	104	SER	2.1

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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.