Full wwPDB X-ray Structure Validation Report

Jan 31, 2016 – 10:12 PM GMT

PDB ID	:	1SME
Title	:	PLASMEPSIN II, A HEMOGLOBIN-DEGRADING ENZYME FROM PLASMODIUM FALCIPARUM, IN COMPLEX WITH PEPSTATIN A
Authors	:	Silva, A.M.; Lee, A.Y.; Gulnik, S.V.; Goldberg, D.E.; Erickson, J.W.
Deposited on	:	1996-06-11
Resolution	:	2.70 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
http://wwpdb.org/validation/2016/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

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

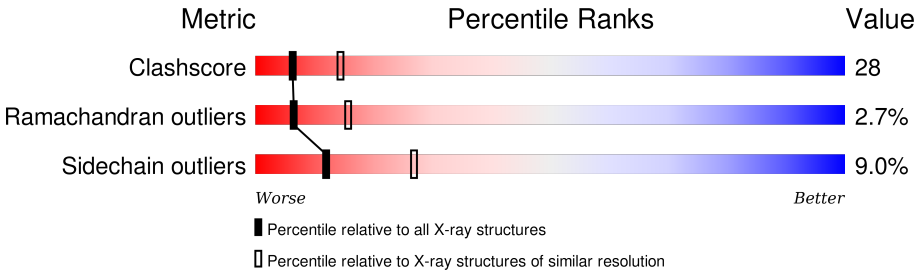
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	102246	2422 (2.70-2.70)
Ramachandran outliers	100387	2382 (2.70-2.70)
Sidechain outliers	100360	2382 (2.70-2.70)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	329	55% 39% 5%
1	B	329	43% 48% 8%
2	C	6	33% 33% 33%
2	D	6	17% 50% 33%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6723 atoms, of which 1298 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 PLASMEPSIN II.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	329	Total	C	H	N	O	S	0	0	0
			3134	1689	527	404	503	11			
1	B	329	Total	C	H	N	O	S	0	0	0
			3134	1689	527	404	503	11			

- Molecule 2 is a protein called Pepstatin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	6	Total	C	H	N	O	0	0	0
			55	34	7	5	9			
2	D	6	Total	C	H	N	O	0	0	0
			55	34	7	5	9			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	77	Total	H	O	0	0
			231	154	77		
3	B	34	Total	H	O	0	0
			102	68	34		
3	C	1	Total	H	O	0	0
			3	2	1		
3	D	3	Total	H	O	0	0
			9	6	3		

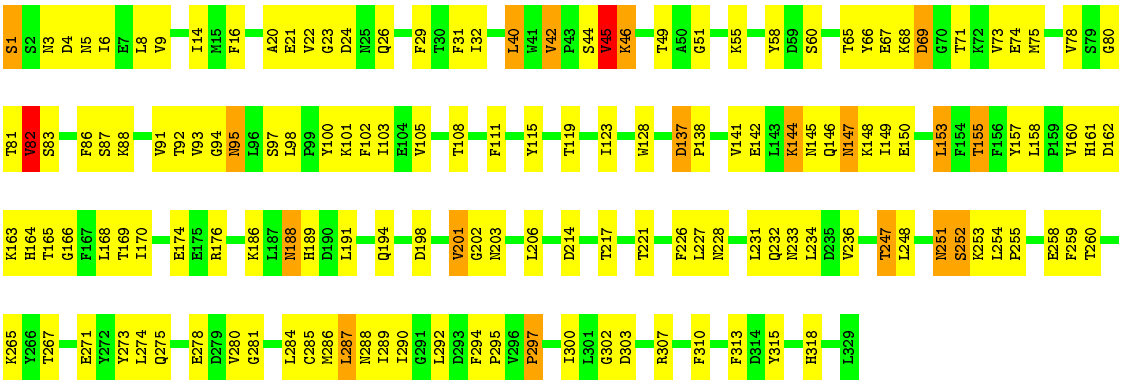
3 Residue-property plots (i)

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.

Note EDS was not executed.

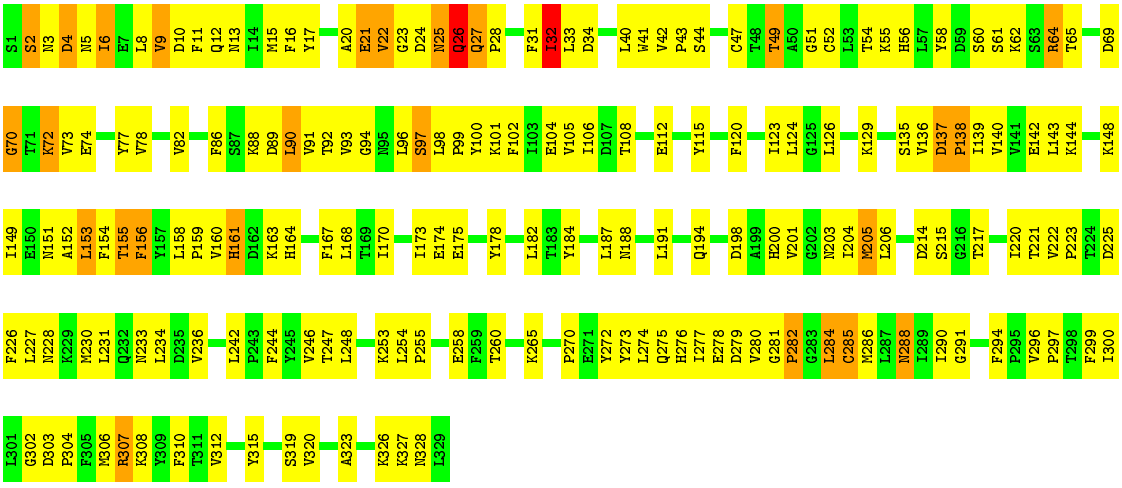
- Molecule 1: PLASMEPSIN II

Chain A:



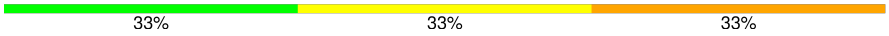
- Molecule 1: PLASMEPSIN II

Chain B:



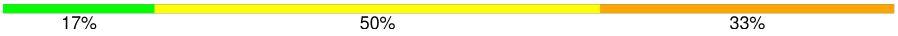
- Molecule 2: Pepstatin

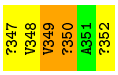
Chain C:



- Molecule 2: Pepstatin

Chain D:





4 Data and refinement statistics (i)

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α, β, γ	142.10Å 142.10Å 97.60Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	8.00 – 2.70	Depositor
% Data completeness (in resolution range)	74.3 (8.00-2.70)	Depositor
R_{merge}	0.94	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.195 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	6723	wwPDB-VP
Average B, all atoms (Å^2)	18.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

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

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	0.56	0/2674	0.96	11/3638 (0.3%)
1	B	0.56	1/2674 (0.0%)	0.89	9/3638 (0.2%)
2	C	0.60	0/17	2.90	1/21 (4.8%)
2	D	0.85	0/17	1.62	0/21
All	All	0.56	1/5382 (0.0%)	0.94	21/7318 (0.3%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
2	C	0	2
2	D	0	3
All	All	0	5

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	307	ARG	CZ-NH2	6.51	1.41	1.33

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	C	348	VAL	CG1-CB-CG2	-10.12	94.71	110.90
1	B	82	VAL	CG1-CB-CG2	-9.82	95.19	110.90
1	A	9	VAL	CG1-CB-CG2	-9.80	95.21	110.90
1	A	42	VAL	CG1-CB-CG2	8.27	124.13	110.90
1	B	9	VAL	CG1-CB-CG2	7.76	123.32	110.90
1	A	82	VAL	CG1-CB-CG2	7.47	122.85	110.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	320	VAL	CG1-CB-CG2	-7.39	99.07	110.90
1	A	45	VAL	CG1-CB-CG2	-7.35	99.14	110.90
1	B	22	VAL	CG1-CB-CG2	-7.03	99.66	110.90
1	A	201	VAL	CG1-CB-CG2	-6.91	99.85	110.90
1	A	280	VAL	CG1-CB-CG2	-6.83	99.97	110.90
1	A	32	ILE	N-CA-C	-6.64	93.07	111.00
1	A	40	LEU	CD1-CG-CD2	-6.62	90.63	110.50
1	A	202	GLY	N-CA-C	-6.53	96.79	113.10
1	A	248	LEU	N-CA-C	-6.36	93.84	111.00
1	B	8	LEU	CD1-CG-CD2	-6.10	92.20	110.50
1	A	153	LEU	CD1-CG-CD2	-5.82	93.04	110.50
1	B	32	ILE	N-CA-C	-5.37	96.49	111.00
1	B	42	VAL	CG1-CB-CG2	5.18	119.18	110.90
1	B	285	CYS	CA-CB-SG	5.06	123.11	114.00
1	B	153	LEU	CD1-CG-CD2	-5.03	95.41	110.50

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	C	350	STA	Mainchain,Peptide
2	D	349	VAL	Peptide
2	D	350	STA	Mainchain,Peptide

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	2607	527	2536	117	0
1	B	2607	527	2536	173	0
2	C	48	7	61	4	0
2	D	48	7	60	10	0
3	A	77	154	0	0	0
3	B	34	68	0	0	0
3	C	1	2	0	0	0
3	D	3	6	0	0	0
All	All	5425	1298	5193	292	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 28.

All (292) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:276:HIS:HA	1:B:285:CYS:HB3	1.34	1.05
1:B:221:THR:HB	1:B:300:ILE:HB	1.42	0.97
1:B:44:SER:HB2	1:B:104:GLU:HG2	1.53	0.90
1:B:234:LEU:HD11	1:B:254:LEU:HD23	1.51	0.90
1:A:6:ILE:HD11	1:A:93:VAL:HG12	1.55	0.88
1:B:20:ALA:HB3	1:B:31:PHE:CE2	2.09	0.85
1:B:49:THR:HG22	1:B:52:CYS:H	1.41	0.84
1:B:276:HIS:HA	1:B:285:CYS:CB	2.10	0.82
1:A:221:THR:HB	1:A:300:ILE:HB	1.61	0.80
1:B:98:LEU:HD22	1:B:143:LEU:HD21	1.63	0.80
1:B:160:VAL:HG11	1:B:163:LYS:HD3	1.63	0.80
1:B:41:TRP:CE3	1:B:105:VAL:HG21	2.18	0.79
1:A:281:GLY:HA3	1:A:284:LEU:HD22	1.65	0.78
1:B:206:LEU:HD23	1:B:299:PHE:HE1	1.47	0.78
1:A:20:ALA:HB3	1:A:31:PHE:CE1	2.18	0.77
1:B:47:CYS:HB2	1:B:106:ILE:HA	1.67	0.76
1:B:247:THR:HG21	1:B:254:LEU:HD21	1.67	0.76
1:A:221:THR:HG23	1:A:292:LEU:HB3	1.69	0.74
1:B:222:VAL:HG23	1:B:299:PHE:CE2	2.21	0.74
1:A:254:LEU:HD13	1:A:274:LEU:HD11	1.66	0.74
1:A:6:ILE:HD11	1:A:93:VAL:CG1	2.18	0.73
1:B:20:ALA:HB3	1:B:31:PHE:HE2	1.51	0.73
1:B:49:THR:HG21	1:B:108:THR:OG1	1.89	0.73
1:A:236:VAL:HG22	1:A:247:THR:HG23	1.73	0.70
1:B:78:VAL:HG22	2:D:352:STA:HD21	1.73	0.70
1:B:164:HIS:HD2	1:B:327:LYS:HG3	1.56	0.70
1:B:223:PRO:HG2	1:B:226:PHE:HB2	1.73	0.70
1:B:2:SER:HB3	1:B:148:LYS:O	1.92	0.69
1:B:6:ILE:HD11	1:B:93:VAL:HG12	1.74	0.69
1:B:44:SER:HB3	1:B:106:ILE:HG22	1.75	0.68
1:B:26:GLN:OE1	1:B:28:PRO:HD3	1.93	0.68
1:A:6:ILE:HD12	1:A:168:LEU:HD23	1.77	0.67
1:B:144:LYS:HA	1:B:149:ILE:HG22	1.76	0.67
1:A:155:THR:HG23	1:A:310:PHE:CE1	2.30	0.67
1:A:75:MET:HB3	1:A:82:VAL:HG12	1.78	0.66
1:B:206:LEU:HD23	1:B:299:PHE:CE1	2.31	0.65
1:B:188:ASN:HB2	1:B:194:GLN:HG3	1.77	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:294:PHE:CZ	1:A:300:ILE:HD11	2.31	0.65
1:A:155:THR:HB	1:A:169:THR:HB	1.79	0.65
1:A:290:ILE:HG13	2:C:347:IVA:HG11	1.79	0.65
1:B:275:GLN:O	1:B:285:CYS:HB2	1.97	0.65
1:A:137:ASP:HB3	1:A:142:GLU:HG2	1.79	0.64
1:B:23:GLY:HA2	1:B:89:ASP:OD1	1.96	0.64
1:B:242:LEU:HB3	1:B:244:PHE:CD2	2.33	0.64
1:A:86:PHE:CD1	1:A:101:LYS:HE2	2.33	0.63
1:A:60:SER:HB2	1:A:66:TYR:CD2	2.33	0.63
1:A:21:GLU:HB2	1:A:26:GLN:O	1.98	0.63
1:A:251:ASN:ND2	1:A:253:LYS:H	1.96	0.63
1:A:82:VAL:HG22	1:A:105:VAL:HG11	1.80	0.63
1:B:88:LYS:HE2	1:B:99:PRO:HB3	1.81	0.63
1:A:214:ASP:O	1:A:302:GLY:HA2	1.97	0.63
1:B:24:ASP:OD2	1:B:64:ARG:HG2	2.00	0.62
1:A:123:ILE:HD13	2:C:350:STA:HD11	1.81	0.61
1:B:25:ASN:HB2	1:B:62:LYS:O	1.99	0.61
1:A:128:TRP:HB2	1:A:191:LEU:O	2.00	0.61
1:A:60:SER:HB2	1:A:66:TYR:CG	2.34	0.61
1:B:242:LEU:HD13	1:B:244:PHE:CE2	2.35	0.61
1:A:174:GLU:HB3	1:A:176:ARG:HG2	1.83	0.60
1:B:149:ILE:HD11	1:B:170:ILE:O	2.00	0.60
1:B:294:PHE:HZ	1:B:300:ILE:HD11	1.67	0.60
1:A:29:PHE:HZ	1:A:58:TYR:HB2	1.67	0.60
1:A:66:TYR:HE1	1:A:87:SER:HB3	1.67	0.60
1:A:146:GLN:O	1:A:148:LYS:HG3	2.01	0.60
1:A:155:THR:HG23	1:A:310:PHE:HE1	1.67	0.59
1:B:160:VAL:HG22	1:B:307:ARG:HG2	1.84	0.59
1:B:6:ILE:O	1:B:167:PHE:HA	2.03	0.59
1:B:191:LEU:HB3	1:B:194:GLN:OE1	2.02	0.59
1:A:217:THR:HA	2:C:348:VAL:O	2.02	0.59
1:B:277:ILE:HG22	1:B:280:VAL:HB	1.84	0.58
1:B:234:LEU:HG	1:B:236:VAL:HG13	1.84	0.58
1:A:80:GLY:HA3	1:A:111:PHE:HD1	1.68	0.58
1:B:90:LEU:HA	1:B:99:PRO:HA	1.84	0.58
1:B:49:THR:HG22	1:B:52:CYS:N	2.17	0.58
1:A:138:PRO:HB2	1:A:141:VAL:HB	1.86	0.58
1:B:277:ILE:O	1:B:277:ILE:HG22	2.03	0.57
1:B:64:ARG:H	1:B:64:ARG:HD3	1.69	0.57
1:B:144:LYS:HD3	1:B:152:ALA:N	2.20	0.57
1:B:155:THR:HG23	1:B:310:PHE:CE1	2.38	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:129:LYS:HE3	1:B:135:SER:HB2	1.86	0.57
1:A:157:TYR:O	1:A:166:GLY:HA3	2.05	0.57
1:B:6:ILE:HB	1:B:168:LEU:HB3	1.87	0.57
1:A:247:THR:HG21	1:A:254:LEU:HD21	1.87	0.56
1:B:281:GLY:HA3	1:B:284:LEU:HD22	1.87	0.56
1:B:260:THR:HG22	1:B:265:LYS:HG3	1.87	0.56
1:B:9:VAL:O	1:B:16:PHE:HA	2.06	0.56
1:B:155:THR:O	1:B:168:LEU:HD12	2.06	0.56
1:B:69:ASP:HB3	1:B:86:PHE:O	2.06	0.56
1:B:270:PRO:HA	1:B:273:TYR:CZ	2.40	0.56
1:B:140:VAL:HB	1:B:315:TYR:HE2	1.71	0.56
1:B:25:ASN:O	1:B:26:GLN:HB3	2.06	0.56
1:A:231:LEU:HD23	1:A:234:LEU:HD12	1.88	0.56
1:B:22:VAL:O	1:B:27:GLN:HB2	2.06	0.56
1:A:188:ASN:ND2	1:A:194:GLN:HE21	2.02	0.56
1:B:204:ILE:CD1	1:B:230:MET:HB3	2.36	0.55
1:A:198:ASP:O	1:A:259:PHE:HA	2.06	0.55
1:B:204:ILE:HD11	1:B:230:MET:HA	1.88	0.55
1:B:43:PRO:HG2	1:B:56:HIS:O	2.07	0.55
1:B:273:TYR:HA	1:B:304:PRO:HB3	1.88	0.55
1:B:140:VAL:HB	1:B:315:TYR:CE2	2.42	0.55
1:B:277:ILE:HG13	1:B:286:MET:HE3	1.89	0.55
1:B:227:LEU:O	1:B:231:LEU:HG	2.06	0.55
1:A:258:GLU:HG3	1:A:267:THR:HG22	1.89	0.55
1:B:242:LEU:HD13	1:B:244:PHE:HE2	1.72	0.54
1:A:160:VAL:HG11	1:A:163:LYS:HE3	1.88	0.54
1:B:101:LYS:HD3	1:B:136:VAL:HG22	1.88	0.54
1:A:3:ASN:HA	1:A:170:ILE:O	2.08	0.54
1:A:260:THR:HG22	1:A:265:LYS:HA	1.90	0.54
1:B:182:LEU:HD12	1:B:323:ALA:HB2	1.88	0.54
1:B:47:CYS:CB	1:B:106:ILE:HA	2.37	0.53
1:B:155:THR:CG2	1:B:156:PHE:N	2.72	0.53
1:A:186:LYS:HD3	1:A:318:HIS:O	2.09	0.53
1:A:78:VAL:HG22	2:C:352:STA:HD21	1.89	0.53
1:B:214:ASP:O	1:B:302:GLY:HA2	2.08	0.53
1:A:273:TYR:O	1:A:288:ASN:HB3	2.09	0.53
1:B:40:LEU:HD13	1:B:124:LEU:HG	1.91	0.53
1:A:21:GLU:HG2	1:A:92:THR:HB	1.91	0.52
1:B:222:VAL:CG1	1:B:227:LEU:HB2	2.39	0.52
1:B:2:SER:HB2	1:B:149:ILE:HA	1.92	0.52
1:B:277:ILE:CG2	1:B:280:VAL:HB	2.40	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:98:LEU:HD13	1:B:143:LEU:HD23	1.92	0.51
1:B:222:VAL:HG22	1:B:226:PHE:HB3	1.91	0.51
1:B:222:VAL:O	1:B:291:GLY:HA2	2.10	0.51
1:A:188:ASN:HD21	1:A:194:GLN:HE21	1.58	0.51
1:B:204:ILE:HD13	1:B:230:MET:HB3	1.91	0.51
1:A:45:VAL:HG23	1:A:58:TYR:O	2.10	0.51
1:A:228:ASN:O	1:A:232:GLN:HG2	2.10	0.51
1:B:223:PRO:HG2	1:B:226:PHE:CB	2.39	0.51
2:D:348:VAL:HG11	2:D:350:STA:HD23	1.92	0.50
1:A:163:LYS:HD2	1:A:164:HIS:CD2	2.46	0.50
1:B:10:ASP:HB2	1:B:159:PRO:HG2	1.92	0.50
1:A:273:TYR:HE1	1:A:287:LEU:HD13	1.77	0.50
1:B:303:ASP:HB3	1:B:307:ARG:NH1	2.27	0.50
1:B:16:PHE:CD2	1:B:158:LEU:HD22	2.46	0.50
1:A:141:VAL:HG23	1:A:315:TYR:CD2	2.47	0.50
1:B:33:LEU:HD23	1:B:124:LEU:HB3	1.94	0.50
1:B:144:LYS:CA	1:B:149:ILE:HG22	2.40	0.49
1:B:86:PHE:HD1	1:B:102:PHE:O	1.95	0.49
1:A:201:VAL:HG23	1:A:226:PHE:HZ	1.77	0.49
1:A:163:LYS:HD2	1:A:164:HIS:HD2	1.78	0.49
1:A:294:PHE:HZ	1:A:300:ILE:HD11	1.78	0.49
1:B:173:ILE:HD11	1:B:312:VAL:HG11	1.95	0.49
1:A:29:PHE:CZ	1:A:58:TYR:HB2	2.48	0.49
1:B:163:LYS:HB3	1:B:327:LYS:HG3	1.94	0.49
1:A:67:GLU:HB2	1:A:88:LYS:HB3	1.95	0.49
1:B:217:THR:HG23	2:D:349:VAL:HG22	1.93	0.49
1:B:72:LYS:N	1:B:72:LYS:HD2	2.28	0.49
1:B:49:THR:HG22	1:B:51:GLY:N	2.28	0.49
1:B:21:GLU:HB2	1:B:26:GLN:O	2.13	0.49
1:B:73:VAL:HG12	1:B:74:GLU:H	1.78	0.48
1:B:277:ILE:HB	1:B:285:CYS:HA	1.94	0.48
1:B:274:LEU:HD13	1:B:285:CYS:SG	2.54	0.48
1:B:32:ILE:HB	1:B:123:ILE:HG12	1.96	0.47
1:A:80:GLY:HA3	1:A:111:PHE:CD1	2.48	0.47
1:A:273:TYR:CE1	1:A:287:LEU:HD13	2.48	0.47
1:B:175:GLU:HA	1:B:178:TYR:CE1	2.49	0.47
1:B:140:VAL:HG21	1:B:154:PHE:CD2	2.49	0.47
1:B:73:VAL:HG12	1:B:74:GLU:N	2.29	0.47
1:B:4:ASP:OD1	1:B:96:LEU:HG	2.14	0.47
1:A:313:PHE:N	1:A:313:PHE:CD1	2.81	0.47
1:A:153:LEU:O	1:A:153:LEU:HD12	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:201:VAL:HG12	1:A:255:PRO:HB2	1.95	0.47
1:B:155:THR:HG23	1:B:156:PHE:N	2.29	0.47
1:A:137:ASP:CB	1:A:142:GLU:HG2	2.45	0.47
1:A:260:THR:HG22	1:A:265:LYS:HG3	1.97	0.47
1:B:92:THR:OG1	1:B:97:SER:HB2	2.15	0.47
1:A:303:ASP:O	1:A:307:ARG:HB2	2.14	0.47
1:B:15:MET:HG3	1:B:16:PHE:H	1.79	0.47
1:B:4:ASP:HB2	1:B:94:GLY:HA3	1.96	0.47
1:B:290:ILE:HD11	2:D:347:IVA:HG21	1.97	0.47
1:A:174:GLU:OE1	1:A:176:ARG:HD3	2.14	0.47
1:B:153:LEU:HD12	1:B:153:LEU:O	2.15	0.47
1:B:15:MET:HG3	1:B:16:PHE:N	2.30	0.46
1:A:201:VAL:CG1	1:A:255:PRO:HB2	2.45	0.46
1:A:201:VAL:HG23	1:A:226:PHE:CZ	2.50	0.46
1:A:40:LEU:O	1:A:103:ILE:HD12	2.15	0.46
1:A:14:ILE:HD11	1:A:275:GLN:CD	2.36	0.46
1:A:221:THR:CG2	1:A:292:LEU:HB3	2.41	0.46
1:B:24:ASP:OD2	1:B:65:THR:HG23	2.15	0.46
1:B:77:TYR:CB	2:D:350:STA:HD13	2.46	0.46
1:B:27:GLN:OE1	1:B:58:TYR:CD1	2.69	0.46
1:B:10:ASP:OD1	1:B:13:ASN:HA	2.16	0.46
1:B:272:TYR:CD1	1:B:308:LYS:HB2	2.50	0.46
1:A:191:LEU:HA	1:A:191:LEU:HD12	1.71	0.46
1:B:164:HIS:HD2	1:B:327:LYS:CG	2.25	0.45
1:B:100:TYR:CE1	1:B:139:ILE:HG12	2.51	0.45
1:B:51:GLY:O	1:B:54:THR:HB	2.15	0.45
1:A:188:ASN:HD22	1:A:189:HIS:CD2	2.33	0.45
1:B:13:ASN:CG	1:B:161:HIS:HB2	2.36	0.45
1:B:47:CYS:SG	1:B:49:THR:HB	2.56	0.45
1:B:273:TYR:O	1:B:288:ASN:HB3	2.16	0.45
1:B:303:ASP:O	1:B:307:ARG:N	2.46	0.45
1:B:64:ARG:CD	1:B:64:ARG:H	2.28	0.45
1:B:70:GLY:N	1:B:86:PHE:O	2.50	0.45
1:B:41:TRP:HE1	1:B:77:TYR:HH	1.63	0.45
1:A:49:THR:HG22	1:A:51:GLY:H	1.82	0.45
1:A:55:LYS:HG3	1:A:115:TYR:OH	2.17	0.45
1:B:294:PHE:CZ	1:B:300:ILE:HD11	2.50	0.44
1:A:231:LEU:HD22	1:A:236:VAL:HG11	1.98	0.44
1:B:277:ILE:HG13	1:B:286:MET:CE	2.46	0.44
1:B:5:ASN:HB3	1:B:167:PHE:CD1	2.53	0.44
1:A:146:GLN:HB3	1:A:148:LYS:HZ2	1.82	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:34:ASP:OD2	2:D:350:STA:HC	2.17	0.44
1:B:203:ASN:C	1:B:205:MET:SD	2.96	0.44
1:A:146:GLN:O	1:A:148:LYS:N	2.51	0.44
1:B:140:VAL:HG21	1:B:154:PHE:HD2	1.83	0.44
1:A:55:LYS:HE3	1:A:115:TYR:CE1	2.52	0.44
1:A:65:THR:HG22	1:A:65:THR:O	2.17	0.44
1:B:55:LYS:HG3	1:B:115:TYR:CZ	2.52	0.44
1:B:155:THR:HG1	1:B:310:PHE:HZ	1.64	0.44
1:A:8:LEU:HD13	1:A:16:PHE:CE1	2.53	0.44
1:B:41:TRP:HB2	1:B:105:VAL:CG2	2.48	0.44
1:B:77:TYR:HB3	2:D:350:STA:HA	2.00	0.44
1:B:2:SER:CB	1:B:149:ILE:HA	2.48	0.44
1:A:94:GLY:O	1:A:95:ASN:HB3	2.18	0.44
1:B:220:ILE:HD13	1:B:220:ILE:HA	1.82	0.44
1:A:274:LEU:HD22	1:A:285:CYS:HB3	1.99	0.43
1:A:146:GLN:HB3	1:A:148:LYS:NZ	2.33	0.43
1:A:287:LEU:HD12	1:A:289:ILE:HD12	1.99	0.43
1:B:228:ASN:HA	1:B:231:LEU:HD12	1.99	0.43
1:A:260:THR:CG2	1:A:265:LYS:HG3	2.48	0.43
1:A:68:LYS:HB2	1:A:68:LYS:HE3	1.51	0.43
1:A:44:SER:OG	1:A:46:LYS:HD3	2.18	0.43
1:B:25:ASN:O	1:B:26:GLN:CB	2.65	0.43
1:A:201:VAL:HG21	1:A:226:PHE:HE1	1.81	0.43
1:B:9:VAL:HG12	1:B:17:TYR:O	2.19	0.43
1:A:251:ASN:HD22	1:A:252:SER:N	2.16	0.43
2:D:352:STA:HD22	2:D:352:STA:HA	1.79	0.43
1:A:155:THR:HG1	1:A:310:PHE:HZ	1.65	0.43
1:A:97:SER:O	1:A:148:LYS:HE2	2.19	0.43
1:A:1:SER:C	1:A:3:ASN:H	2.20	0.43
1:B:200:HIS:HB2	1:B:258:GLU:HB2	1.99	0.43
1:A:69:ASP:HB3	1:A:86:PHE:O	2.19	0.43
1:B:44:SER:CB	1:B:104:GLU:HG2	2.35	0.43
1:A:74:GLU:HG3	1:A:81:THR:CG2	2.49	0.43
1:B:108:THR:HG21	1:B:115:TYR:CE2	2.53	0.43
1:A:294:PHE:CE1	1:A:300:ILE:HD11	2.53	0.43
1:B:22:VAL:HG22	1:B:91:VAL:HG22	2.01	0.43
1:B:253:LYS:HD3	1:B:253:LYS:HA	1.73	0.43
1:A:22:VAL:HG12	1:A:23:GLY:N	2.34	0.43
1:B:248:LEU:HA	1:B:248:LEU:HD12	1.89	0.43
1:B:17:TYR:CE1	1:B:120:PHE:HB3	2.54	0.42
1:B:201:VAL:HG13	1:B:255:PRO:HB2	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:144:LYS:HE3	1:B:151:ASN:HA	2.00	0.42
1:B:51:GLY:O	1:B:55:LYS:HG2	2.19	0.42
1:B:55:LYS:HG3	1:B:115:TYR:OH	2.19	0.42
1:B:182:LEU:HA	1:B:323:ALA:HB2	2.01	0.42
1:B:182:LEU:HD12	1:B:323:ALA:CB	2.49	0.42
1:A:91:VAL:HG21	1:A:100:TYR:HB3	2.01	0.42
1:B:215:SER:HB3	1:B:306:MET:SD	2.58	0.42
1:A:91:VAL:CG2	1:A:100:TYR:HB3	2.49	0.42
1:A:119:THR:HG22	1:A:119:THR:O	2.19	0.42
1:B:34:ASP:O	1:B:126:LEU:N	2.52	0.42
1:B:137:ASP:HA	1:B:138:PRO:HD2	1.62	0.42
1:B:55:LYS:HD2	1:B:120:PHE:O	2.19	0.42
1:B:222:VAL:HG21	1:B:226:PHE:CD1	2.55	0.42
1:A:227:LEU:HD11	1:A:231:LEU:HD11	2.02	0.42
1:B:3:ASN:HA	1:B:170:ILE:O	2.19	0.42
1:A:157:TYR:HD2	1:A:166:GLY:HA2	1.84	0.42
1:A:149:ILE:O	1:A:149:ILE:HG13	2.20	0.42
1:B:175:GLU:O	1:B:175:GLU:HG2	2.20	0.42
1:A:275:GLN:HB2	1:A:286:MET:HG3	2.02	0.42
1:A:251:ASN:C	1:A:251:ASN:HD22	2.22	0.42
1:A:155:THR:CG2	1:A:310:PHE:HE1	2.31	0.41
1:A:108:THR:HG21	1:A:115:TYR:CE2	2.55	0.41
1:B:290:ILE:HG13	2:D:347:IVA:HG11	2.02	0.41
1:B:270:PRO:HA	1:B:273:TYR:CE2	2.56	0.41
1:B:11:PHE:O	1:B:13:ASN:N	2.54	0.41
1:B:184:TYR:HD1	1:B:319:SER:HB3	1.85	0.41
1:B:129:LYS:HD2	1:B:137:ASP:OD2	2.21	0.41
1:A:294:PHE:HB3	1:A:295:PRO:CD	2.51	0.41
1:A:144:LYS:O	1:A:145:ASN:C	2.59	0.41
1:B:278:GLU:OE2	1:B:282:PRO:HA	2.21	0.41
1:B:246:VAL:HG22	1:B:286:MET:HE1	2.03	0.41
1:B:78:VAL:HG12	1:B:78:VAL:O	2.21	0.41
1:B:194:GLN:HG2	1:B:194:GLN:H	1.72	0.41
1:B:27:GLN:OE1	1:B:58:TYR:HD1	2.03	0.41
1:A:16:PHE:CE2	1:A:158:LEU:HD22	2.56	0.41
1:A:221:THR:HB	1:A:300:ILE:CB	2.41	0.41
1:A:221:THR:HG21	1:A:292:LEU:HD23	2.02	0.41
1:B:303:ASP:O	1:B:306:MET:N	2.54	0.41
1:B:144:LYS:CB	1:B:152:ALA:HB2	2.50	0.41
1:B:6:ILE:CD1	1:B:93:VAL:HG12	2.47	0.41
1:A:83:SER:O	1:A:105:VAL:HG13	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:234:LEU:HA	1:A:253:LYS:HD2	2.02	0.41
1:B:100:TYR:CZ	1:B:139:ILE:HG12	2.56	0.41
1:B:277:ILE:HA	1:B:277:ILE:HD13	1.68	0.40
1:A:188:ASN:O	1:A:189:HIS:CG	2.74	0.40
1:B:290:ILE:CD1	2:D:347:IVA:HG21	2.51	0.40
1:A:22:VAL:HG11	1:A:102:PHE:CE2	2.56	0.40
1:B:294:PHE:HB2	1:B:296:VAL:O	2.22	0.40
1:A:29:PHE:HB2	1:A:31:PHE:CE1	2.57	0.40
1:A:278:GLU:O	1:A:278:GLU:HG3	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	327/329 (99%)	289 (88%)	30 (9%)	8 (2%)	7 19
1	B	327/329 (99%)	288 (88%)	29 (9%)	10 (3%)	5 12
2	C	3/6 (50%)	3 (100%)	0	0	100 100
2	D	3/6 (50%)	3 (100%)	0	0	100 100
All	All	660/670 (98%)	583 (88%)	59 (9%)	18 (3%)	6 16

All (18) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	95	ASN
1	A	161	HIS
1	B	161	HIS
1	A	69	ASP
1	A	147	ASN
1	A	297	PRO

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Mol	Chain	Res	Type
1	B	12	GLN
1	B	49	THR
1	A	4	ASP
1	B	25	ASN
1	B	26	GLN
1	B	61	SER
1	B	70	GLY
1	A	5	ASN
1	A	162	ASP
1	B	27	GLN
1	B	138	PRO
1	B	297	PRO

5.3.2 Protein sidechains (i)

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	294/294 (100%)	269 (92%)	25 (8%)	13 30
1	B	294/294 (100%)	266 (90%)	28 (10%)	11 24
2	C	2/2 (100%)	2 (100%)	0	100 100
2	D	2/2 (100%)	2 (100%)	0	100 100
All	All	592/592 (100%)	539 (91%)	53 (9%)	12 27

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	SER
1	A	24	ASP
1	A	42	VAL
1	A	45	VAL
1	A	46	LYS
1	A	71	THR
1	A	73	VAL
1	A	82	VAL
1	A	98	LEU

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Mol	Chain	Res	Type
1	A	137	ASP
1	A	144	LYS
1	A	147	ASN
1	A	150	GLU
1	A	155	THR
1	A	165	THR
1	A	188	ASN
1	A	203	ASN
1	A	206	LEU
1	A	233	ASN
1	A	247	THR
1	A	251	ASN
1	A	252	SER
1	A	271	GLU
1	A	287	LEU
1	A	297	PRO
1	B	2	SER
1	B	4	ASP
1	B	6	ILE
1	B	21	GLU
1	B	26	GLN
1	B	32	ILE
1	B	60	SER
1	B	64	ARG
1	B	72	LYS
1	B	90	LEU
1	B	97	SER
1	B	112	GLU
1	B	137	ASP
1	B	142	GLU
1	B	155	THR
1	B	156	PHE
1	B	174	GLU
1	B	187	LEU
1	B	198	ASP
1	B	205	MET
1	B	225	ASP
1	B	233	ASN
1	B	279	ASP
1	B	282	PRO
1	B	284	LEU
1	B	288	ASN

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Mol	Chain	Res	Type
1	B	326	LYS
1	B	328	ASN

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

Mol	Chain	Res	Type
1	A	189	HIS
1	A	194	GLN
1	A	233	ASN
1	A	251	ASN
1	B	3	ASN
1	B	5	ASN
1	B	25	ASN
1	B	200	HIS
1	B	275	GLN
1	B	288	ASN

5.3.3 RNA

There are no RNA molecules in this entry.

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

4 non-standard protein/DNA/RNA residues are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
2	STA	C	350	2	10,10,11	0.95	1 (10%)	10,12,14	1.42	2 (20%)
2	STA	C	352	2	8,11,11	2.09	2 (25%)	8,14,14	1.66	2 (25%)
2	STA	D	350	2	10,10,11	1.55	2 (20%)	10,12,14	1.54	2 (20%)
2	STA	D	352	2	8,11,11	0.94	0	8,14,14	1.09	1 (12%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	STA	C	350	2	-	0/11/11/12	0/0/0/0
2	STA	C	352	2	-	0/10/12/12	0/0/0/0
2	STA	D	350	2	-	0/11/11/12	0/0/0/0
2	STA	D	352	2	-	0/10/12/12	0/0/0/0

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	350	STA	OH-CH	2.50	1.49	1.43
2	C	350	STA	CH-CA	2.79	1.56	1.53
2	C	352	STA	OH-CH	3.60	1.51	1.43
2	D	350	STA	CH-CA	3.96	1.57	1.53
2	C	352	STA	CH-CA	4.58	1.57	1.53

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	C	350	STA	OH-CH-CM	-3.12	102.09	109.34
2	C	352	STA	CM-CH-CA	-2.53	108.83	112.68
2	C	350	STA	CB-CA-N	-2.21	102.77	109.13
2	D	350	STA	CD2-CG-CD1	-2.16	99.69	110.55
2	D	352	STA	OH-CH-CM	2.54	114.14	109.73
2	C	352	STA	OH-CH-CM	3.13	115.18	109.73
2	D	350	STA	CG-CB-CA	3.68	123.90	115.77

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	350	STA	1	0
2	C	352	STA	1	0
2	D	350	STA	4	0
2	D	352	STA	2	0

5.5 Carbohydrates

There are no carbohydrates in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates

EDS was not executed - this section will therefore be empty.

6.4 Ligands

EDS was not executed - this section will therefore be empty.

6.5 Other polymers

EDS was not executed - this section will therefore be empty.