wwPDB EM Map/Model Validation Report (i)

May 2, 2016 – 10:46 PM EDT

PDB ID	:	3EPF
EMDB ID:	:	EMD-1563
Title	:	CryoEM structure of poliovirus receptor bound to poliovirus type 2
Authors	:	Zhang, P.; Mueller, S.; Morais, M.C.; Bator, C.M.; Bowman, V.D.; Hafenstein, S.; Wimmer, E.; Rossmann, M.G.
Deposited on	:	2008-09-29
Resolution	:	9.00 Å(reported)

This is a wwPDB EM Map/Model Validation Report for a publicly released PDB/EMDB entry.
For rigid body fitted models, validation errors reported here could
stem from errors in the original structure(s) used in the fitting.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
http://wwpdb.org/validation/2016/EMValidationReportHelp

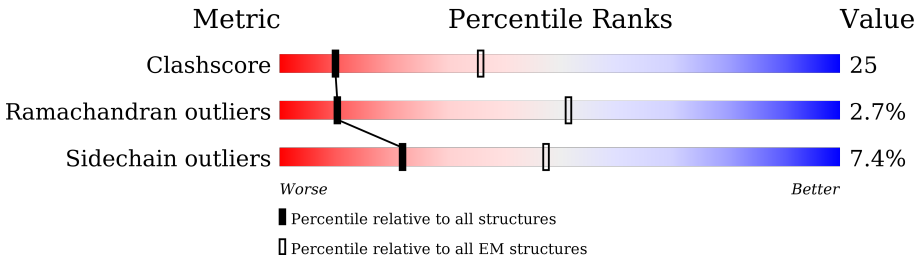
MolProbity	:	4.02b-467
Mogul	:	1.7.1 (RC1), CSD as537be (2016)
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP)	:	rb-20027457

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
ELECTRON MICROSCOPY

The reported resolution of this entry is 9.00 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)	EM structures (#Entries)
Clashscore	114402	924
Ramachandran outliers	111179	726
Sidechain outliers	111093	686

The table below summarises the geometric issues observed across the polymeric chains. The red, orange, yellow and green segments on the 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%

Mol	Chain	Length	Quality of chain
1	R	213	36% 43% 16% .
2	1	278	73% 24% .
3	2	262	83% 15% .
4	4	68	78% 18% .
5	3	235	87% 10% .

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 8461 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, 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 Poliovirus receptor.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	R	213	Total	C	N	O	S	0	0
			1638	1038	281	310	9		

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
R	105	ASP	ASN	ENGINEERED	UNP P15151
R	120	SER	ASN	ENGINEERED	UNP P15151
R	188	GLN	ASN	ENGINEERED	UNP P15151
R	218	GLN	ASN	ENGINEERED	UNP P15151
R	237	SER	ASN	ENGINEERED	UNP P15151

- Molecule 2 is a protein called Protein VP1.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	1	272	Total	C	N	O	S	0	0
			2129	1361	363	400	5		

- Molecule 3 is a protein called Protein VP2.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	2	262	Total	C	N	O	S	0	0
			2042	1298	346	384	14		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
2	11	VAL	ASP	CONFLICT	UNP P06210

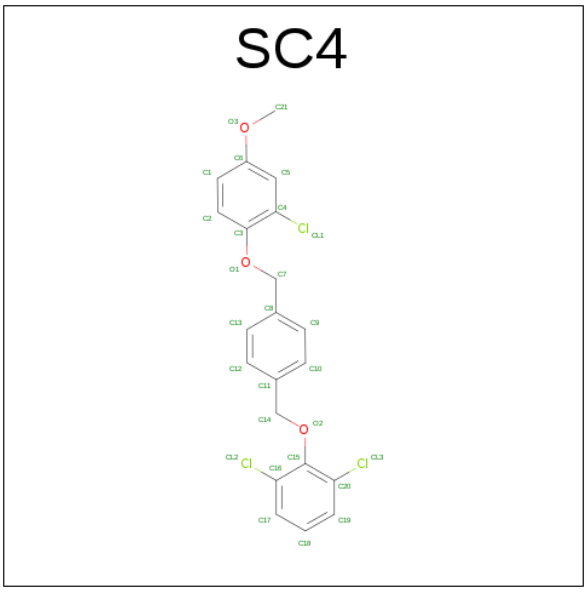
- Molecule 4 is a protein called Protein VP4.

Mol	Chain	Residues	Atoms				AltConf	Trace
4	4	68	Total 518	C 318	N 91	O 109	0	0

- Molecule 5 is a protein called Protein VP3.

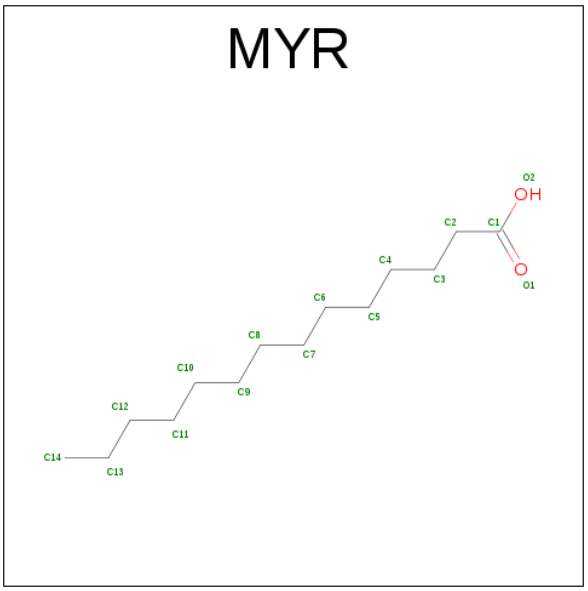
Mol	Chain	Residues	Atoms					AltConf	Trace
5	3	235	Total 1825	C 1156	N 303	O 349	S 17	0	0

- Molecule 6 is 1[2-CHLORO-4-METHOXY-PHENYL-OXYMETHYL]-4-[2,6-DICHLORO-PHENYL-OXYMETHYL]-BENZENE (three-letter code: SC4) (formula: $C_{21}H_{17}Cl_3O_3$).



Mol	Chain	Residues	Atoms				AltConf
6	1	1	Total 27	C 21	Cl 3	O 3	0

- Molecule 7 is MYRISTIC ACID (three-letter code: MYR) (formula: $C_{14}H_{28}O_2$).



Mol	Chain	Residues	Atoms			AltConf
7	4	1	Total	C	O	0
			11	10	1	

- Molecule 8 is water.

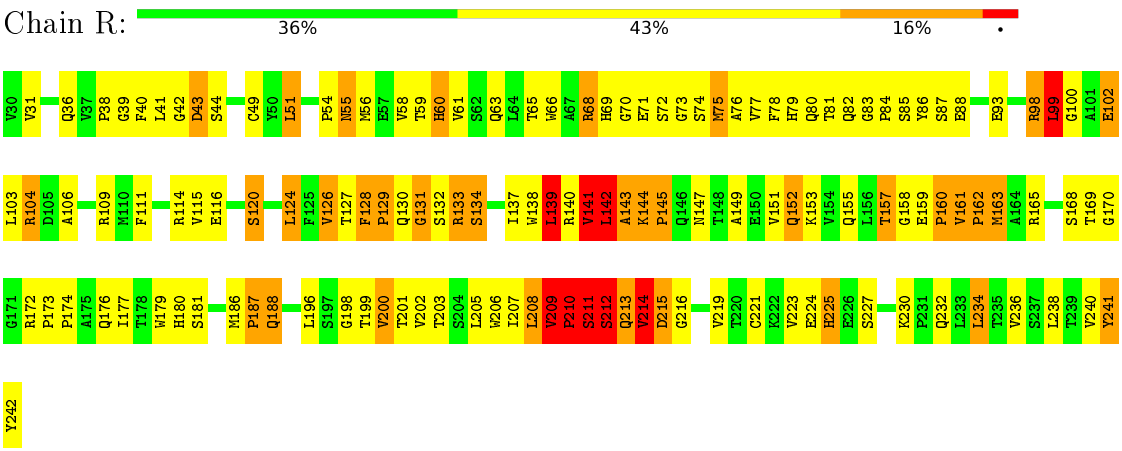
Mol	Chain	Residues	Atoms		AltConf
8	1	136	Total	O	0
			136	136	
8	2	95	Total	O	0
			95	95	
8	4	25	Total	O	0
			25	25	
8	3	15	Total	O	0
			15	15	

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. 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. 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.

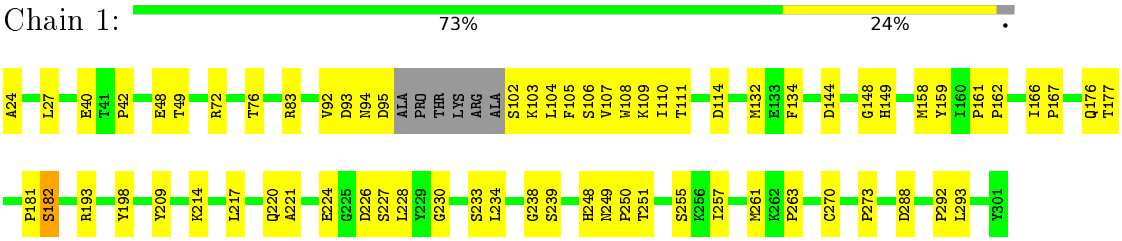
- Molecule 1: Poliovirus receptor

Chain R:



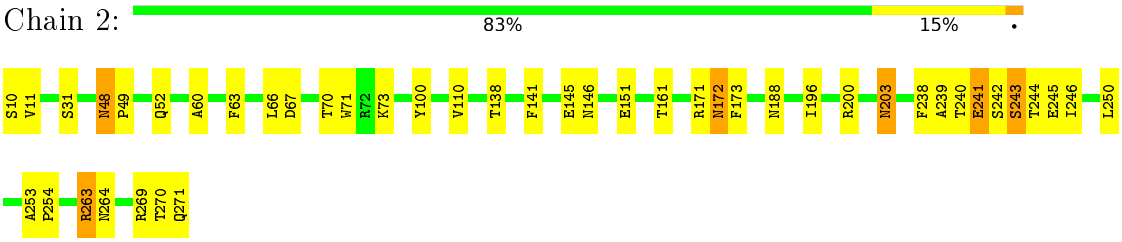
- Molecule 2: Protein VP1

Chain 1:



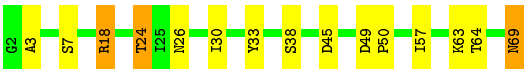
- Molecule 3: Protein VP2

Chain 2:



- Molecule 4: Protein VP4

Chain 4:



- Molecule 5: Protein VP3

Chain 3:

4 Experimental information (i)

Property	Value	Source
Reconstruction method	Not provided	Depositor
Imposed symmetry	POINT, I	Depositor
Number of images	Not provided	Depositor
Resolution determination method	Not provided	Depositor
CTF correction method	Not provided	Depositor
Microscope	Not provided	Depositor
Voltage (kV)	Not provided	Depositor
Electron dose ($e^-/Å^2$)	Not provided	Depositor
Minimum defocus (nm)	Not provided	Depositor
Maximum defocus (nm)	Not provided	Depositor
Magnification	Not provided	Depositor
Image detector	Not provided	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: SC4, MYR

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	$\#\|Z\| > 2$	RMSZ	$\#\|Z\| > 2$
1	R	0.41	0/1678	0.77	4/2289 (0.2%)
2	1	0.43	0/2193	0.69	1/2996 (0.0%)
3	2	0.44	0/2098	0.72	1/2863 (0.0%)
4	4	0.52	0/527	0.68	0/713
5	3	0.43	0/1869	0.67	1/2549 (0.0%)
All	All	0.44	0/8365	0.71	7/11410 (0.1%)

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
1	R	1	6
2	1	0	1
All	All	1	7

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	R	211	SER	O-C-N	-8.56	109.01	122.70
1	R	99	LEU	CA-CB-CG	6.25	129.67	115.30
2	1	288	ASP	N-CA-C	5.95	127.06	111.00
1	R	211	SER	CA-C-N	5.90	130.19	117.20
3	2	203	ASN	CB-CA-C	-5.76	98.89	110.40
1	R	209	VAL	N-CA-C	5.28	125.25	111.00
5	3	5	LEU	CA-CB-CG	5.19	127.24	115.30

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	R	143	ALA	CA

All (7) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	1	198	TYR	Sidechain
1	R	141	VAL	Peptide
1	R	142	LEU	Peptide
1	R	210	PRO	Peptide
1	R	211	SER	Peptide
1	R	212	SER	Mainchain
1	R	213	GLN	Peptide

5.2 Too-close contacts (i)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	R	1638	0	1617	322	0
2	1	2129	0	2068	155	0
3	2	2042	0	1969	36	0
4	4	518	0	495	12	0
5	3	1825	0	1800	54	0
6	1	27	0	17	1	0
7	4	11	0	16	0	0
8	1	136	0	0	8	0
8	2	95	0	0	1	0
8	3	15	0	0	1	0
8	4	25	0	0	1	0
All	All	8461	0	7982	410	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 25.

All (410) 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:R:130:GLN:CG	2:1:106:SER:HA	1.13	1.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:R:130:GLN:CG	2:1:106:SER:CA	1.86	1.54
1:R:115:VAL:HG11	1:R:142:LEU:CD1	1.28	1.53
1:R:130:GLN:NE2	2:1:105:PHE:CE1	1.76	1.52
1:R:115:VAL:CG1	1:R:142:LEU:HD11	1.03	1.50
1:R:130:GLN:CB	2:1:106:SER:HA	1.38	1.49
1:R:115:VAL:CG1	1:R:142:LEU:CD1	1.79	1.48
1:R:130:GLN:HG3	2:1:106:SER:CA	1.42	1.43
1:R:128:PHE:HE2	2:1:108:TRP:NE1	1.17	1.41
1:R:83:GLY:CA	2:1:226:ASP:O	1.70	1.37
1:R:41:LEU:CB	1:R:143:ALA:HB3	1.59	1.33
1:R:41:LEU:HB3	1:R:143:ALA:CB	1.60	1.31
1:R:128:PHE:CE2	2:1:108:TRP:NE1	2.02	1.25
1:R:99:LEU:HG	2:1:226:ASP:OD2	1.37	1.24
1:R:128:PHE:CE1	2:1:114:ASP:OD1	1.69	1.22
1:R:132:SER:OG	2:1:107:VAL:HG11	1.39	1.22
1:R:130:GLN:HB2	2:1:106:SER:CA	1.72	1.20
1:R:128:PHE:CD2	8:1:397:HOH:O	1.65	1.18
1:R:40:PHE:CZ	1:R:144:LYS:HB3	1.79	1.18
1:R:130:GLN:CA	2:1:107:VAL:H	1.58	1.16
1:R:130:GLN:HG2	2:1:106:SER:C	1.66	1.15
1:R:162:PRO:HD2	1:R:163:MET:HA	1.30	1.14
1:R:128:PHE:CG	8:1:397:HOH:O	1.65	1.14
1:R:161:VAL:HB	1:R:163:MET:HB2	1.21	1.13
1:R:99:LEU:CG	2:1:226:ASP:OD2	1.94	1.13
1:R:43:ASP:HB2	1:R:44:SER:CA	1.79	1.12
1:R:130:GLN:HB2	2:1:106:SER:CB	1.79	1.12
1:R:130:GLN:HA	2:1:107:VAL:H	0.95	1.10
1:R:115:VAL:HG13	1:R:142:LEU:HD12	1.34	1.09
1:R:132:SER:OG	2:1:107:VAL:CG1	1.99	1.08
1:R:82:GLN:HB2	2:1:228:LEU:H	1.16	1.08
1:R:142:LEU:O	1:R:173:PRO:CD	2.01	1.08
1:R:83:GLY:HA3	2:1:226:ASP:C	1.74	1.08
1:R:115:VAL:HG13	1:R:142:LEU:CD1	1.68	1.07
1:R:43:ASP:CB	1:R:44:SER:HA	1.85	1.06
1:R:82:GLN:CB	2:1:228:LEU:H	1.68	1.06
1:R:73:GLY:N	5:3:97:HIS:NE2	2.05	1.05
1:R:130:GLN:HA	2:1:107:VAL:N	1.71	1.04
1:R:130:GLN:HG3	2:1:106:SER:N	1.71	1.04
1:R:129:PRO:O	2:1:107:VAL:O	1.76	1.04
1:R:98:ARG:HD3	1:R:104:ARG:HH21	1.21	1.03
1:R:132:SER:CB	2:1:166:ILE:HD13	1.89	1.03

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:R:79:HIS:CD2	2:1:234:LEU:HD12	1.94	1.02
1:R:98:ARG:HD3	1:R:104:ARG:NH2	1.75	1.01
1:R:130:GLN:HG2	2:1:107:VAL:N	1.74	1.01
1:R:83:GLY:N	2:1:227:SER:HA	1.74	1.01
1:R:130:GLN:CB	2:1:106:SER:CA	2.16	1.01
1:R:132:SER:HB3	2:1:166:ILE:HD13	1.40	0.99
1:R:116:GLU:CD	5:3:59:ASN:ND2	2.16	0.99
1:R:54:PRO:HA	1:R:55:ASN:HB2	1.43	0.99
1:R:132:SER:CB	2:1:107:VAL:HG11	1.92	0.99
1:R:82:GLN:CB	2:1:228:LEU:N	2.26	0.99
1:R:177:ILE:HD12	1:R:205:LEU:HB2	1.45	0.99
1:R:83:GLY:HA3	2:1:226:ASP:O	0.80	0.97
1:R:73:GLY:CA	5:3:97:HIS:NE2	2.28	0.97
1:R:73:GLY:HA3	5:3:97:HIS:CD2	2.00	0.96
1:R:128:PHE:CE2	8:1:397:HOH:O	1.96	0.96
1:R:116:GLU:CD	5:3:59:ASN:HD21	1.66	0.96
1:R:40:PHE:CZ	1:R:144:LYS:CB	2.47	0.96
1:R:82:GLN:HB3	2:1:228:LEU:N	1.82	0.95
1:R:130:GLN:HB2	2:1:106:SER:HA	1.32	0.95
1:R:116:GLU:HA	5:3:59:ASN:ND2	1.81	0.95
1:R:143:ALA:O	1:R:225:HIS:NE2	1.99	0.94
1:R:128:PHE:CD1	8:1:397:HOH:O	1.93	0.94
1:R:116:GLU:CG	5:3:59:ASN:HD21	1.80	0.94
1:R:73:GLY:HA3	5:3:97:HIS:NE2	1.82	0.94
1:R:130:GLN:CG	2:1:106:SER:C	2.27	0.93
1:R:68:ARG:HD2	1:R:76:ALA:HB3	1.52	0.92
1:R:79:HIS:HD2	2:1:234:LEU:HD12	1.30	0.90
1:R:83:GLY:H	2:1:227:SER:HA	1.32	0.90
1:R:142:LEU:O	1:R:173:PRO:HD3	1.71	0.90
1:R:40:PHE:HZ	1:R:144:LYS:HB3	1.33	0.90
1:R:86:TYR:CE1	2:1:224:GLU:OE1	2.25	0.89
1:R:162:PRO:CD	1:R:163:MET:HA	2.03	0.88
1:R:141:VAL:HG13	1:R:142:LEU:N	1.89	0.88
1:R:130:GLN:HG3	2:1:105:PHE:C	1.94	0.87
1:R:43:ASP:HB2	1:R:44:SER:HA	0.91	0.87
1:R:128:PHE:CZ	2:1:111:THR:CG2	2.58	0.87
1:R:128:PHE:CD2	2:1:108:TRP:CD1	2.63	0.86
1:R:128:PHE:O	2:1:108:TRP:HA	1.74	0.86
1:R:130:GLN:HG3	2:1:106:SER:HA	0.93	0.85
1:R:99:LEU:CD1	2:1:226:ASP:OD2	2.24	0.85
1:R:141:VAL:HG13	1:R:142:LEU:HB2	1.56	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:R:128:PHE:CE2	2:1:108:TRP:CD1	2.63	0.85
1:R:40:PHE:HZ	1:R:144:LYS:CB	1.85	0.85
1:R:130:GLN:NE2	2:1:105:PHE:HE1	1.44	0.85
1:R:141:VAL:HG13	1:R:142:LEU:CB	2.07	0.84
1:R:82:GLN:NE2	2:1:228:LEU:HD13	1.19	0.84
2:1:24:ALA:HB2	2:1:76:THR:HG21	1.58	0.84
3:2:11:VAL:HA	8:4:145:HOH:O	1.78	0.84
1:R:128:PHE:HE1	2:1:114:ASP:OD1	1.55	0.84
1:R:81:THR:N	2:1:234:LEU:HD11	1.93	0.83
1:R:81:THR:N	2:1:234:LEU:CD1	2.41	0.83
1:R:81:THR:HG21	2:1:234:LEU:HA	1.60	0.83
1:R:116:GLU:OE1	5:3:59:ASN:ND2	2.11	0.82
2:1:176:GLN:HE22	5:3:233:GLN:NE2	1.77	0.82
1:R:54:PRO:CA	1:R:55:ASN:HB2	2.11	0.81
1:R:142:LEU:O	1:R:173:PRO:CG	2.28	0.80
2:1:148:GLY:HA3	2:1:251:THR:HG23	1.63	0.80
1:R:130:GLN:HB2	2:1:106:SER:HB3	1.61	0.80
1:R:142:LEU:HD23	1:R:172:ARG:HG2	1.63	0.80
1:R:79:HIS:CD2	2:1:234:LEU:CD1	2.65	0.80
1:R:130:GLN:CB	2:1:107:VAL:H	1.94	0.79
2:1:233:SER:HB3	8:3:244:HOH:O	1.82	0.79
1:R:130:GLN:CG	2:1:107:VAL:N	2.44	0.78
1:R:128:PHE:CZ	8:1:397:HOH:O	2.22	0.78
1:R:81:THR:CG2	2:1:234:LEU:HA	2.10	0.77
1:R:81:THR:H	2:1:234:LEU:HD11	1.48	0.76
1:R:215:ASP:O	1:R:238:LEU:HB3	1.85	0.76
3:2:263:ARG:HG2	5:3:136:PRO:HD2	1.68	0.76
3:2:48:ASN:HB3	3:2:49:PRO:HD3	1.66	0.76
1:R:63:GLN:HG2	1:R:81:THR:HA	1.69	0.75
1:R:130:GLN:HE21	2:1:107:VAL:HG23	1.51	0.75
1:R:82:GLN:HE22	2:1:228:LEU:CD1	1.46	0.75
1:R:173:PRO:HG2	1:R:225:HIS:HE1	1.52	0.74
1:R:129:PRO:C	2:1:107:VAL:O	2.25	0.74
1:R:41:LEU:CA	1:R:143:ALA:HB3	2.17	0.74
1:R:86:TYR:HE1	2:1:224:GLU:OE1	1.71	0.74
1:R:130:GLN:NE2	2:1:105:PHE:CZ	2.53	0.73
1:R:130:GLN:HE21	2:1:105:PHE:HE1	1.15	0.73
1:R:130:GLN:NE2	2:1:105:PHE:CD1	2.53	0.73
1:R:99:LEU:HD12	2:1:226:ASP:OD2	1.87	0.73
1:R:82:GLN:NE2	2:1:228:LEU:HD12	1.21	0.73
5:3:51:THR:HG21	5:3:99:MET:H	1.53	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:R:100:GLY:CA	3:2:138:THR:HG21	2.19	0.72
1:R:80:GLN:HG2	1:R:98:ARG:HG2	1.71	0.72
2:1:177:THR:HG21	2:1:182:SER:OG	1.89	0.72
1:R:142:LEU:CD2	1:R:172:ARG:HG2	2.20	0.71
1:R:130:GLN:CB	2:1:107:VAL:N	2.52	0.71
1:R:165:ARG:HG3	1:R:206:TRP:HB3	1.73	0.71
1:R:41:LEU:HB3	1:R:143:ALA:HB3	0.77	0.71
1:R:130:GLN:CA	2:1:107:VAL:N	2.42	0.71
1:R:82:GLN:HE21	2:1:228:LEU:CD1	1.36	0.71
1:R:128:PHE:HB3	1:R:129:PRO:HD3	1.70	0.71
1:R:132:SER:HB2	2:1:107:VAL:HG11	1.73	0.71
1:R:124:LEU:HA	1:R:134:SER:HB2	1.73	0.70
1:R:128:PHE:HZ	2:1:111:THR:CG2	2.04	0.70
1:R:141:VAL:HG11	1:R:142:LEU:HD13	1.72	0.70
1:R:40:PHE:HZ	1:R:144:LYS:CD	2.05	0.70
1:R:40:PHE:HZ	1:R:144:LYS:HD3	1.56	0.70
1:R:75:MET:HE2	5:3:92:ASP:OD2	1.76	0.70
1:R:116:GLU:CA	5:3:59:ASN:ND2	2.55	0.70
1:R:40:PHE:CE1	1:R:227:SER:OG	2.44	0.70
1:R:68:ARG:HD2	1:R:76:ALA:CB	2.21	0.70
2:1:109:LYS:HA	2:1:239:SER:HB3	1.73	0.69
1:R:104:ARG:HD2	1:R:106:ALA:HB2	1.74	0.69
1:R:130:GLN:HG3	2:1:105:PHE:O	1.90	0.69
1:R:141:VAL:HG13	1:R:142:LEU:CA	2.22	0.69
1:R:116:GLU:CA	5:3:59:ASN:HD21	2.05	0.69
1:R:132:SER:OG	2:1:107:VAL:HG12	1.92	0.69
1:R:128:PHE:CZ	2:1:111:THR:HG21	2.26	0.69
1:R:72:SER:C	5:3:97:HIS:NE2	2.46	0.68
1:R:100:GLY:HA2	3:2:138:THR:HG21	1.75	0.68
1:R:116:GLU:HA	5:3:59:ASN:HD21	1.55	0.68
1:R:82:GLN:HE22	2:1:228:LEU:HD11	1.00	0.67
1:R:128:PHE:CE1	8:1:397:HOH:O	2.20	0.67
1:R:173:PRO:HG2	1:R:225:HIS:CE1	2.30	0.67
1:R:40:PHE:CE1	1:R:227:SER:CB	2.77	0.67
1:R:142:LEU:HD23	1:R:172:ARG:CG	2.24	0.67
1:R:143:ALA:O	1:R:225:HIS:CE1	2.47	0.67
1:R:82:GLN:HB2	2:1:228:LEU:N	1.97	0.67
1:R:98:ARG:HG3	1:R:99:LEU:H	1.60	0.66
3:2:145:GLU:OE2	3:2:269:ARG:HD2	1.94	0.66
1:R:132:SER:HB3	2:1:166:ILE:CD1	2.23	0.66
1:R:82:GLN:HE21	2:1:228:LEU:HD12	1.06	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:R:128:PHE:HD2	2:1:108:TRP:CD1	2.10	0.66
1:R:59:THR:HG23	1:R:127:THR:HG23	1.78	0.66
1:R:240:VAL:O	1:R:241:TYR:HD2	1.79	0.66
2:1:102:SER:O	2:1:104:LEU:N	2.29	0.65
1:R:157:THR:OG1	1:R:158:GLY:HA2	1.97	0.65
1:R:40:PHE:CZ	1:R:144:LYS:HD3	2.32	0.65
1:R:141:VAL:HG11	1:R:142:LEU:CD1	2.26	0.65
1:R:141:VAL:CG1	1:R:142:LEU:CD1	2.75	0.65
2:1:158:MET:SD	2:1:177:THR:HG23	2.36	0.65
1:R:84:PRO:O	2:1:214:LYS:NZ	2.30	0.65
1:R:40:PHE:HE1	1:R:227:SER:OG	1.80	0.65
1:R:128:PHE:HD2	2:1:108:TRP:HD1	1.43	0.64
1:R:99:LEU:CB	2:1:226:ASP:OD2	2.45	0.64
1:R:159:GLU:N	1:R:160:PRO:HD3	2.13	0.64
1:R:215:ASP:O	1:R:238:LEU:CB	2.45	0.64
1:R:40:PHE:HE1	1:R:227:SER:CB	2.10	0.64
1:R:73:GLY:CA	5:3:97:HIS:CD2	2.76	0.64
1:R:133:ARG:HB2	2:1:167:PRO:O	1.98	0.64
1:R:83:GLY:CA	2:1:227:SER:HA	2.28	0.63
1:R:149:ALA:HB2	1:R:234:LEU:HD12	1.81	0.63
1:R:142:LEU:O	1:R:173:PRO:HG3	1.99	0.62
1:R:81:THR:N	2:1:234:LEU:HD13	2.11	0.62
1:R:141:VAL:CG1	1:R:142:LEU:N	2.59	0.62
1:R:85:SER:HA	2:1:214:LYS:HZ3	1.64	0.62
1:R:132:SER:HB2	2:1:166:ILE:HD13	1.79	0.62
1:R:100:GLY:HA2	3:2:138:THR:CG2	2.30	0.62
1:R:130:GLN:NE2	2:1:107:VAL:HG23	2.14	0.62
1:R:66:TRP:HB2	1:R:78:PHE:HB3	1.81	0.61
1:R:82:GLN:NE2	2:1:228:LEU:CD1	0.78	0.61
1:R:99:LEU:HD23	3:2:141:PHE:CE1	2.35	0.61
1:R:139:LEU:HD22	1:R:140:ARG:HG3	1.80	0.61
1:R:162:PRO:HD3	1:R:208:LEU:H	1.66	0.60
1:R:215:ASP:HB3	1:R:216:GLY:CA	2.30	0.60
1:R:151:VAL:HA	1:R:163:MET:O	2.01	0.60
1:R:162:PRO:HD3	1:R:209:VAL:O	2.02	0.59
1:R:128:PHE:CZ	2:1:111:THR:HG22	2.35	0.59
1:R:141:VAL:CG2	1:R:142:LEU:HD12	2.32	0.59
5:3:64:THR:O	5:3:67:MET:HG2	2.03	0.59
1:R:116:GLU:HA	5:3:59:ASN:CG	2.22	0.59
1:R:142:LEU:O	1:R:173:PRO:HD2	2.01	0.59
1:R:181:SER:HB3	1:R:219:VAL:HG23	1.85	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:R:82:GLN:NE2	2:1:228:LEU:HD11	0.65	0.59
1:R:41:LEU:HB3	1:R:143:ALA:CA	2.29	0.58
1:R:131:GLY:H	2:1:107:VAL:HB	1.67	0.58
2:1:27:LEU:HD21	2:1:72:ARG:HG3	1.85	0.58
1:R:68:ARG:HG2	1:R:70:GLY:HA3	1.85	0.58
3:2:110:VAL:HG22	3:2:250:LEU:HD12	1.86	0.58
1:R:215:ASP:HB3	1:R:216:GLY:HA2	1.85	0.58
5:3:198:VAL:HG11	5:3:208:MET:CE	2.34	0.58
1:R:214:VAL:HG22	1:R:215:ASP:H	1.67	0.58
1:R:75:MET:O	5:3:93:PRO:HD3	2.04	0.58
1:R:116:GLU:CB	5:3:59:ASN:HD21	2.16	0.57
1:R:41:LEU:CA	1:R:143:ALA:CB	2.81	0.57
1:R:151:VAL:HG21	1:R:238:LEU:HA	1.86	0.57
4:4:18:ARG:HG3	4:4:24:THR:HG21	1.87	0.57
1:R:188:GLN:HB3	1:R:208:LEU:O	2.03	0.57
3:2:60:ALA:O	3:2:254:PRO:HG2	2.04	0.57
3:2:66:LEU:HD12	3:2:250:LEU:HD23	1.86	0.57
3:2:100:TYR:CE1	5:3:137:PRO:HG2	2.40	0.57
1:R:172:ARG:HA	1:R:201:THR:H	1.70	0.57
3:2:73:LYS:HE2	3:2:242:SER:HA	1.87	0.56
1:R:65:THR:HB	1:R:124:LEU:HG	1.86	0.56
1:R:128:PHE:CE1	2:1:111:THR:HG21	2.40	0.56
2:1:24:ALA:HB2	2:1:76:THR:CG2	2.33	0.56
1:R:131:GLY:HA2	1:R:133:ARG:HD3	1.86	0.56
1:R:141:VAL:CG1	1:R:142:LEU:HD12	2.35	0.56
1:R:161:VAL:HB	1:R:163:MET:CB	2.14	0.56
1:R:82:GLN:HB3	2:1:228:LEU:CA	2.31	0.56
1:R:141:VAL:CG1	1:R:142:LEU:HB2	2.34	0.56
3:2:263:ARG:HD3	5:3:136:PRO:O	2.06	0.56
2:1:249:ASN:OD1	2:1:250:PRO:HD2	2.07	0.55
1:R:40:PHE:CZ	1:R:144:LYS:HB2	2.41	0.55
1:R:85:SER:HA	2:1:214:LYS:NZ	2.21	0.55
1:R:130:GLN:HG2	2:1:107:VAL:HG23	1.87	0.55
1:R:42:GLY:N	1:R:43:ASP:HA	2.22	0.55
1:R:81:THR:HG21	2:1:234:LEU:CA	2.33	0.55
3:2:270:THR:O	3:2:271:GLN:HG2	2.07	0.54
2:1:132:MET:HG2	2:1:261:MET:CE	2.37	0.54
1:R:100:GLY:HA3	3:2:138:THR:HG21	1.89	0.54
1:R:120:SER:HB3	1:R:138:TRP:CG	2.43	0.54
1:R:41:LEU:HD22	1:R:42:GLY:H	1.72	0.54
1:R:147:ASN:ND2	1:R:223:VAL:HG21	2.23	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:R:99:LEU:HD23	3:2:171:ARG:CZ	2.36	0.54
1:R:99:LEU:HB3	2:1:226:ASP:OD2	2.07	0.54
1:R:152:GLN:O	1:R:163:MET:HB3	2.08	0.54
1:R:73:GLY:HA3	5:3:97:HIS:CE1	2.43	0.54
2:1:132:MET:HG2	2:1:261:MET:HE2	1.90	0.54
2:1:40:GLU:HB3	4:4:64:THR:HB	1.91	0.53
1:R:80:GLN:CG	1:R:98:ARG:HG2	2.38	0.53
1:R:36:GLN:NE2	1:R:137:ILE:HG13	2.23	0.53
1:R:40:PHE:CE1	1:R:227:SER:HB3	2.43	0.53
1:R:41:LEU:H	1:R:143:ALA:HB2	1.73	0.53
2:1:132:MET:HE1	6:1:999:SC4:H71	1.90	0.53
1:R:132:SER:HB2	2:1:166:ILE:HG21	1.90	0.53
1:R:87:SER:OG	5:3:181:ASN:ND2	2.42	0.53
1:R:49:CYS:HB2	1:R:66:TRP:HZ2	1.73	0.53
2:1:217:LEU:O	2:1:220:GLN:HG2	2.09	0.53
1:R:127:THR:HG22	1:R:129:PRO:HD2	1.91	0.53
1:R:130:GLN:CD	2:1:105:PHE:CE1	2.71	0.52
2:1:193:ARG:NH1	5:3:8:PRO:HG2	2.23	0.52
1:R:130:GLN:HA	2:1:107:VAL:CA	2.38	0.52
2:1:176:GLN:HE22	5:3:233:GLN:HE22	1.54	0.52
1:R:80:GLN:O	1:R:98:ARG:NH2	2.42	0.52
1:R:73:GLY:N	1:R:74:SER:HA	2.25	0.52
3:2:172:ASN:ND2	3:2:173:PHE:H	2.07	0.52
1:R:98:ARG:C	1:R:100:GLY:H	2.13	0.52
2:1:93:ASP:C	2:1:95:ASP:H	2.14	0.51
2:1:209:TYR:O	2:1:230:GLY:HA2	2.11	0.51
3:2:10:SER:HB2	4:4:69:ASN:O	2.10	0.51
1:R:43:ASP:CB	1:R:44:SER:CA	2.66	0.51
3:2:241:GLU:HB3	3:2:243:SER:O	2.11	0.51
1:R:40:PHE:CE1	1:R:144:LYS:HB2	2.46	0.51
1:R:41:LEU:C	1:R:143:ALA:CB	2.79	0.51
1:R:71:GLU:OE1	5:3:97:HIS:CG	2.63	0.51
2:1:273:PRO:HB3	3:2:188:ASN:HB3	1.93	0.50
5:3:198:VAL:HG11	5:3:208:MET:HE1	1.93	0.50
1:R:169:THR:HA	1:R:202:VAL:HG23	1.93	0.50
2:1:132:MET:CG	2:1:261:MET:HE2	2.42	0.50
1:R:158:GLY:HA3	1:R:159:GLU:HB3	1.94	0.49
1:R:187:PRO:O	1:R:188:GLN:HG3	2.11	0.49
8:1:348:HOH:O	5:3:15:THR:HG23	2.11	0.49
3:2:63:PHE:CD1	3:2:253:ALA:HB2	2.47	0.49
1:R:98:ARG:CD	1:R:104:ARG:NH2	2.64	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:R:93:GLU:HB3	1:R:109:ARG:HB3	1.92	0.49
1:R:155:GLN:NE2	1:R:212:SER:H	2.10	0.49
5:3:149:MET:HG2	5:3:149:MET:O	2.13	0.49
1:R:81:THR:HG23	2:1:234:LEU:HA	1.88	0.49
1:R:100:GLY:CA	3:2:138:THR:CG2	2.87	0.49
2:1:182:SER:H	5:3:15:THR:CG2	2.25	0.49
1:R:102:GLU:HG3	1:R:104:ARG:NH1	2.28	0.49
1:R:71:GLU:OE1	5:3:97:HIS:CD2	2.66	0.48
1:R:114:ARG:CD	5:3:59:ASN:HD22	2.26	0.48
1:R:180:HIS:HA	1:R:186:MET:HG3	1.96	0.48
2:1:158:MET:SD	2:1:177:THR:CG2	3.02	0.48
3:2:241:GLU:CD	3:2:241:GLU:N	2.66	0.48
2:1:49:THR:O	4:4:57:ILE:HD11	2.13	0.48
3:2:146:ASN:HB3	3:2:172:ASN:ND2	2.29	0.48
1:R:60:HIS:O	1:R:128:PHE:N	2.26	0.48
2:1:221:ALA:O	2:1:224:GLU:HB2	2.13	0.48
1:R:145:PRO:HA	1:R:170:GLY:O	2.12	0.48
1:R:240:VAL:O	1:R:241:TYR:CD2	2.64	0.48
1:R:99:LEU:CD2	3:2:141:PHE:CE1	2.97	0.48
1:R:40:PHE:CE1	1:R:144:LYS:CB	2.96	0.47
2:1:261:MET:HE1	2:1:263:PRO:HG3	1.97	0.47
1:R:41:LEU:C	1:R:143:ALA:HB3	2.33	0.47
3:2:11:VAL:HG23	4:4:69:ASN:HB3	1.94	0.47
2:1:273:PRO:HB3	3:2:188:ASN:CB	2.44	0.47
2:1:193:ARG:CZ	5:3:8:PRO:HG2	2.44	0.47
1:R:179:TRP:CZ3	1:R:219:VAL:HG22	2.49	0.47
2:1:159:TYR:CE2	2:1:161:PRO:HG3	2.49	0.47
1:R:116:GLU:CD	5:3:59:ASN:CG	2.72	0.47
1:R:224:GLU:O	1:R:225:HIS:CB	2.62	0.47
2:1:181:PRO:HA	5:3:15:THR:HG22	1.97	0.47
1:R:141:VAL:HG13	1:R:142:LEU:CD1	2.44	0.47
1:R:54:PRO:CA	1:R:55:ASN:CB	2.87	0.47
2:1:257:ILE:HD12	2:1:257:ILE:N	2.29	0.47
1:R:104:ARG:HA	1:R:104:ARG:HD3	1.57	0.47
1:R:170:GLY:HA2	1:R:200:VAL:HG13	1.96	0.47
1:R:98:ARG:O	1:R:99:LEU:HD22	2.15	0.47
1:R:152:GLN:HB2	1:R:153:LYS:HA	1.97	0.47
1:R:73:GLY:CA	5:3:97:HIS:CE1	2.98	0.47
2:1:48:GLU:HA	3:2:196:ILE:HB	1.95	0.46
1:R:141:VAL:CG1	1:R:142:LEU:HD13	2.40	0.46
1:R:186:MET:O	1:R:188:GLN:N	2.48	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:2:71:TRP:CE3	3:2:246:ILE:HD11	2.50	0.46
1:R:115:VAL:HG13	1:R:141:VAL:HG21	1.98	0.46
1:R:188:GLN:O	1:R:207:ILE:HG22	2.16	0.46
1:R:40:PHE:HD1	1:R:227:SER:HG	1.48	0.46
2:1:94:ASN:OD1	2:1:248:HIS:HB3	2.16	0.46
2:1:176:GLN:HE22	5:3:233:GLN:HE21	1.58	0.46
1:R:38:PRO:HA	1:R:39:GLY:HA3	1.66	0.45
1:R:161:VAL:HG12	1:R:208:LEU:HB2	1.97	0.45
2:1:110:ILE:HG13	2:1:239:SER:HA	1.99	0.45
5:3:87:LEU:HD11	5:3:114:LEU:HD12	1.98	0.45
2:1:176:GLN:NE2	5:3:233:GLN:NE2	2.55	0.45
1:R:210:PRO:HB2	1:R:211:SER:H	1.68	0.45
4:4:57:ILE:HA	4:4:57:ILE:HD13	1.78	0.44
1:R:157:THR:OG1	1:R:158:GLY:CA	2.64	0.44
1:R:81:THR:HG21	2:1:234:LEU:HD23	0.45	0.44
1:R:172:ARG:O	1:R:174:PRO:HD3	2.17	0.44
1:R:207:ILE:HA	1:R:208:LEU:HB3	2.00	0.44
1:R:162:PRO:HB2	1:R:238:LEU:HG	2.00	0.44
1:R:31:VAL:HG12	1:R:51:LEU:HD11	1.98	0.44
3:2:238:PHE:O	3:2:240:THR:N	2.51	0.44
1:R:73:GLY:HA2	5:3:230:HIS:NE2	2.32	0.44
1:R:41:LEU:CB	1:R:143:ALA:CB	2.48	0.44
1:R:177:ILE:HG22	1:R:223:VAL:HG22	2.00	0.44
1:R:114:ARG:HD3	5:3:59:ASN:HD22	1.83	0.44
1:R:130:GLN:HE22	1:R:133:ARG:HD2	1.82	0.44
1:R:141:VAL:HG21	1:R:142:LEU:HD12	2.00	0.44
1:R:241:TYR:HA	1:R:242:TYR:C	2.37	0.44
4:4:33:TYR:HB2	4:4:38:SER:HB2	2.00	0.44
1:R:65:THR:HA	1:R:78:PHE:O	2.18	0.43
2:1:134:PHE:O	2:1:193:ARG:HA	2.19	0.43
1:R:130:GLN:CG	2:1:105:PHE:O	2.64	0.43
1:R:162:PRO:CD	1:R:208:LEU:H	2.32	0.43
3:2:110:VAL:HG22	3:2:250:LEU:CD1	2.47	0.43
3:2:263:ARG:NH1	5:3:137:PRO:O	2.52	0.43
1:R:82:GLN:O	1:R:98:ARG:HG3	2.18	0.43
3:2:67:ASP:N	8:2:336:HOH:O	2.51	0.43
1:R:116:GLU:OE2	5:3:59:ASN:OD1	2.37	0.42
4:4:7:SER:HA	4:4:26:ASN:HA	2.01	0.42
1:R:40:PHE:HZ	1:R:144:LYS:CG	2.31	0.42
2:1:42:PRO:HA	4:4:63:LYS:O	2.18	0.42
1:R:84:PRO:C	2:1:214:LYS:NZ	2.71	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:R:69:HIS:N	1:R:70:GLY:HA3	2.33	0.42
1:R:132:SER:HB2	2:1:166:ILE:CG2	2.49	0.42
1:R:31:VAL:HG11	1:R:134:SER:HA	2.01	0.42
3:2:263:ARG:HB2	3:2:264:ASN:H	1.68	0.42
1:R:144:LYS:HG3	1:R:145:PRO:N	2.34	0.42
2:1:92:VAL:HG23	2:1:255:SER:HB2	2.02	0.42
4:4:49:ASP:HA	4:4:50:PRO:HD3	1.85	0.42
1:R:128:PHE:HB3	1:R:129:PRO:CD	2.43	0.42
1:R:81:THR:HA	2:1:234:LEU:HD22	1.63	0.42
2:1:148:GLY:HA3	2:1:251:THR:CG2	2.43	0.42
1:R:128:PHE:HE2	2:1:108:TRP:HE1	0.48	0.42
1:R:142:LEU:HD23	1:R:172:ARG:CB	2.50	0.42
5:3:195:GLN:HA	5:3:195:GLN:OE1	2.20	0.41
1:R:71:GLU:OE1	5:3:97:HIS:CE1	2.73	0.41
1:R:144:LYS:CG	1:R:145:PRO:N	2.83	0.41
1:R:168:SER:HB3	1:R:203:THR:HG23	2.01	0.41
1:R:98:ARG:CG	1:R:99:LEU:H	2.24	0.41
1:R:85:SER:CA	2:1:214:LYS:NZ	2.84	0.41
8:1:348:HOH:O	5:3:15:THR:CG2	2.67	0.41
2:1:24:ALA:HB3	4:4:45:ASP:O	2.21	0.41
3:2:141:PHE:CE2	3:2:171:ARG:HG2	2.56	0.41
1:R:157:THR:CB	1:R:158:GLY:HA2	2.50	0.41
1:R:159:GLU:H	1:R:160:PRO:HD3	1.82	0.41
5:3:124:MET:HG3	5:3:124:MET:O	2.18	0.41
5:3:44:MET:O	5:3:48:GLU:HG3	2.20	0.41
4:4:3:ALA:HA	4:4:30:ILE:HG12	2.01	0.41
1:R:161:VAL:HA	1:R:162:PRO:HD3	1.94	0.41
1:R:76:ALA:H	1:R:77:VAL:HG23	1.85	0.41
1:R:41:LEU:O	1:R:143:ALA:CB	2.69	0.41
2:1:162:PRO:HD3	2:1:238:GLY:CA	2.50	0.41
1:R:126:VAL:HA	1:R:132:SER:HA	2.03	0.41
1:R:221:CYS:HB3	1:R:234:LEU:HG	2.02	0.41
1:R:40:PHE:CZ	1:R:144:LYS:CD	2.92	0.41
1:R:186:MET:H	1:R:187:PRO:CD	2.34	0.41
5:3:143:LYS:HA	5:3:143:LYS:HD3	1.75	0.41
2:1:292:PRO:HG2	2:1:293:LEU:HD12	2.02	0.40
2:1:94:ASN:O	2:1:95:ASP:HB2	2.21	0.40
1:R:116:GLU:CG	5:3:59:ASN:ND2	2.62	0.40
1:R:139:LEU:HB3	1:R:140:ARG:H	1.57	0.40
1:R:157:THR:HG21	1:R:211:SER:OG	2.22	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 PDB entries followed by that with respect to all EM entries.

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	R	211/213 (99%)	137 (65%)	50 (24%)	24 (11%)	0 10
2	1	268/278 (96%)	250 (93%)	16 (6%)	2 (1%)	26 71
3	2	260/262 (99%)	245 (94%)	13 (5%)	2 (1%)	24 69
4	4	66/68 (97%)	62 (94%)	4 (6%)	0	100 100
5	3	233/235 (99%)	226 (97%)	7 (3%)	0	100 100
All	All	1038/1056 (98%)	920 (89%)	90 (9%)	28 (3%)	10 45

All (28) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	R	55	ASN
1	R	99	LEU
1	R	141	VAL
1	R	188	GLN
1	R	210	PRO
1	R	212	SER
2	1	103	LYS
3	2	239	ALA
1	R	139	LEU
1	R	198	GLY
1	R	56	MET
1	R	88	GLU
1	R	131	GLY
1	R	134	SER
1	R	143	ALA
1	R	160	PRO
1	R	161	VAL
1	R	187	PRO
1	R	215	ASP
3	2	48	ASN
1	R	145	PRO

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Mol	Chain	Res	Type
1	R	225	HIS
1	R	232	GLN
1	R	162	PRO
1	R	209	VAL
1	R	214	VAL
1	R	129	PRO
2	1	270	CYS

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 PDB entries followed by that with respect to all EM entries.

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	R	185/185 (100%)	149 (80%)	36 (20%)	2 12
2	1	231/235 (98%)	227 (98%)	4 (2%)	68 87
3	2	221/221 (100%)	208 (94%)	13 (6%)	24 61
4	4	56/56 (100%)	53 (95%)	3 (5%)	27 64
5	3	207/207 (100%)	196 (95%)	11 (5%)	28 64
All	All	900/904 (100%)	833 (93%)	67 (7%)	22 54

All (67) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	R	43	ASP
1	R	51	LEU
1	R	58	VAL
1	R	60	HIS
1	R	61	VAL
1	R	68	ARG
1	R	75	MET
1	R	98	ARG
1	R	99	LEU
1	R	102	GLU
1	R	103	LEU
1	R	104	ARG

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Mol	Chain	Res	Type
1	R	111	PHE
1	R	120	SER
1	R	124	LEU
1	R	126	VAL
1	R	128	PHE
1	R	133	ARG
1	R	139	LEU
1	R	142	LEU
1	R	144	LYS
1	R	152	GLN
1	R	157	THR
1	R	163	MET
1	R	176	GLN
1	R	196	LEU
1	R	199	THR
1	R	200	VAL
1	R	208	LEU
1	R	212	SER
1	R	213	GLN
1	R	214	VAL
1	R	230	LYS
1	R	234	LEU
1	R	236	VAL
1	R	241	TYR
2	1	83	ARG
2	1	144	ASP
2	1	149	HIS
2	1	182	SER
3	2	31	SER
3	2	52	GLN
3	2	70	THR
3	2	151	GLU
3	2	161	THR
3	2	172	ASN
3	2	200	ARG
3	2	203	ASN
3	2	241	GLU
3	2	243	SER
3	2	244	THR
3	2	245	GLU
3	2	263	ARG
4	4	18	ARG

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Mol	Chain	Res	Type
4	4	24	THR
4	4	69	ASN
5	3	5	LEU
5	3	15	THR
5	3	51	THR
5	3	99	MET
5	3	124	MET
5	3	149	MET
5	3	160	LEU
5	3	208	MET
5	3	218	ASN
5	3	224	LEU
5	3	227	ASP

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

Mol	Chain	Res	Type
1	R	36	GLN
1	R	55	ASN
1	R	69	HIS
1	R	80	GLN
1	R	130	GLN
1	R	152	GLN
1	R	155	GLN
1	R	232	GLN
2	1	62	GLN
2	1	65	HIS
2	1	149	HIS
2	1	220	GLN
3	2	48	ASN
3	2	172	ASN
3	2	203	ASN
4	4	69	ASN
5	3	59	ASN
5	3	181	ASN
5	3	218	ASN
5	3	233	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 carbohydrates in this entry.

5.6 Ligand geometry (i)

2 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
6	SC4	1	999	-	29,29,29	1.31	5 (17%)	39,39,39	1.34	5 (12%)
7	MYR	4	1	4	10,10,15	0.67	0	9,9,15	0.77	0

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
6	SC4	1	999	-	-	0/12/12/12	0/3/3/3
7	MYR	4	1	4	-	0/7/8/13	0/0/0/0

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	1	999	SC4	C3-C4	2.06	1.43	1.39
6	1	999	SC4	C18-C17	2.07	1.42	1.38
6	1	999	SC4	C18-C19	2.29	1.43	1.38
6	1	999	SC4	C15-C16	2.41	1.45	1.40
6	1	999	SC4	C5-C6	3.28	1.44	1.38

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
6	1	999	SC4	O1-C3-C2	-2.95	117.99	124.03
6	1	999	SC4	C15-C20-CL3	2.46	121.42	118.44
6	1	999	SC4	C15-C16-CL2	2.59	121.58	118.44
6	1	999	SC4	C21-O3-C6	2.61	123.56	117.51
6	1	999	SC4	O1-C3-C4	4.06	121.89	116.40

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	1	999	SC4	1	0

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.