Full wwPDB X-ray Structure Validation Report ⓘ

May 12, 2020 – 11:48 pm BST

PDB ID	:	1MHE
Title	:	THE HUMAN NON-CLASSICAL MAJOR HISTOCOMPATIBILITY COMPLEX MOLECULE HLA-E
Authors	:	O'Callaghan, C.A.; Tormo, J.; Willcox, B.E.; Braud, V.B.; Jakobsen, B.K.; Stuart, D.I.; Mcmichael, A.J.; Bell, J.I.; Jones, E.Y.
Deposited on	:	1998-08-24
Resolution	:	2.85 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 following versions of software and data (see references ⓘ) were used in the production of this report:

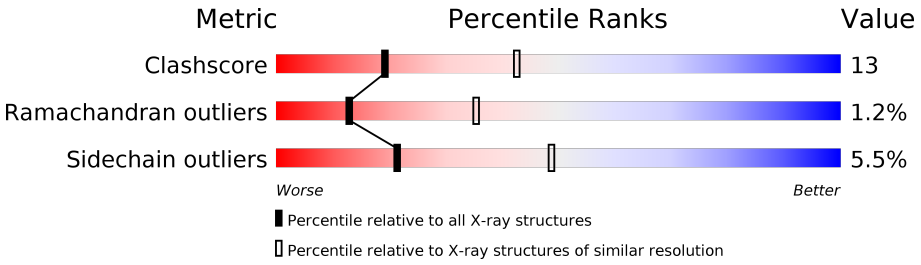
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

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 2.85 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	141614	3438 (2.90-2.82)
Ramachandran outliers	138981	3348 (2.90-2.82)
Sidechain outliers	138945	3351 (2.90-2.82)

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 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 <=5%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	274	68% 25% 5% •
1	C	274	71% 24% •
2	B	100	80% 18% •
2	D	100	76% 23% •
3	P	9	44% 56%
3	Q	9	78% 22%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 7752 atoms, of which 1459 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 I HISTOCOMPATIBILITY ANTIGEN HLA-E.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	273	Total	C	H	N	O	S	44	0	0
			2748	1396	514	401	430	7			
1	C	273	Total	C	H	N	O	S	0	0	0
			2748	1396	514	401	430	7			

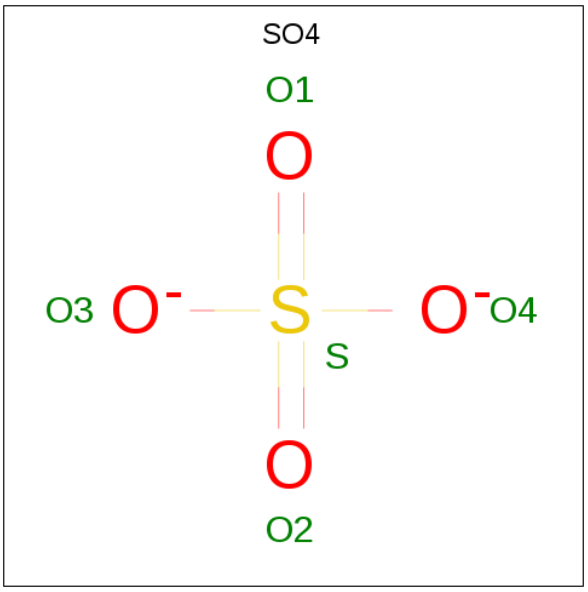
- Molecule 2 is a protein called BETA-2-MICROGLOBULIN.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	100	Total	C	H	N	O	S	0	0	0
			1029	533	192	141	159	4			
2	D	100	Total	C	H	N	O	S	0	0	0
			1029	533	192	141	159	4			

- Molecule 3 is a protein called PEPTIDE (VMAPRTVLL).

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
3	P	9	Total	C	H	N	O	S	0	0	0
			85	45	16	12	11	1			
3	Q	9	Total	C	H	N	O	S	0	0	0
			85	45	16	12	11	1			

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total O S 5 4 1	0	0

- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	3	Total H O 9 6 3	0	0
5	C	5	Total H O 14 9 5	0	0

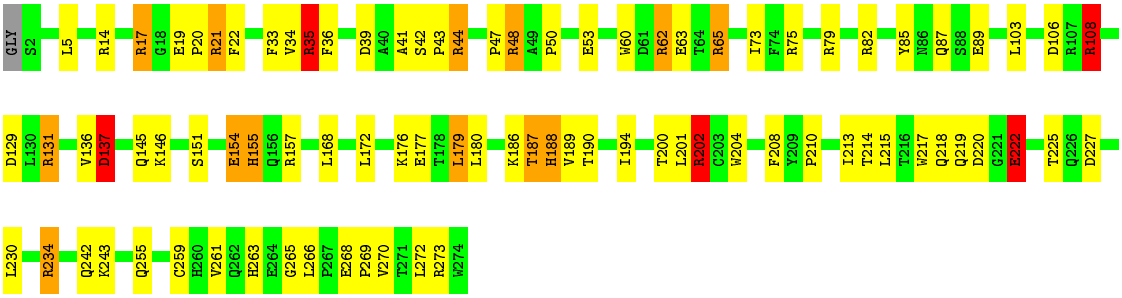
3 Residue-property plots (i)

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

Note EDS was not executed.

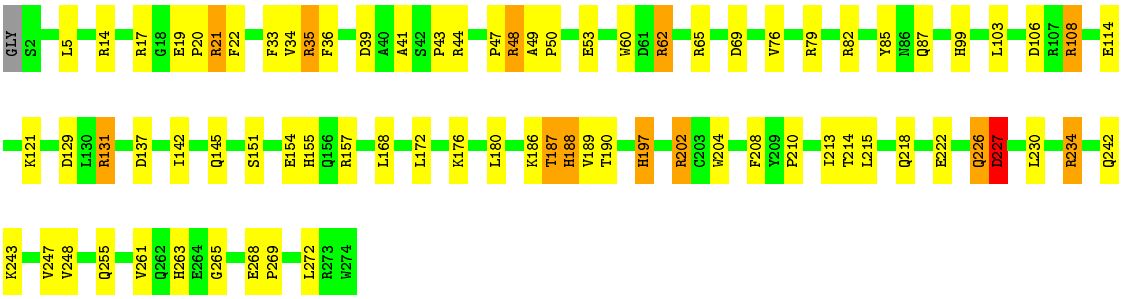
- Molecule 1: HLA CLASS I HISTOCOMPATIBILITY ANTIGEN HLA-E

Chain A:



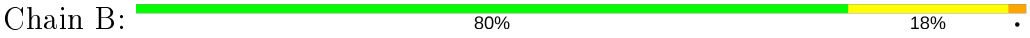
- Molecule 1: HLA CLASS I HISTOCOMPATIBILITY ANTIGEN HLA-E

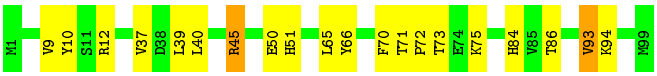
Chain C:



- Molecule 2: BETA-2-MICROGLOBULIN

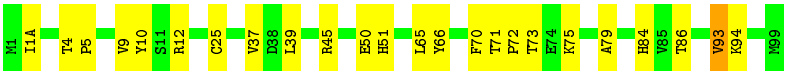
Chain B:





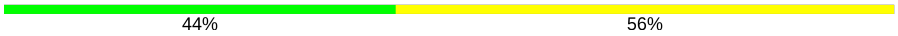
- Molecule 2: BETA-2-MICROGLOBULIN

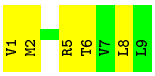
Chain D:



- Molecule 3: PEPTIDE (VMAPRTVLL)

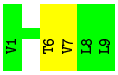
Chain P:





- Molecule 3: PEPTIDE (VMAPRTVLL)

Chain Q:



4 Data and refinement statistics (i)

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

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α, β, γ	182.20Å 182.20Å 88.40Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	25.00 – 2.85	Depositor
% Data completeness (in resolution range)	91.0 (25.00-2.85)	Depositor
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.202 , 0.235	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	7752	wwPDB-VP
Average B, all atoms (Å^2)	51.0	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: SO4

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	0.52	0/2300	1.11	20/3127 (0.6%)
1	C	0.55	0/2300	1.08	20/3127 (0.6%)
2	B	0.56	0/860	0.84	2/1162 (0.2%)
2	D	0.53	0/860	0.88	3/1162 (0.3%)
3	P	0.55	0/69	0.83	0/92
3	Q	0.61	0/69	0.81	0/92
All	All	0.54	0/6458	1.03	45/8762 (0.5%)

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	A	0	1

There are no bond length outliers.

All (45) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	35	ARG	NE-CZ-NH2	-15.74	112.43	120.30
1	A	35	ARG	NE-CZ-NH1	13.53	127.06	120.30
1	C	35	ARG	NE-CZ-NH2	-13.14	113.73	120.30
1	C	108	ARG	NE-CZ-NH2	-13.09	113.75	120.30
1	A	108	ARG	NE-CZ-NH2	-12.27	114.17	120.30
1	C	108	ARG	NE-CZ-NH1	11.96	126.28	120.30
1	C	35	ARG	NE-CZ-NH1	11.84	126.22	120.30
1	A	131	ARG	NE-CZ-NH2	-11.77	114.42	120.30
2	D	45	ARG	NE-CZ-NH2	-11.73	114.44	120.30
2	D	45	ARG	NE-CZ-NH1	11.59	126.10	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	44	ARG	NE-CZ-NH1	11.46	126.03	120.30
1	C	131	ARG	NE-CZ-NH2	-11.40	114.60	120.30
1	A	65	ARG	NE-CZ-NH2	-11.37	114.61	120.30
1	C	48	ARG	NE-CZ-NH2	-11.34	114.63	120.30
1	A	65	ARG	NE-CZ-NH1	11.32	125.96	120.30
1	A	108	ARG	NE-CZ-NH1	11.25	125.92	120.30
1	A	48	ARG	NE-CZ-NH2	-11.23	114.69	120.30
1	C	44	ARG	NE-CZ-NH2	-11.11	114.75	120.30
2	B	45	ARG	NE-CZ-NH2	-11.00	114.80	120.30
1	C	131	ARG	NE-CZ-NH1	10.92	125.76	120.30
1	A	234	ARG	NE-CZ-NH2	-10.74	114.93	120.30
2	B	45	ARG	NE-CZ-NH1	10.66	125.63	120.30
1	C	202	ARG	NE-CZ-NH2	-10.62	114.99	120.30
1	A	44	ARG	NE-CZ-NH2	-10.60	115.00	120.30
1	A	234	ARG	NE-CZ-NH1	10.48	125.54	120.30
1	A	48	ARG	NE-CZ-NH1	10.31	125.45	120.30
1	A	44	ARG	NE-CZ-NH1	10.18	125.39	120.30
1	A	202	ARG	NE-CZ-NH2	-10.17	115.22	120.30
1	C	48	ARG	NE-CZ-NH1	10.16	125.38	120.30
1	A	131	ARG	NE-CZ-NH1	9.66	125.13	120.30
1	C	65	ARG	NE-CZ-NH2	-9.25	115.67	120.30
1	A	202	ARG	NE-CZ-NH1	9.03	124.82	120.30
1	C	65	ARG	NE-CZ-NH1	9.03	124.81	120.30
1	C	202	ARG	NE-CZ-NH1	8.88	124.74	120.30
1	C	35	ARG	CG-CD-NE	-8.57	93.81	111.80
1	C	234	ARG	NE-CZ-NH1	6.99	123.80	120.30
1	C	35	ARG	CD-NE-CZ	6.93	133.30	123.60
1	A	35	ARG	CD-NE-CZ	6.39	132.55	123.60
1	C	108	ARG	CD-NE-CZ	6.39	132.54	123.60
1	C	44	ARG	CD-NE-CZ	5.72	131.61	123.60
2	D	45	ARG	CD-NE-CZ	5.65	131.51	123.60
1	A	65	ARG	CD-NE-CZ	5.63	131.49	123.60
1	A	108	ARG	CD-NE-CZ	5.54	131.36	123.60
1	A	137	ASP	N-CA-C	-5.24	96.85	111.00
1	C	234	ARG	NE-CZ-NH2	-5.24	117.68	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	35	ARG	Sidechain

5.2 Too-close contacts (i)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2234	514	2072	71	0
1	C	2234	514	2072	61	0
2	B	837	192	803	17	0
2	D	837	192	803	17	0
3	P	69	16	83	7	0
3	Q	69	16	83	1	0
4	A	5	0	0	0	0
5	A	3	6	0	0	0
5	C	5	9	0	0	0
All	All	6293	1459	5916	156	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 13.

All (156) 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:C:226:GLN:O	1:C:227:ASP:HB2	1.77	0.83
1:A:202:ARG:HG2	1:A:204:TRP:NE1	1.98	0.79
1:A:35:ARG:HH11	1:A:48:ARG:HH11	1.32	0.76
1:A:17:ARG:HD3	1:A:17:ARG:N	2.05	0.71
1:C:218:GLN:HB2	1:C:222:GLU:O	1.90	0.71
1:C:142:ILE:HD12	1:C:142:ILE:H	1.56	0.70
1:A:35:ARG:HG2	1:A:48:ARG:HD3	1.74	0.69
1:C:213:ILE:HG12	1:C:214:THR:N	2.10	0.66
1:C:47:PRO:HG3	1:C:60:TRP:CZ2	2.30	0.66
1:A:213:ILE:HG12	1:A:214:THR:N	2.11	0.65
1:C:202:ARG:HG2	1:C:204:TRP:NE1	2.12	0.65
1:A:5:LEU:HB2	1:A:168:LEU:HD13	1.79	0.65
1:A:35:ARG:HH11	1:A:48:ARG:NH1	1.94	0.64
1:A:47:PRO:HG3	1:A:60:TRP:CZ2	2.32	0.64
1:A:234:ARG:HH11	2:B:10:TYR:CB	2.11	0.64
1:C:213:ILE:HG12	1:C:214:THR:H	1.64	0.63
1:C:268:GLU:HB2	1:C:269:PRO:HD2	1.81	0.62
1:A:186:LYS:N	1:A:186:LYS:HD2	2.15	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:17:ARG:HH11	1:A:17:ARG:H	1.46	0.61
1:C:255:GLN:H	1:C:255:GLN:CD	2.03	0.61
1:C:5:LEU:HB2	1:C:168:LEU:HD13	1.81	0.61
1:A:172:LEU:HD23	1:A:179:LEU:HD23	1.83	0.60
1:A:41:ALA:O	1:A:43:PRO:HD3	2.02	0.60
1:A:234:ARG:NH1	2:B:10:TYR:CB	2.65	0.59
1:C:176:LYS:HG3	1:C:180:LEU:HD12	1.84	0.59
1:A:155:HIS:NE2	3:P:5:ARG:HD3	2.17	0.59
1:A:35:ARG:NH1	1:A:48:ARG:NH1	2.50	0.59
1:A:62:ARG:HH11	1:A:62:ARG:CB	2.15	0.59
1:A:146:LYS:HE3	1:C:69:ASP:OD2	2.02	0.59
1:C:47:PRO:HG3	1:C:60:TRP:CH2	2.38	0.58
1:C:186:LYS:HD2	1:C:186:LYS:N	2.17	0.58
1:A:213:ILE:HG12	1:A:214:THR:H	1.69	0.58
1:C:14:ARG:HD2	1:C:17:ARG:HD3	1.86	0.57
1:C:62:ARG:CB	1:C:62:ARG:HH11	2.18	0.56
1:C:76:VAL:O	1:C:79:ARG:HG2	2.04	0.56
1:A:35:ARG:NH1	1:A:48:ARG:HH11	2.04	0.56
1:C:263:HIS:CD2	1:C:265:GLY:H	2.24	0.55
2:D:51:HIS:HA	2:D:65:LEU:O	2.07	0.55
1:C:79:ARG:HA	1:C:82:ARG:NH1	2.22	0.55
1:C:47:PRO:O	1:C:48:ARG:HD3	2.07	0.54
1:A:234:ARG:NH1	2:B:10:TYR:HB3	2.21	0.54
1:A:129:ASP:OD2	1:A:131:ARG:HB2	2.09	0.53
1:C:142:ILE:HD12	1:C:142:ILE:N	2.23	0.53
1:A:106:ASP:OD2	1:A:108:ARG:HB2	2.08	0.53
1:A:266:LEU:HD13	1:A:270:VAL:HG23	1.90	0.53
2:B:51:HIS:HA	2:B:65:LEU:O	2.09	0.53
1:C:197:HIS:N	1:C:197:HIS:ND1	2.57	0.53
1:C:215:LEU:HD22	1:C:261:VAL:HG22	1.90	0.53
2:D:84:HIS:CE1	2:D:86:THR:HG23	2.44	0.53
1:A:89:GLU:O	1:A:89:GLU:HG2	2.09	0.53
1:A:79:ARG:HA	1:A:82:ARG:NH1	2.24	0.52
1:C:41:ALA:O	1:C:43:PRO:HD3	2.08	0.52
1:A:234:ARG:HD2	2:B:10:TYR:CZ	2.44	0.52
1:A:210:PRO:O	1:A:263:HIS:HE1	1.92	0.52
1:C:131:ARG:HD3	1:C:157:ARG:NH1	2.23	0.52
1:A:215:LEU:HD22	1:A:261:VAL:HG22	1.91	0.52
1:A:47:PRO:HG3	1:A:60:TRP:CH2	2.44	0.52
2:B:84:HIS:CE1	2:B:86:THR:HG23	2.44	0.52
1:A:230:LEU:HD11	1:A:243:LYS:HE3	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:33:PHE:CD2	1:C:34:VAL:HG13	2.45	0.51
1:C:230:LEU:HD11	1:C:243:LYS:HE3	1.93	0.51
1:A:208:PHE:CE1	1:A:242:GLN:HA	2.46	0.51
1:A:187:THR:HG21	1:A:272:LEU:HD21	1.93	0.51
1:A:65:ARG:NH2	1:C:145:GLN:NE2	2.59	0.50
1:C:234:ARG:HH21	2:D:10:TYR:CB	2.24	0.50
1:C:103:LEU:HG	1:C:168:LEU:HD23	1.92	0.50
1:A:136:VAL:O	1:A:137:ASP:HB2	2.10	0.50
1:A:272:LEU:HD22	1:A:272:LEU:N	2.27	0.50
1:A:33:PHE:CD2	1:A:34:VAL:HG13	2.47	0.50
1:C:121:LYS:HE2	2:D:1(A):ILE:HD11	1.92	0.50
1:A:75:ARG:HH12	1:C:79:ARG:NH2	2.10	0.49
2:B:37:VAL:HG21	2:B:66:TYR:CE1	2.47	0.49
1:C:62:ARG:HB3	1:C:62:ARG:HH11	1.77	0.49
1:C:210:PRO:O	1:C:263:HIS:HE1	1.95	0.49
1:A:172:LEU:HA	1:A:179:LEU:HD23	1.95	0.49
1:A:136:VAL:HG12	1:A:136:VAL:O	2.12	0.49
1:A:62:ARG:HB3	1:A:62:ARG:HH11	1.77	0.48
1:A:103:LEU:HG	1:A:168:LEU:HD23	1.94	0.48
1:A:187:THR:O	1:A:188:HIS:HB3	2.12	0.48
1:A:263:HIS:CD2	1:A:265:GLY:H	2.30	0.48
1:C:202:ARG:HG2	1:C:204:TRP:CD1	2.49	0.48
1:C:187:THR:O	1:C:188:HIS:HB3	2.14	0.48
1:C:99:HIS:HB3	1:C:114:GLU:HG3	1.96	0.48
1:A:194:ILE:HD11	1:A:200:THR:OG1	2.13	0.48
1:C:208:PHE:CE1	1:C:242:GLN:HA	2.48	0.47
1:A:234:ARG:HD2	2:B:10:TYR:CE2	2.50	0.47
1:A:202:ARG:HG2	1:A:204:TRP:HE1	1.74	0.47
2:D:51:HIS:HB3	2:D:66:TYR:CD2	2.50	0.47
1:A:268:GLU:HB2	1:A:269:PRO:HD2	1.96	0.47
1:A:155:HIS:NE2	3:P:5:ARG:CD	2.77	0.46
1:A:35:ARG:O	1:A:35:ARG:HG3	2.14	0.46
1:A:186:LYS:N	1:A:186:LYS:CD	2.78	0.46
1:A:201:LEU:HA	1:A:201:LEU:HD23	1.74	0.46
1:C:129:ASP:OD2	1:C:131:ARG:HB2	2.15	0.46
1:A:208:PHE:HE1	1:A:242:GLN:HA	1.81	0.46
1:A:47:PRO:O	1:A:48:ARG:HD2	2.15	0.46
2:B:73:THR:CG2	2:B:75:LYS:HG2	2.46	0.46
1:A:177:GLU:CD	1:A:177:GLU:H	2.19	0.46
2:D:73:THR:CG2	2:D:75:LYS:HG2	2.45	0.46
2:D:37:VAL:HG21	2:D:66:TYR:CE1	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:63:GLU:CD	3:P:1:VAL:HB	2.36	0.45
2:D:39:LEU:HD23	2:D:39:LEU:HA	1.73	0.45
2:D:94:LYS:HB3	2:D:94:LYS:HE2	1.77	0.45
1:A:73:ILE:HG21	3:P:6:THR:HG23	1.98	0.45
3:P:2:MET:HB2	3:P:2:MET:HE3	1.80	0.45
1:A:50:PRO:HA	1:A:53:GLU:OE2	2.17	0.45
2:B:39:LEU:HD23	2:B:39:LEU:HA	1.73	0.45
1:C:234:ARG:HH21	2:D:10:TYR:HB2	1.82	0.44
2:B:73:THR:HG22	2:B:75:LYS:HG2	2.00	0.44
2:D:73:THR:HG22	2:D:75:LYS:HG2	1.98	0.44
1:A:217:TRP:CZ3	1:A:259:CYS:HB2	2.52	0.44
2:D:9:VAL:CG2	2:D:93:VAL:HG22	2.47	0.44
1:C:50:PRO:HA	1:C:53:GLU:OE2	2.18	0.44
1:C:227:ASP:CG	1:C:248:VAL:HB	2.38	0.44
1:C:142:ILE:CD1	1:C:142:ILE:H	2.29	0.43
1:C:227:ASP:O	1:C:247:VAL:HA	2.19	0.43
1:A:218:GLN:HB3	1:A:222:GLU:O	2.19	0.43
1:C:208:PHE:HE1	1:C:242:GLN:HA	1.84	0.43
1:C:272:LEU:HD22	1:C:272:LEU:N	2.33	0.43
1:C:85:TYR:HB2	1:C:87:GLN:NE2	2.34	0.43
1:A:21:ARG:HD3	1:A:22:PHE:N	2.33	0.43
2:B:94:LYS:HB3	2:B:94:LYS:HE2	1.77	0.43
2:B:9:VAL:CG2	2:B:93:VAL:HG22	2.49	0.43
1:C:226:GLN:O	1:C:227:ASP:CB	2.59	0.43
1:C:21:ARG:HG3	1:C:21:ARG:HH11	1.82	0.43
1:C:106:ASP:OD2	1:C:108:ARG:HB2	2.19	0.42
1:C:180:LEU:HD23	1:C:180:LEU:HA	1.69	0.42
1:A:145:GLN:HG3	1:A:146:LYS:N	2.35	0.42
1:A:73:ILE:HG21	3:P:6:THR:CG2	2.50	0.42
2:B:71:THR:HA	2:B:72:PRO:HD2	1.80	0.42
1:A:63:GLU:OE2	3:P:1:VAL:HB	2.20	0.42
2:D:79:ALA:HA	2:D:94:LYS:HA	2.02	0.42
1:C:234:ARG:NH2	2:D:10:TYR:CB	2.83	0.42
3:Q:6:THR:HG23	3:Q:7:VAL:N	2.35	0.42
1:A:36:PHE:C	1:A:36:PHE:CD1	2.93	0.41
1:A:85:TYR:HB2	1:A:87:GLN:NE2	2.35	0.41
2:B:51:HIS:HB3	2:B:66:TYR:CD2	2.55	0.41
1:C:36:PHE:C	1:C:36:PHE:CD1	2.93	0.41
2:D:4:THR:CG2	2:D:5:PRO:HD2	2.51	0.41
1:C:189:VAL:HG12	1:C:190:THR:N	2.36	0.41
2:B:40:LEU:HD23	2:B:45:ARG:HA	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:154:GLU:O	1:A:157:ARG:HB3	2.21	0.41
1:A:176:LYS:HA	1:A:180:LEU:HD12	2.03	0.41
1:C:255:GLN:N	1:C:255:GLN:CD	2.73	0.41
2:D:71:THR:HA	2:D:72:PRO:HD2	1.80	0.41
1:A:19:GLU:HG3	1:A:20:PRO:HD2	2.01	0.41
1:C:121:LYS:HB2	1:C:121:LYS:HE3	1.86	0.41
1:C:186:LYS:N	1:C:186:LYS:CD	2.82	0.41
1:C:172:LEU:HD23	1:C:172:LEU:HA	1.80	0.40
1:C:21:ARG:HD3	1:C:22:PHE:N	2.36	0.40
1:C:49:ALA:HA	1:C:50:PRO:HD3	1.87	0.40
2:D:25:CYS:O	2:D:65:LEU:HD12	2.21	0.40
1:A:255:GLN:O	1:A:273:ARG:NE	2.54	0.40
1:A:189:VAL:HG12	1:A:190:THR:N	2.36	0.40
2:B:40:LEU:CD2	2:B:45:ARG:HA	2.51	0.40
1:C:19:GLU:HG3	1:C:20:PRO:HD2	2.03	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	271/274 (99%)	234 (86%)	32 (12%)	5 (2%)	8 25
1	C	271/274 (99%)	236 (87%)	31 (11%)	4 (2%)	10 30
2	B	98/100 (98%)	93 (95%)	5 (5%)	0	100 100
2	D	98/100 (98%)	93 (95%)	5 (5%)	0	100 100
3	P	7/9 (78%)	7 (100%)	0	0	100 100
3	Q	7/9 (78%)	4 (57%)	3 (43%)	0	100 100
All	All	752/766 (98%)	667 (89%)	76 (10%)	9 (1%)	13 35

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	225	THR
1	A	227	ASP
1	C	227	ASP
1	A	137	ASP
1	A	188	HIS
1	C	226	GLN
1	A	222	GLU
1	C	188	HIS
1	C	137	ASP

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	236/236 (100%)	218 (92%)	18 (8%)	13 33
1	C	236/236 (100%)	226 (96%)	10 (4%)	30 60
2	B	95/95 (100%)	91 (96%)	4 (4%)	30 60
2	D	95/95 (100%)	91 (96%)	4 (4%)	30 60
3	P	8/8 (100%)	7 (88%)	1 (12%)	4 12
3	Q	8/8 (100%)	8 (100%)	0	100 100
All	All	678/678 (100%)	641 (94%)	37 (6%)	21 49

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	ARG
1	A	17	ARG
1	A	21	ARG
1	A	35	ARG
1	A	39	ASP
1	A	42	SER
1	A	44	ARG
1	A	62	ARG
1	A	108	ARG
1	A	151	SER

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Mol	Chain	Res	Type
1	A	154	GLU
1	A	155	HIS
1	A	179	LEU
1	A	187	THR
1	A	202	ARG
1	A	219	GLN
1	A	220	ASP
1	A	222	GLU
2	B	12	ARG
2	B	50	GLU
2	B	70	PHE
2	B	93	VAL
1	C	21	ARG
1	C	35	ARG
1	C	39	ASP
1	C	62	ARG
1	C	151	SER
1	C	154	GLU
1	C	155	HIS
1	C	187	THR
1	C	197	HIS
1	C	227	ASP
2	D	12	ARG
2	D	50	GLU
2	D	70	PHE
2	D	93	VAL
3	P	8	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	72	GLN
1	A	87	GLN
1	A	115	GLN
1	A	127	ASN
1	A	260	HIS
1	A	263	HIS
1	C	87	GLN
1	C	115	GLN
1	C	127	ASN
1	C	145	GLN
1	C	260	HIS

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Mol	Chain	Res	Type
1	C	263	HIS

5.3.3 RNA

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates

There are no carbohydrates in this entry.

5.6 Ligand geometry

1 ligand is modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
4	SO4	A	1002	-	4,4,4	0.84	0	6,6,6	0.58	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

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 is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.