



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 12:27 AM UTC

PDB ID : 6GL1 / pdb_00006gl1
Title : HLA-E*01:03 in complex with the HIV epitope, RL9HIV
Authors : Walters, L.C.; Gillespie, G.M.; McMichael, A.J.; Rozbesky, D.; Jones, E.Y.; Harlos, K.
Deposited on : 2018-05-22
Resolution : 2.62 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:

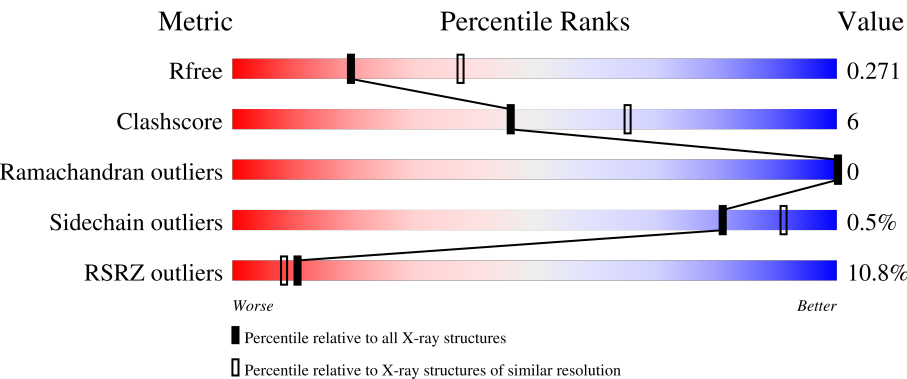
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

1 Overall quality at a glance i

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

The reported resolution of this entry is 2.62 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	4951 (2.64-2.60)
Clashscore	190562	5303 (2.64-2.60)
Ramachandran outliers	187476	5217 (2.64-2.60)
Sidechain outliers	187428	5217 (2.64-2.60)
RSRZ outliers	180081	4950 (2.64-2.60)

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	274	
1	C	274	
1	E	274	
1	G	274	
2	B	100	

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Mol	Chain	Length	Quality of chain
2	D	100	
2	F	100	
2	H	100	
3	P	9	
3	Q	9	
3	R	9	
3	T	9	

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 12700 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 MHC class I antigen.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	274	Total	C	N	O	S	0	0	0
			2231	1394	399	431	7			
1	C	273	Total	C	N	O	S	0	0	0
			2227	1392	398	430	7			
1	E	274	Total	C	N	O	S	0	0	0
			2231	1394	399	431	7			
1	G	249	Total	C	N	O	S	0	0	0
			2030	1271	363	389	7			

- Molecule 2 is a protein called Beta-2-microglobulin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	100	Total	C	N	O	S	0	0	0
			837	533	141	159	4			
2	D	100	Total	C	N	O	S	0	0	0
			837	533	141	159	4			
2	F	100	Total	C	N	O	S	0	0	0
			837	533	141	159	4			
2	H	100	Total	C	N	O	S	0	0	0
			837	533	141	159	4			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	1	MET	-	initiating methionine	UNP P61769
D	1	MET	-	initiating methionine	UNP P61769
F	1	MET	-	initiating methionine	UNP P61769
H	1	MET	-	initiating methionine	UNP P61769

- Molecule 3 is a protein called ARG-MET-TYR-SER-PRO-THR-SER-ILE-LEU.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	P	9	Total	C	N	O	S	0	0	0
			74	47	12	14	1			
3	R	9	Total	C	N	O	S	0	0	0
			74	47	12	14	1			
3	T	9	Total	C	N	O	S	0	0	0
			74	47	12	14	1			
3	Q	9	Total	C	N	O	S	0	0	0
			74	47	12	14	1			

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

- Molecule 5 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	G	1	Total 1	Zn 1	0	0

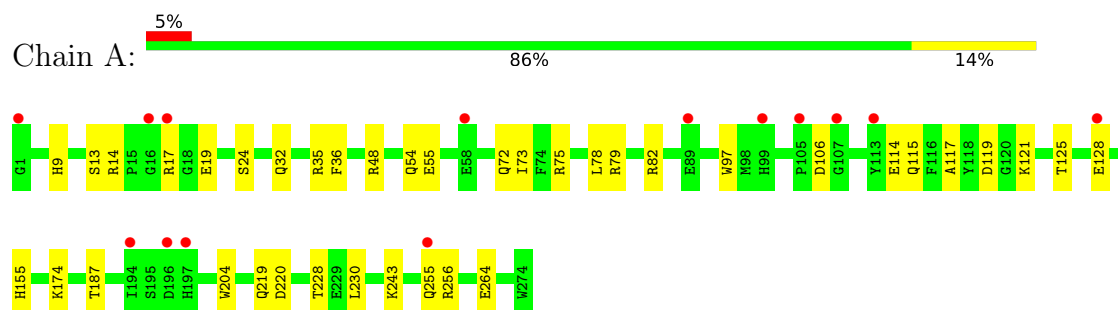
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	57	Total 57	O 57	0	0
6	B	25	Total 25	O 25	0	0
6	C	57	Total 57	O 57	0	0
6	D	27	Total 27	O 27	0	0
6	E	43	Total 43	O 43	0	0
6	F	19	Total 19	O 19	0	0
6	G	40	Total 40	O 40	0	0
6	H	29	Total 29	O 29	0	0
6	P	3	Total 3	O 3	0	0
6	R	4	Total 4	O 4	0	0
6	T	1	Total 1	O 1	0	0
6	Q	1	Total 1	O 1	0	0

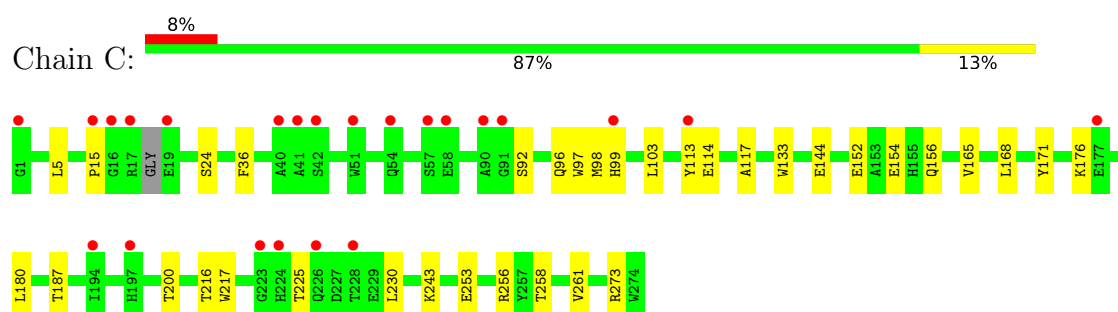
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

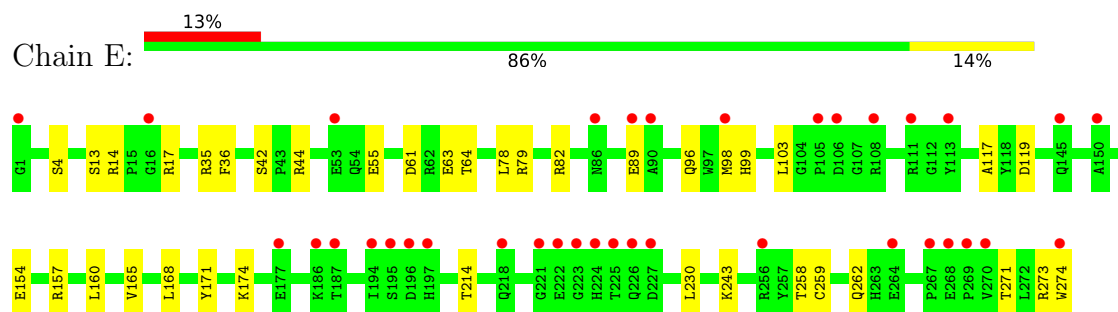
- Molecule 1: MHC class I antigen



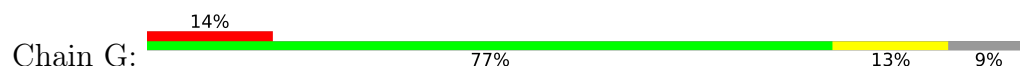
- Molecule 1: MHC class I antigen

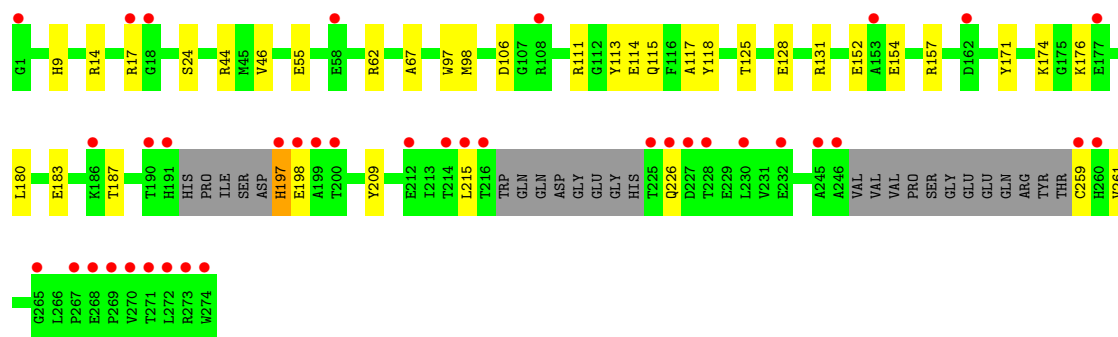


- Molecule 1: MHC class I antigen

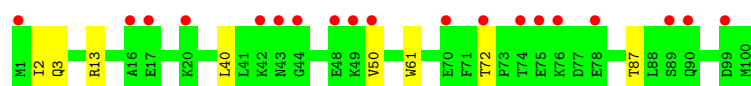


- Molecule 1: MHC class I antigen

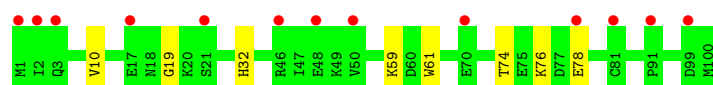




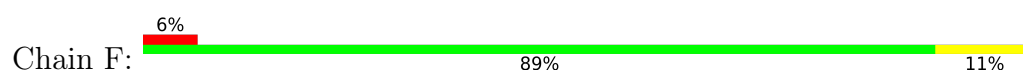
• Molecule 2: Beta-2-microglobulin



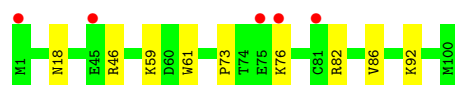
• Molecule 2: Beta-2-microglobulin



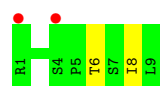
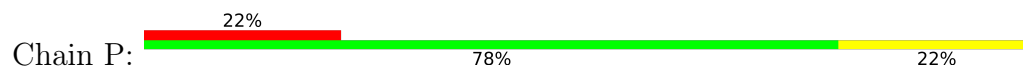
• Molecule 2: Beta-2-microglobulin



• Molecule 2: Beta-2-microglobulin

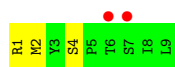


• Molecule 3: ARG-MET-TYR-SER-PRO-THR-SER-ILE-LEU

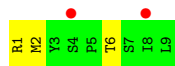


• Molecule 3: ARG-MET-TYR-SER-PRO-THR-SER-ILE-LEU





- Molecule 3: ARG-MET-TYR-SER-PRO-THR-SER-ILE-LEU



- Molecule 3: ARG-MET-TYR-SER-PRO-THR-SER-ILE-LEU



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	244.82Å 48.22Å 153.02Å 90.00° 117.38° 90.00°	Depositor
Resolution (Å)	135.88 – 2.62 135.88 – 2.62	Depositor EDS
% Data completeness (in resolution range)	85.3 (135.88-2.62) 80.5 (135.88-2.62)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.79 (at 2.62Å)	Xtriage
Refinement program	PHENIX (1.11.1_2575: ???)	Depositor
R, R_{free}	0.223 , 0.273 0.221 , 0.271	Depositor DCC
R_{free} test set	2112 reflections (4.34%)	wwPDB-VP
Wilson B-factor (Å ²)	16.4	Xtriage
Anisotropy	0.202	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 55.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	12700	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 29.91 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.4308e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, 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	0.15	0/2297	0.40	0/3123
1	C	0.15	0/2292	0.37	0/3115
1	E	0.15	0/2297	0.35	0/3123
1	G	0.17	0/2086	0.40	0/2829
2	B	0.11	0/860	0.34	0/1162
2	D	0.12	0/860	0.34	0/1162
2	F	0.11	0/860	0.32	0/1162
2	H	0.12	0/860	0.33	0/1162
3	P	0.13	0/75	0.41	0/99
3	Q	0.14	0/75	0.40	0/99
3	R	0.14	0/75	0.40	0/99
3	T	0.14	0/75	0.37	0/99
All	All	0.15	0/12712	0.37	0/17234

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	2231	0	2068	31	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	2227	0	2064	23	0
1	E	2231	0	2068	29	0
1	G	2030	0	1891	31	0
2	B	837	0	803	7	0
2	D	837	0	803	8	0
2	F	837	0	803	7	0
2	H	837	0	803	9	0
3	P	74	0	79	2	0
3	Q	74	0	79	3	0
3	R	74	0	79	4	0
3	T	74	0	79	4	0
4	A	10	0	0	0	0
4	C	5	0	0	0	0
4	E	10	0	0	1	0
4	H	5	0	0	0	0
5	G	1	0	0	0	0
6	A	57	0	0	9	0
6	B	25	0	0	2	0
6	C	57	0	0	0	1
6	D	27	0	0	3	0
6	E	43	0	0	7	1
6	F	19	0	0	1	0
6	G	40	0	0	8	0
6	H	29	0	0	3	0
6	P	3	0	0	0	0
6	Q	1	0	0	1	0
6	R	4	0	0	2	0
6	T	1	0	0	0	0
All	All	12700	0	11619	138	1

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 6.

All (138) 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:E:98:MET:SD	6:E:443:HOH:O	2.07	1.10
2:H:76:LYS:O	6:H:301:HOH:O	1.84	0.94
1:C:176:LYS:HG3	1:C:180:LEU:HD12	1.51	0.92
2:F:35:ASP:O	6:F:201:HOH:O	1.91	0.87
2:D:78:GLU:O	6:D:201:HOH:O	1.93	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:115:GLN:NE2	6:G:402:HOH:O	2.08	0.80
1:A:228:THR:OG1	6:A:401:HOH:O	1.99	0.80
1:C:258:THR:HG22	1:C:273:ARG:HE	1.49	0.77
1:G:118:TYR:O	6:G:401:HOH:O	2.02	0.77
1:A:72:GLN:NE2	6:A:404:HOH:O	2.08	0.75
3:R:4:SER:O	6:R:101:HOH:O	2.04	0.75
1:G:176:LYS:HG3	1:G:180:LEU:HD12	1.69	0.74
1:A:17:ARG:O	6:A:403:HOH:O	2.05	0.74
1:A:19:GLU:OE2	6:A:402:HOH:O	2.05	0.73
2:F:48:GLU:HG2	2:F:49:LYS:HG3	1.69	0.72
4:E:301:SO4:O1	6:E:403:HOH:O	2.06	0.72
1:E:274:TRP:OXT	6:E:404:HOH:O	2.09	0.70
1:G:98:MET:HE1	2:H:59:LYS:HA	1.73	0.69
2:H:92:LYS:NZ	6:H:303:HOH:O	2.22	0.68
1:A:75:ARG:HE	1:A:79:ARG:HH22	1.41	0.67
1:E:63:GLU:OE1	6:E:405:HOH:O	2.12	0.67
1:E:273:ARG:NH2	6:E:406:HOH:O	2.28	0.67
2:H:46:ARG:HE	2:H:82:ARG:HH12	1.42	0.66
1:E:14:ARG:HD2	1:E:17:ARG:HH11	1.59	0.66
1:C:152:GLU:OE2	3:Q:6:THR:OG1	2.14	0.64
3:R:1:ARG:NH1	6:R:103:HOH:O	2.30	0.63
2:B:13:ARG:NH2	6:B:201:HOH:O	2.18	0.62
1:G:187:THR:HG21	1:G:261:VAL:HG21	1.82	0.61
2:D:74:THR:HG22	2:D:76:LYS:H	1.67	0.60
2:D:10:VAL:O	6:D:202:HOH:O	2.16	0.60
1:E:258:THR:HG22	1:E:273:ARG:HE	1.66	0.60
1:E:4:SER:OG	6:E:401:HOH:O	1.91	0.60
1:G:115:GLN:HG2	1:G:125:THR:HG23	1.83	0.60
1:E:154:GLU:HG3	1:E:157:ARG:HH22	1.67	0.60
1:G:97:TRP:HE1	1:G:114:GLU:HG2	1.68	0.59
6:E:420:HOH:O	3:R:2:MET:HE3	2.03	0.59
1:A:119:ASP:OD2	6:A:405:HOH:O	2.16	0.58
1:G:67:ALA:HA	3:T:2:MET:HE1	1.86	0.58
1:E:214:THR:HB	1:E:262:GLN:HB2	1.85	0.57
2:B:3:GLN:HB3	2:B:87:THR:HG22	1.87	0.57
1:A:117:ALA:HB2	2:B:61:TRP:CE2	2.39	0.57
1:C:230:LEU:HD22	1:C:243:LYS:HE3	1.88	0.56
1:C:98:MET:HE1	2:D:59:LYS:HA	1.87	0.56
1:E:82:ARG:NE	1:E:89:GLU:HG2	2.22	0.55
1:G:117:ALA:HB2	2:H:61:TRP:CE2	2.41	0.55
1:C:103:LEU:HG	1:C:168:LEU:HD23	1.88	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:230:LEU:HD22	1:E:243:LYS:HE3	1.90	0.54
1:C:97:TRP:HE1	1:C:114:GLU:HG2	1.73	0.53
2:D:32:HIS:ND1	6:D:205:HOH:O	2.33	0.53
1:E:154:GLU:N	1:E:154:GLU:OE1	2.42	0.52
1:A:75:ARG:HE	1:A:79:ARG:NH2	2.08	0.52
1:G:44:ARG:HH21	1:G:46:VAL:HG12	1.74	0.52
1:A:97:TRP:HE1	1:A:114:GLU:HG2	1.75	0.51
1:G:226:GLN:O	6:G:403:HOH:O	2.19	0.51
1:C:96:GLN:HB2	1:C:117:ALA:HB3	1.91	0.51
1:G:9:HIS:CE1	3:T:2:MET:HE3	2.46	0.51
1:E:35:ARG:HG2	1:E:36:PHE:N	2.26	0.51
1:G:131:ARG:NH1	6:G:404:HOH:O	2.22	0.51
1:E:96:GLN:HB2	1:E:117:ALA:HB3	1.93	0.51
1:G:97:TRP:HE1	1:G:114:GLU:CG	2.24	0.50
1:A:9:HIS:ND1	1:A:24:SER:OG	2.33	0.50
1:A:79:ARG:HG2	1:A:82:ARG:NH2	2.27	0.49
1:C:117:ALA:HB2	2:D:61:TRP:CE2	2.47	0.49
1:G:113:TYR:HB2	6:G:415:HOH:O	2.12	0.49
1:A:230:LEU:HD11	1:A:243:LYS:HE3	1.95	0.49
1:C:253:GLU:HB3	1:C:256:ARG:HD3	1.95	0.49
1:G:62:ARG:NH2	6:G:407:HOH:O	2.44	0.49
1:C:187:THR:HG21	1:C:261:VAL:HG21	1.94	0.49
1:E:14:ARG:HD2	1:E:17:ARG:NH1	2.27	0.49
1:E:44:ARG:HA	1:E:64:THR:HG23	1.95	0.48
1:E:171:TYR:OH	3:R:1:ARG:N	2.45	0.48
1:A:54:GLN:HG3	2:D:19:GLY:HA3	1.95	0.48
1:C:24:SER:HB2	1:C:36:PHE:HB3	1.95	0.48
1:E:154:GLU:HG3	1:E:157:ARG:NH2	2.27	0.48
2:B:72:THR:N	6:B:205:HOH:O	2.46	0.48
1:G:215:LEU:HD22	1:G:261:VAL:HG22	1.95	0.48
1:E:103:LEU:HD21	1:E:165:VAL:HG13	1.94	0.47
2:H:46:ARG:NE	2:H:82:ARG:HH12	2.10	0.47
1:A:75:ARG:NH1	6:A:402:HOH:O	2.40	0.47
1:A:115:GLN:HG2	1:A:125:THR:HG23	1.95	0.47
1:E:55:GLU:OE2	1:E:174:LYS:HE3	2.14	0.47
1:A:220:ASP:OD1	1:A:256:ARG:HB3	2.15	0.47
2:F:46:ARG:HE	2:F:82:ARG:HH21	1.63	0.47
1:C:171:TYR:OH	3:Q:1:ARG:N	2.42	0.46
1:E:117:ALA:HB2	2:F:61:TRP:CE2	2.50	0.46
1:E:79:ARG:HG2	1:E:82:ARG:HH22	1.80	0.46
1:A:121:LYS:HB2	2:B:2:ILE:HD12	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:197:HIS:N	6:G:409:HOH:O	2.49	0.45
1:G:9:HIS:ND1	1:G:24:SER:OG	2.34	0.45
1:G:117:ALA:HB2	2:H:61:TRP:CD2	2.51	0.45
1:G:154:GLU:OE1	1:G:154:GLU:N	2.50	0.45
1:A:264:GLU:HG3	6:A:409:HOH:O	2.16	0.45
1:E:119:ASP:HB3	2:F:1:MET:HG3	1.97	0.45
1:A:14:ARG:HB3	1:A:17:ARG:HB2	1.99	0.45
1:A:155:HIS:ND1	3:P:6:THR:HG21	2.32	0.45
1:C:5:LEU:HB2	1:C:168:LEU:HD13	1.98	0.45
1:C:217:TRP:O	1:C:225:THR:OG1	2.29	0.45
1:G:171:TYR:OH	3:T:1:ARG:N	2.35	0.45
1:A:35:ARG:HG2	1:A:36:PHE:N	2.31	0.45
1:C:103:LEU:HD21	1:C:165:VAL:HG13	1.99	0.44
1:E:14:ARG:HB3	1:E:17:ARG:HB2	1.98	0.44
1:E:13:SER:HB3	1:E:78:LEU:HD13	1.99	0.44
1:A:97:TRP:HE1	1:A:114:GLU:CG	2.30	0.44
1:E:99:HIS:CD2	1:E:160:LEU:HD21	2.53	0.43
2:H:18:ASN:O	6:H:302:HOH:O	2.21	0.43
1:A:128:GLU:HG2	6:A:440:HOH:O	2.17	0.43
1:E:103:LEU:HG	1:E:168:LEU:HD23	2.00	0.43
1:A:32:GLN:NE2	1:A:48:ARG:HG3	2.33	0.43
1:A:13:SER:HB3	1:A:78:LEU:HD13	2.01	0.43
3:Q:4:SER:O	6:Q:101:HOH:O	2.21	0.43
1:E:44:ARG:NH2	1:E:61:ASP:OD1	2.52	0.43
1:C:97:TRP:HE1	1:C:114:GLU:CG	2.32	0.42
1:A:187:THR:HA	1:A:204:TRP:O	2.19	0.42
1:A:106:ASP:OD1	1:A:106:ASP:N	2.52	0.42
1:A:117:ALA:HB2	2:B:61:TRP:CD2	2.55	0.42
1:E:259:CYS:O	1:E:271:THR:HA	2.19	0.42
1:G:55:GLU:OE2	1:G:174:LYS:HE3	2.19	0.42
1:C:98:MET:HG2	1:C:113:TYR:HE1	1.85	0.42
1:C:117:ALA:HB2	2:D:61:TRP:CD2	2.54	0.42
1:G:183:GLU:HB3	1:G:209:TYR:N	2.34	0.42
1:G:259:CYS:N	6:G:410:HOH:O	2.52	0.42
1:C:99:HIS:NE2	1:C:156:GLN:OE1	2.53	0.42
1:G:152:GLU:OE2	3:T:6:THR:OG1	2.32	0.42
1:A:255:GLN:HB2	6:A:426:HOH:O	2.20	0.41
1:G:14:ARG:HD2	1:G:17:ARG:HH12	1.85	0.41
2:F:7:LYS:O	2:F:28:VAL:HA	2.21	0.41
1:A:73:ILE:HG23	3:P:8:ILE:HD13	2.02	0.41
1:C:133:TRP:HB2	1:C:144:GLU:HG3	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:111:ARG:NH1	1:G:128:GLU:HG3	2.36	0.41
1:C:15:PRO:HD3	1:C:92:SER:HB2	2.02	0.41
1:C:154:GLU:OE1	1:C:154:GLU:N	2.53	0.41
1:A:55:GLU:OE2	1:A:174:LYS:HE3	2.21	0.41
2:B:40:LEU:HB2	2:B:50:VAL:HG11	2.03	0.41
1:G:183:GLU:HB3	1:G:209:TYR:H	1.86	0.41
2:H:18:ASN:HA	2:H:73:PRO:O	2.21	0.41
1:G:106:ASP:N	1:G:106:ASP:OD1	2.54	0.41
2:F:74:THR:OG1	2:F:77:ASP:OD2	2.35	0.40
1:G:154:GLU:HG3	1:G:157:ARG:NH2	2.37	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:C:417:HOH:O	6:E:440:HOH:O[3_545]	2.13	0.07

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	272/274 (99%)	268 (98%)	4 (2%)	0	100	100
1	C	269/274 (98%)	264 (98%)	5 (2%)	0	100	100
1	E	272/274 (99%)	268 (98%)	4 (2%)	0	100	100
1	G	241/274 (88%)	236 (98%)	5 (2%)	0	100	100
2	B	98/100 (98%)	97 (99%)	1 (1%)	0	100	100
2	D	98/100 (98%)	97 (99%)	1 (1%)	0	100	100
2	F	98/100 (98%)	98 (100%)	0	0	100	100
2	H	98/100 (98%)	96 (98%)	2 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	P	7/9 (78%)	7 (100%)	0	0	100	100
3	Q	7/9 (78%)	7 (100%)	0	0	100	100
3	R	7/9 (78%)	7 (100%)	0	0	100	100
3	T	7/9 (78%)	7 (100%)	0	0	100	100
All	All	1474/1532 (96%)	1452 (98%)	22 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	235/235 (100%)	234 (100%)	1 (0%)	84	92
1	C	235/235 (100%)	233 (99%)	2 (1%)	70	86
1	E	235/235 (100%)	234 (100%)	1 (0%)	84	92
1	G	213/235 (91%)	211 (99%)	2 (1%)	70	86
2	B	95/95 (100%)	95 (100%)	0	100	100
2	D	95/95 (100%)	95 (100%)	0	100	100
2	F	95/95 (100%)	95 (100%)	0	100	100
2	H	95/95 (100%)	94 (99%)	1 (1%)	65	83
3	P	9/9 (100%)	9 (100%)	0	100	100
3	Q	9/9 (100%)	9 (100%)	0	100	100
3	R	9/9 (100%)	9 (100%)	0	100	100
3	T	9/9 (100%)	9 (100%)	0	100	100
All	All	1334/1356 (98%)	1327 (100%)	7 (0%)	81	91

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	219	GLN

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Mol	Chain	Res	Type
1	C	200	THR
1	C	216	THR
1	E	42	SER
1	G	197	HIS
1	G	198	GLU
2	H	86	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	262	GLN
2	D	52	HIS
1	E	192	HIS
1	E	255	GLN
1	G	72	GLN
2	H	90	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](#)

Of 7 ligands modelled in this entry, 1 is monoatomic - leaving 6 for Mogul analysis.

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	E	302	-	4,4,4	0.23	0	6,6,6	0.05	0
4	SO4	A	302	-	4,4,4	0.24	0	6,6,6	0.09	0
4	SO4	H	201	-	4,4,4	0.24	0	6,6,6	1.41	0
4	SO4	C	301	-	4,4,4	0.23	0	6,6,6	0.12	0
4	SO4	E	301	-	4,4,4	0.28	0	6,6,6	0.10	0
4	SO4	A	301	-	4,4,4	0.21	0	6,6,6	0.14	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.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	E	301	SO4	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.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	274/274 (100%)	0.52	14 (5%) 33 28	15, 30, 54, 77	0
1	C	273/274 (99%)	0.78	23 (8%) 17 13	17, 35, 59, 75	0
1	E	274/274 (100%)	0.96	36 (13%) 7 5	13, 41, 66, 74	0
1	G	249/274 (90%)	0.94	38 (15%) 5 4	17, 37, 71, 99	0
2	B	100/100 (100%)	1.11	19 (19%) 3 2	20, 44, 71, 84	0
2	D	100/100 (100%)	0.86	13 (13%) 7 6	22, 40, 60, 73	0
2	F	100/100 (100%)	0.77	6 (6%) 27 23	17, 41, 72, 79	0
2	H	100/100 (100%)	0.50	5 (5%) 34 29	15, 33, 54, 73	0
3	P	9/9 (100%)	1.29	2 (22%) 2 2	32, 41, 51, 60	0
3	Q	9/9 (100%)	1.59	2 (22%) 2 2	28, 43, 55, 62	0
3	R	9/9 (100%)	1.39	2 (22%) 2 2	33, 42, 51, 59	0
3	T	9/9 (100%)	1.15	2 (22%) 2 2	28, 40, 46, 51	0
All	All	1506/1532 (98%)	0.82	162 (10%) 11 8	13, 36, 65, 99	0

All (162) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	16	GLY	6.0
1	E	223	GLY	5.6
1	G	216	THR	5.0
1	E	221	GLY	4.9
1	C	1	GLY	4.9
1	A	197	HIS	4.8
1	G	197	HIS	4.6
3	Q	4	SER	4.5
1	G	260	HIS	4.5
1	E	224	HIS	4.2
2	D	46	ARG	4.1

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Mol	Chain	Res	Type	RSRZ
1	G	200	THR	4.1
1	C	40	ALA	4.1
1	C	41	ALA	4.1
1	C	223	GLY	3.8
2	B	48	GLU	3.7
2	D	1	MET	3.7
1	G	58	GLU	3.7
3	R	6	THR	3.6
2	D	48	GLU	3.6
1	C	177	GLU	3.6
1	G	273	ARG	3.6
2	F	18	ASN	3.6
1	G	225	THR	3.5
1	C	91	GLY	3.5
1	G	18	GLY	3.4
1	A	255	GLN	3.4
2	D	78	GLU	3.3
1	C	17	ARG	3.3
1	E	227	ASP	3.3
1	C	58	GLU	3.2
1	G	271	THR	3.2
1	E	268	GLU	3.2
2	B	16	ALA	3.2
1	A	128	GLU	3.2
1	G	269	PRO	3.2
1	G	198	GLU	3.1
2	B	74	THR	3.1
1	G	274	TRP	3.1
1	G	1	GLY	3.1
1	C	113	TYR	3.0
1	G	267	PRO	3.0
2	H	76	LYS	3.0
1	E	1	GLY	3.0
1	G	199	ALA	3.0
1	G	191	HIS	3.0
1	G	226	GLN	3.0
2	D	21	SER	2.9
2	B	76	LYS	2.9
3	P	4	SER	2.9
2	B	1	MET	2.9
2	D	99	ASP	2.9
1	E	264	GLU	2.9

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Mol	Chain	Res	Type	RSRZ
2	F	75	GLU	2.9
2	B	44	GLY	2.9
2	H	45	GLU	2.9
1	G	246	ALA	2.9
1	G	268	GLU	2.8
1	A	194	ILE	2.8
1	E	106	ASP	2.8
1	E	113	TYR	2.8
1	E	90	ALA	2.8
1	E	222	GLU	2.8
1	G	17	ARG	2.8
2	B	99	ASP	2.8
2	B	90	GLN	2.8
2	F	20	LYS	2.8
1	E	105	PRO	2.7
2	D	2	ILE	2.7
1	A	1	GLY	2.7
2	B	43	ASN	2.7
1	G	186	LYS	2.7
2	B	49	LYS	2.7
1	A	89	GLU	2.7
1	E	177	GLU	2.7
1	G	232	GLU	2.7
3	P	1	ARG	2.7
1	G	270	VAL	2.7
2	D	3	GLN	2.7
1	E	196	ASP	2.7
3	T	8	ILE	2.6
1	G	177	GLU	2.6
1	G	265	GLY	2.6
2	B	20	LYS	2.6
3	R	7	SER	2.6
1	E	225	THR	2.6
1	G	214	THR	2.6
1	A	105	PRO	2.6
1	G	245	ALA	2.6
3	T	4	SER	2.6
1	G	190	THR	2.5
1	E	89	GLU	2.5
1	C	54	GLN	2.5
1	C	228	THR	2.5
1	A	196	ASP	2.5

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Mol	Chain	Res	Type	RSRZ
1	E	256	ARG	2.5
1	G	230	LEU	2.5
1	A	58	GLU	2.5
1	E	53	GLU	2.5
1	G	212	GLU	2.5
1	C	15	PRO	2.5
1	C	90	ALA	2.5
1	E	226	GLN	2.5
2	D	17	GLU	2.5
2	D	70	GLU	2.5
1	C	42	SER	2.4
1	E	197	HIS	2.4
2	F	76	LYS	2.4
1	C	224	HIS	2.4
1	C	226	GLN	2.4
1	G	228	THR	2.4
1	C	57	SER	2.4
2	B	17	GLU	2.4
2	H	75	GLU	2.4
1	E	218	GLN	2.4
2	D	50	VAL	2.4
1	C	99	HIS	2.3
1	A	107	GLY	2.3
2	F	19	GLY	2.3
1	E	86	ASN	2.3
2	H	1	MET	2.3
1	E	269	PRO	2.3
2	B	78	GLU	2.3
2	D	91	PRO	2.3
1	C	51	TRP	2.3
1	G	215	LEU	2.3
1	G	108	ARG	2.3
1	E	98	MET	2.2
2	B	70	GLU	2.2
2	B	89	SER	2.2
2	B	50	VAL	2.2
1	E	150	ALA	2.2
1	C	19	GLU	2.2
1	E	270	VAL	2.2
2	F	50	VAL	2.2
1	E	194	ILE	2.2
1	E	16	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
1	G	272	LEU	2.2
1	G	227	ASP	2.2
1	C	194	ILE	2.2
1	C	197	HIS	2.2
1	A	17	ARG	2.1
1	G	162	ASP	2.1
1	E	195	SER	2.1
1	E	274	TRP	2.1
1	E	187	THR	2.1
3	Q	5	PRO	2.1
1	E	145	GLN	2.1
1	E	111	ARG	2.1
2	B	72	THR	2.1
1	E	186	LYS	2.1
2	D	81	CYS	2.1
2	H	81	CYS	2.1
2	B	42	LYS	2.1
2	B	75	GLU	2.0
1	E	267	PRO	2.0
1	A	16	GLY	2.0
1	A	113	TYR	2.0
1	G	259	CYS	2.0
1	G	153	ALA	2.0
1	A	99	HIS	2.0
1	E	108	ARG	2.0

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

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

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	SO4	C	301	5/5	0.71	0.28	77,79,81,83	0
4	SO4	E	302	5/5	0.74	0.20	65,66,67,70	0
4	SO4	A	302	5/5	0.84	0.13	61,63,64,68	0
4	SO4	E	301	5/5	0.88	0.13	69,69,73,77	0
4	SO4	H	201	5/5	0.93	0.21	51,52,55,55	5
4	SO4	A	301	5/5	0.94	0.12	55,56,61,61	0
5	ZN	G	301	1/1	0.98	0.12	34,34,34,34	1

6.5 Other polymers [i](#)

There are no such residues in this entry.