



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 05:33 AM UTC

PDB ID : 4JFE / pdb_00004jfe
Title : Preservation of peptide specificity during TCR-MHC contact dominated affinity enhancement of a melanoma-specific TCR
Authors : Rizkallah, P.J.; Cole, D.K.; Madura, F.; Sewell, A.K.
Deposited on : 2013-02-28
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

<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

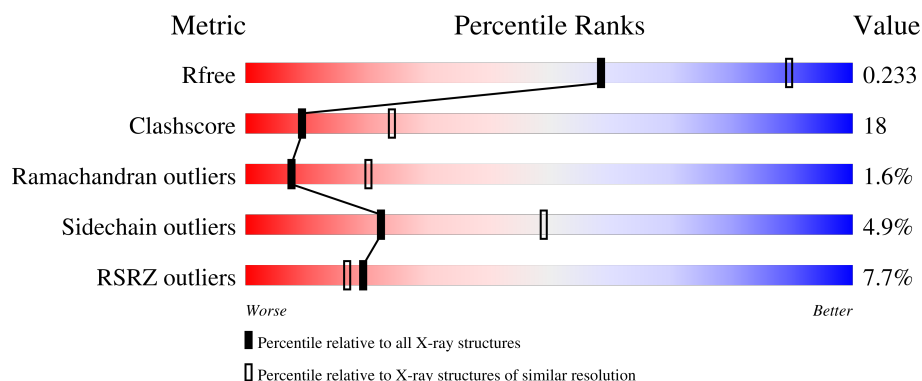
1 Overall quality at a glance

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))
R_{free}	180053	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (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 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	276	14% 71% 24% 5%
2	B	100	% 84% 15%
3	C	10	50% 40% 10%
4	D	197	9% 69% 22% 8%
5	E	245	2% 77% 20%

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 6765 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called HLA class I histocompatibility antigen, A-2 alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	276	Total	C	N	O	S	0	0	0
			2253	1408	410	426	9			

- 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			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	0	MET	-	initiating methionine	UNP P61769

- Molecule 3 is a protein called Melanoma peptide L7A.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	10	Total	C	N	O	0	0	0
			66	42	10	14			

- Molecule 4 is a protein called High Affinity TCR Alpha Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	197	Total	C	N	O	S	0	0	0
			1536	956	256	316	8			

- Molecule 5 is a protein called High Affinity TCR Beta Chain.

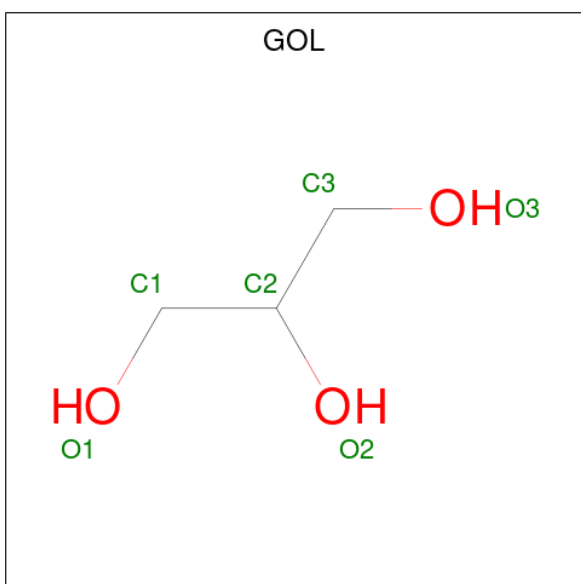
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	E	245	Total	C	N	O	S	0	3	0
			1975	1253	344	371	7			

- Molecule 6 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

- Molecule 7 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	D	1	Total	C	O	0	0
			6	3	3		
7	E	1	Total	C	O	0	0
			6	3	3		

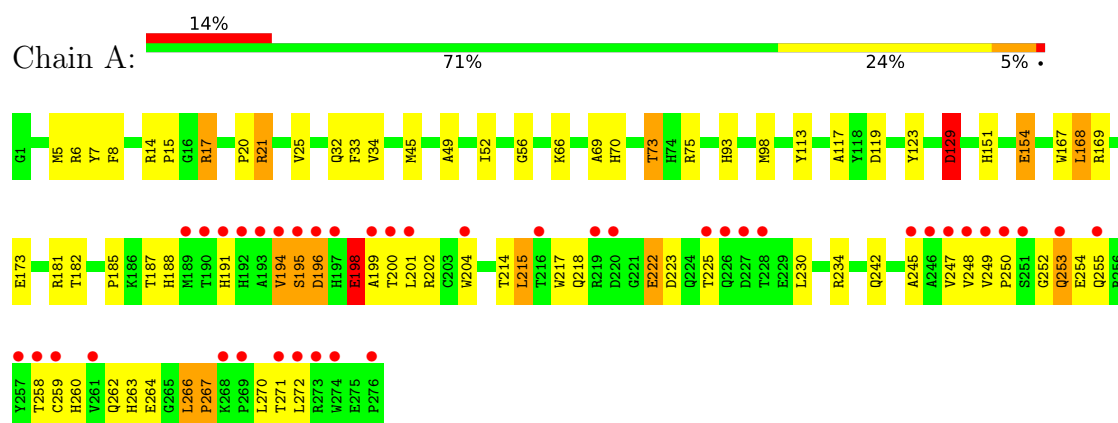
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	16	Total	O	0	0
			16	16		
8	B	10	Total	O	0	0
			10	10		
8	D	17	Total	O	0	0
			17	17		
8	E	23	Total	O	0	0
			23	23		

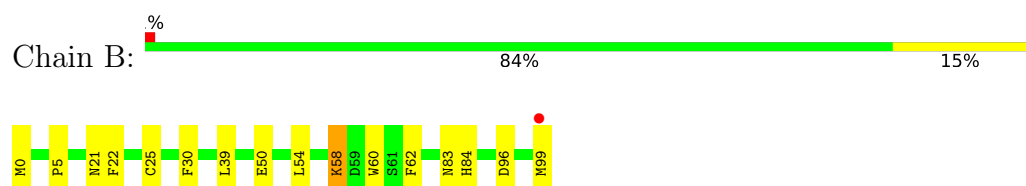
3 Residue-property plots [i](#)

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: HLA class I histocompatibility antigen, A-2 alpha chain



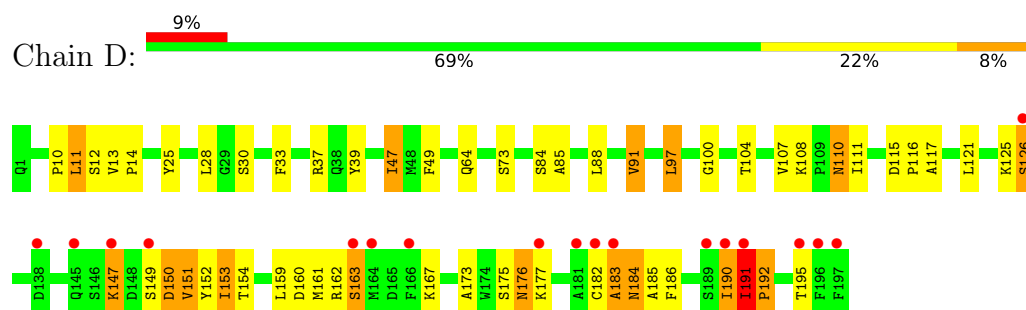
- Molecule 2: Beta-2-microglobulin



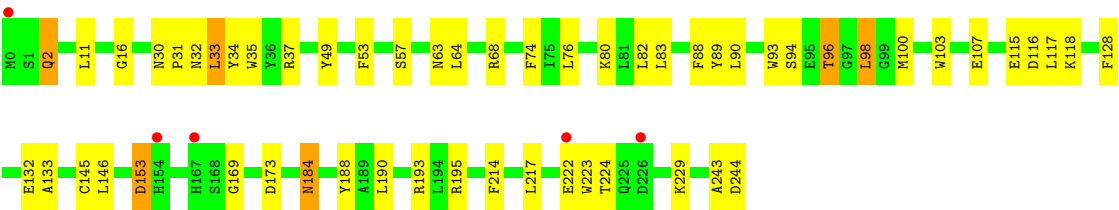
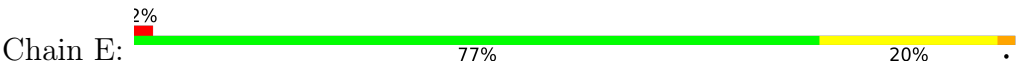
- Molecule 3: Melanoma peptide L7A



- Molecule 4: High Affinity TCR Alpha Chain



● Molecule 5: High Affinity TCR Beta Chain



4 Data and refinement statistics

Property	Value	Source
Space group	P 43	Depositor
Cell constants a, b, c, α , β , γ	121.61Å 121.61Å 82.25Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	68.13 – 2.70 68.13 – 2.70	Depositor EDS
% Data completeness (in resolution range)	100.0 (68.13-2.70) 99.9 (68.13-2.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.69 (at 2.69Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.202 , 0.250 (Not available) , 0.233	Depositor DCC
R_{free} test set	1674 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	54.0	Xtriage
Anisotropy	0.212	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 61.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.037 for h,-k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6765	wwPDB-VP
Average B, all atoms (Å ²)	62.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.06% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, 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	1.14	4/2319 (0.2%)	0.77	3/3149 (0.1%)
2	B	0.98	0/860	0.73	0/1162
3	C	1.87	1/65 (1.5%)	0.84	0/86
4	D	1.26	3/1568 (0.2%)	0.84	4/2123 (0.2%)
5	E	1.15	2/2036 (0.1%)	0.79	1/2776 (0.0%)
All	All	1.16	10/6848 (0.1%)	0.79	8/9296 (0.1%)

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	4	GLY	C-O	-8.69	1.18	1.24
1	A	123	TYR	CA-C	6.75	1.58	1.52
1	A	151	HIS	CA-C	-5.64	1.46	1.53
4	D	191	ILE	CA-C	5.59	1.58	1.52
1	A	98	MET	CA-C	-5.51	1.45	1.52
4	D	191	ILE	CA-CB	5.38	1.61	1.54
1	A	168	LEU	C-O	-5.29	1.18	1.24
5	E	11	LEU	C-O	-5.28	1.17	1.23
5	E	214	PHE	CA-C	-5.11	1.46	1.52
4	D	47	ILE	CA-CB	-5.03	1.48	1.54

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	129	ASP	N-CA-C	-7.70	97.52	109.76
5	E	153	ASP	N-CA-C	6.23	119.58	112.97
4	D	183	ALA	N-CA-C	-6.06	105.72	113.23
4	D	191	ILE	CA-C-N	5.76	127.05	119.84
4	D	191	ILE	C-N-CA	5.76	127.05	119.84
4	D	100	GLY	N-CA-C	-5.72	103.84	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	56	GLY	CA-C-N	5.33	125.15	119.28
1	A	56	GLY	C-N-CA	5.33	125.15	119.28

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2253	0	2103	112	0
2	B	837	0	803	13	0
3	C	66	0	73	7	0
4	D	1536	0	1448	79	0
5	E	1975	0	1872	52	0
6	A	10	0	0	1	0
6	D	5	0	0	0	0
6	E	5	0	0	0	0
7	D	6	0	8	0	0
7	E	6	0	8	0	0
8	A	16	0	0	0	0
8	B	10	0	0	0	0
8	D	17	0	0	0	0
8	E	23	0	0	0	0
All	All	6765	0	6315	235	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 18.

All (235) 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:A:263:HIS:O	1:A:266:LEU:HD22	1.27	1.26
1:A:215:LEU:HD12	1:A:215:LEU:O	1.44	1.18
4:D:190:ILE:HD13	4:D:191:ILE:N	1.65	1.11
2:B:21:ASN:HD22	2:B:22:PHE:N	1.47	1.10

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:21:ASN:ND2	2:B:22:PHE:H	1.47	1.09
4:D:150:ASP:O	4:D:151:VAL:HG22	1.52	1.08
1:A:258:THR:HG23	1:A:272:LEU:O	1.53	1.08
1:A:187:THR:HB	1:A:272:LEU:HD21	1.34	1.03
4:D:190:ILE:CD1	4:D:191:ILE:H	1.71	1.02
1:A:69:ALA:O	1:A:73:THR:HG22	1.63	0.96
4:D:190:ILE:HD13	4:D:191:ILE:H	0.80	0.94
4:D:183:ALA:C	4:D:184:ASN:HD22	1.78	0.91
1:A:263:HIS:O	1:A:266:LEU:CD2	2.18	0.90
1:A:17:ARG:HG3	1:A:17:ARG:HH11	1.34	0.89
4:D:161:MET:HE2	5:E:195:ARG:HB3	1.52	0.88
1:A:270:LEU:HD23	1:A:271:THR:N	1.89	0.88
4:D:190:ILE:HG23	4:D:192:PRO:HD3	1.55	0.87
1:A:199:ALA:O	1:A:250:PRO:CG	2.24	0.86
4:D:190:ILE:O	4:D:192:PRO:HD2	1.73	0.86
1:A:266:LEU:HD23	1:A:266:LEU:O	1.75	0.85
4:D:147:LYS:HA	4:D:147:LYS:NZ	1.92	0.85
1:A:199:ALA:O	1:A:250:PRO:HG2	1.75	0.84
1:A:167:TRP:CE2	3:C:1:GLU:HG3	2.13	0.84
1:A:222:GLU:OE2	1:A:223:ASP:N	2.08	0.84
1:A:187:THR:CB	1:A:272:LEU:HD21	2.08	0.83
4:D:49:PHE:CD2	5:E:100:MET:HE2	2.14	0.83
1:A:201:LEU:HB2	1:A:247:VAL:CG2	2.10	0.82
4:D:150:ASP:OD2	4:D:151:VAL:N	2.12	0.81
1:A:258:THR:HG23	1:A:272:LEU:C	2.07	0.79
4:D:159:LEU:HD11	5:E:169:GLY:HA2	1.66	0.78
1:A:255:GLN:HE21	1:A:255:GLN:HA	1.49	0.76
1:A:249:VAL:N	1:A:250:PRO:HD3	2.02	0.74
5:E:16:GLY:HA2	5:E:80:LYS:HG3	1.69	0.74
1:A:247:VAL:HG23	1:A:247:VAL:O	1.85	0.73
1:A:191:HIS:HB2	1:A:201:LEU:HD23	1.69	0.73
4:D:184:ASN:HD22	4:D:184:ASN:N	1.85	0.73
1:A:201:LEU:HB2	1:A:247:VAL:HG22	1.72	0.70
4:D:147:LYS:HA	4:D:147:LYS:HZ1	1.54	0.70
4:D:190:ILE:C	4:D:192:PRO:HD3	2.16	0.70
2:B:21:ASN:ND2	2:B:22:PHE:N	2.21	0.70
4:D:150:ASP:O	4:D:151:VAL:CG2	2.38	0.70
1:A:215:LEU:O	1:A:215:LEU:CD1	2.33	0.69
2:B:21:ASN:HD22	2:B:22:PHE:H	0.74	0.68
4:D:190:ILE:C	4:D:192:PRO:CD	2.66	0.68
1:A:254:GLU:N	1:A:254:GLU:OE1	2.27	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:E:30:ASN:O	5:E:96:THR:HB	1.94	0.67
1:A:191:HIS:CB	1:A:201:LEU:HD23	2.23	0.67
1:A:270:LEU:HD22	1:A:272:LEU:CD1	2.25	0.67
1:A:201:LEU:HB2	1:A:247:VAL:HG21	1.76	0.66
1:A:230:LEU:C	1:A:230:LEU:HD12	2.20	0.66
4:D:110:ASN:C	4:D:110:ASN:HD22	2.03	0.66
4:D:159:LEU:HD11	5:E:169:GLY:CA	2.26	0.66
1:A:270:LEU:HD23	1:A:271:THR:H	1.59	0.65
4:D:159:LEU:HD11	5:E:169:GLY:C	2.21	0.65
1:A:187:THR:HB	1:A:272:LEU:CD2	2.19	0.65
1:A:248:VAL:HG12	1:A:248:VAL:O	1.97	0.65
1:A:187:THR:HG21	1:A:272:LEU:HD11	1.78	0.65
1:A:25:VAL:HG23	1:A:32:GLN:HE21	1.61	0.64
1:A:66:LYS:HE2	3:C:2:LEU:HB2	1.77	0.64
4:D:195:THR:HG21	5:E:133:ALA:CB	2.28	0.64
1:A:194:VAL:O	1:A:195:SER:CB	2.44	0.64
4:D:161:MET:CE	5:E:195:ARG:HD3	2.28	0.64
4:D:190:ILE:O	4:D:192:PRO:CD	2.45	0.63
1:A:249:VAL:N	1:A:250:PRO:CD	2.61	0.63
1:A:33:PHE:CD2	1:A:34:VAL:HG13	2.34	0.63
1:A:188:HIS:HB3	1:A:204:TRP:HB2	1.80	0.62
1:A:270:LEU:HD22	1:A:272:LEU:HD12	1.81	0.62
4:D:162:ARG:O	4:D:163:SER:CB	2.46	0.62
1:A:215:LEU:HD11	1:A:245:ALA:HB2	1.80	0.62
5:E:184:ASN:C	5:E:184:ASN:HD22	2.07	0.62
1:A:198:GLU:OE2	1:A:198:GLU:O	2.17	0.62
4:D:195:THR:HG21	5:E:133:ALA:HB1	1.82	0.62
4:D:176:ASN:HD22	4:D:177:LYS:N	1.99	0.61
1:A:258:THR:HG21	1:A:271:THR:OG1	2.01	0.61
1:A:195:SER:O	1:A:196:ASP:HB3	2.00	0.61
1:A:17:ARG:HG3	1:A:17:ARG:NH1	2.13	0.60
2:B:83:ASN:HD22	2:B:84:HIS:H	1.50	0.59
1:A:191:HIS:HB2	1:A:200:THR:O	2.02	0.59
4:D:147:LYS:HA	4:D:147:LYS:HZ2	1.66	0.59
4:D:152:TYR:O	4:D:173:ALA:HA	2.03	0.59
1:A:7:TYR:OH	3:C:1:GLU:N	2.23	0.58
1:A:17:ARG:HH11	1:A:17:ARG:CG	2.12	0.58
4:D:159:LEU:HD12	4:D:160:ASP:N	2.19	0.58
1:A:182:THR:HG23	1:A:182:THR:O	2.03	0.58
4:D:153:ILE:HD13	4:D:186:PHE:CE2	2.38	0.58
1:A:266:LEU:CD2	1:A:266:LEU:O	2.52	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:96:ASP:HB3	2:B:99:MET:HB3	1.86	0.57
2:B:5:PRO:HB3	2:B:30:PHE:HB3	1.87	0.57
4:D:49:PHE:CD2	5:E:100:MET:CE	2.86	0.57
1:A:247:VAL:CG2	1:A:247:VAL:O	2.52	0.57
1:A:117:ALA:HB2	2:B:60:TRP:CE2	2.40	0.57
1:A:270:LEU:CD2	1:A:272:LEU:HD12	2.34	0.56
1:A:169:ARG:O	1:A:173:GLU:HG3	2.05	0.56
5:E:243:ALA:O	5:E:244:ASP:OD1	2.23	0.56
1:A:255:GLN:HA	1:A:255:GLN:NE2	2.20	0.56
4:D:161:MET:HE2	5:E:195:ARG:CB	2.29	0.56
1:A:234:ARG:HE	1:A:242:GLN:HE21	1.52	0.56
5:E:117:LEU:HD22	5:E:217:LEU:HD21	1.89	0.55
4:D:131:VAL:HG12	4:D:132:CYS:N	2.21	0.55
1:A:258:THR:CG2	1:A:271:THR:OG1	2.54	0.55
5:E:224:THR:O	5:E:224:THR:HG22	2.06	0.55
4:D:37:ARG:HB2	4:D:47:ILE:HD11	1.89	0.55
1:A:201:LEU:HD11	1:A:254:GLU:HB3	1.89	0.55
1:A:217:TRP:CZ2	1:A:259:CYS:SG	3.00	0.55
5:E:173:ASP:HB2	5:E:190:LEU:HD11	1.89	0.54
1:A:25:VAL:HG23	1:A:32:GLN:NE2	2.22	0.54
4:D:49:PHE:CD2	4:D:49:PHE:C	2.85	0.54
1:A:194:VAL:HG22	1:A:199:ALA:HA	1.88	0.54
4:D:190:ILE:CG2	4:D:192:PRO:HD3	2.32	0.54
5:E:83:LEU:HD11	5:E:115:GLU:HG2	1.89	0.54
1:A:49:ALA:O	1:A:52:ILE:HG22	2.07	0.54
5:E:116:ASP:CG	5:E:118:LYS:HG2	2.33	0.53
1:A:230:LEU:HD12	1:A:230:LEU:O	2.09	0.53
1:A:70:HIS:HA	1:A:73:THR:CG2	2.38	0.53
5:E:2:GLN:HG2	5:E:93:TRP:CH2	2.44	0.53
5:E:173:ASP:HB2	5:E:190:LEU:CD1	2.39	0.53
4:D:121:LEU:N	4:D:121:LEU:HD12	2.24	0.53
5:E:49:TYR:HB3	5:E:57:SER:HB3	1.91	0.52
4:D:88:LEU:N	4:D:88:LEU:HD12	2.24	0.52
4:D:110:ASN:HD22	4:D:111:ILE:N	2.06	0.52
1:A:5:MET:HB2	1:A:168:LEU:HD13	1.91	0.51
1:A:6:ARG:NE	1:A:113:TYR:OH	2.43	0.51
1:A:202:ARG:HG2	1:A:204:TRP:NE1	2.26	0.51
4:D:131:VAL:HG23	5:E:128:PHE:CE2	2.44	0.51
4:D:161:MET:HE3	5:E:195:ARG:HD3	1.92	0.51
1:A:69:ALA:CB	5:E:98:LEU:HD11	2.40	0.51
1:A:248:VAL:C	1:A:250:PRO:HD3	2.35	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:75:ARG:HD2	5:E:53:PHE:CZ	2.45	0.51
1:A:225:THR:O	1:A:225:THR:HG23	2.10	0.51
1:A:8:PHE:HB2	1:A:25:VAL:HG12	1.93	0.50
1:A:198:GLU:HA	1:A:250:PRO:HB2	1.92	0.50
4:D:153:ILE:HD13	4:D:186:PHE:HE2	1.73	0.50
4:D:11:LEU:HD23	4:D:12:SER:N	2.27	0.50
1:A:270:LEU:HD22	1:A:272:LEU:HD11	1.93	0.50
1:A:194:VAL:O	1:A:195:SER:OG	2.28	0.50
5:E:35:TRP:CD1	5:E:74:PHE:CE2	2.99	0.50
1:A:199:ALA:O	1:A:250:PRO:HG3	2.06	0.50
1:A:266:LEU:O	1:A:267:PRO:C	2.55	0.50
5:E:33:LEU:HD22	5:E:68:ARG:CZ	2.42	0.50
4:D:184:ASN:N	4:D:184:ASN:ND2	2.54	0.49
1:A:217:TRP:CZ3	1:A:258:THR:C	2.90	0.49
1:A:218:GLN:HB2	1:A:260:HIS:NE2	2.27	0.49
1:A:215:LEU:HD12	1:A:215:LEU:C	2.28	0.49
1:A:187:THR:CG2	1:A:270:LEU:HD13	2.43	0.49
1:A:254:GLU:N	1:A:254:GLU:CD	2.71	0.49
1:A:14:ARG:O	1:A:15:PRO:C	2.54	0.49
4:D:159:LEU:CD1	5:E:169:GLY:C	2.86	0.49
4:D:97:LEU:HD13	5:E:103:TRP:CZ3	2.48	0.48
4:D:191:ILE:O	4:D:192:PRO:O	2.31	0.48
4:D:30:SER:HB3	4:D:91:VAL:HG13	1.96	0.48
5:E:145:CYS:C	5:E:146:LEU:HD12	2.39	0.48
1:A:195:SER:O	1:A:196:ASP:CB	2.62	0.48
4:D:153:ILE:CD1	4:D:186:PHE:HE2	2.27	0.48
1:A:187:THR:CG2	1:A:272:LEU:HD11	2.44	0.47
1:A:272:LEU:HD12	1:A:272:LEU:N	2.28	0.47
4:D:190:ILE:HD13	4:D:191:ILE:HG12	1.95	0.47
1:A:194:VAL:O	1:A:195:SER:HB2	2.14	0.47
4:D:162:ARG:O	4:D:163:SER:OG	2.32	0.47
4:D:125:LYS:O	4:D:126:SER:C	2.58	0.47
5:E:30:ASN:N	5:E:31:PRO:CD	2.76	0.47
1:A:253:GLN:N	1:A:253:GLN:OE1	2.48	0.47
1:A:167:TRP:CD2	3:C:1:GLU:HG3	2.50	0.46
1:A:234:ARG:HE	1:A:242:GLN:NE2	2.13	0.46
1:A:230:LEU:C	1:A:230:LEU:CD1	2.88	0.46
1:A:263:HIS:CD2	1:A:264:GLU:O	2.68	0.46
1:A:270:LEU:CD2	1:A:272:LEU:CD1	2.93	0.46
5:E:153:ASP:HB3	5:E:188:TYR:CE2	2.51	0.46
1:A:20:PRO:HG2	1:A:75:ARG:HG3	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:218:GLN:HB3	1:A:222:GLU:O	2.16	0.46
4:D:159:LEU:HD12	4:D:160:ASP:H	1.80	0.46
4:D:162:ARG:O	4:D:163:SER:HB2	2.15	0.46
4:D:176:ASN:H	4:D:176:ASN:ND2	2.14	0.46
4:D:117:ALA:HA	4:D:192:PRO:HB3	1.97	0.46
4:D:151:VAL:HG12	4:D:175:SER:HB2	1.98	0.46
5:E:193:ARG:HD2	5:E:193:ARG:N	2.31	0.46
4:D:88:LEU:N	4:D:88:LEU:CD1	2.78	0.46
5:E:224:THR:O	5:E:224:THR:CG2	2.64	0.45
4:D:176:ASN:HD22	4:D:177:LYS:H	1.61	0.45
4:D:161:MET:CE	5:E:195:ARG:CG	2.94	0.45
4:D:190:ILE:CD1	4:D:191:ILE:N	2.49	0.45
1:A:201:LEU:O	1:A:247:VAL:HG22	2.17	0.45
1:A:129:ASP:O	1:A:129:ASP:CG	2.59	0.45
1:A:187:THR:HG21	1:A:270:LEU:HD13	1.98	0.45
2:B:83:ASN:HD22	2:B:84:HIS:N	2.15	0.45
4:D:33:PHE:N	4:D:33:PHE:CD2	2.85	0.44
1:A:258:THR:CG2	1:A:272:LEU:N	2.80	0.44
1:A:75:ARG:HD2	5:E:53:PHE:CE1	2.53	0.44
4:D:13:VAL:CG2	4:D:107:VAL:HG22	2.48	0.44
5:E:32:ASN:HB2	5:E:94:SER:OG	2.17	0.44
1:A:45:MET:CE	3:C:2:LEU:HD11	2.47	0.44
4:D:10:PRO:HA	4:D:104:THR:O	2.18	0.44
1:A:17:ARG:NH1	1:A:17:ARG:CG	2.77	0.44
1:A:191:HIS:CB	1:A:201:LEU:CD2	2.95	0.43
1:A:21:ARG:NH2	6:A:301:SO4:O1	2.51	0.43
2:B:58:LYS:CD	2:B:58:LYS:H	2.30	0.43
4:D:115:ASP:N	4:D:116:PRO:CD	2.82	0.43
3:C:7:ALA:O	3:C:8:LEU:HD23	2.18	0.43
5:E:30:ASN:N	5:E:31:PRO:HD3	2.32	0.43
1:A:154:GLU:CD	1:A:154:GLU:H	2.26	0.43
4:D:37:ARG:HD3	4:D:39:TYR:CZ	2.54	0.43
4:D:64:GLN:HB2	4:D:73:SER:HB2	2.00	0.42
1:A:202:ARG:HG2	1:A:204:TRP:CE2	2.55	0.42
5:E:35:TRP:CE2	5:E:76:LEU:HB2	2.54	0.42
5:E:88:PHE:CE2	5:E:90:LEU:HD21	2.53	0.42
2:B:58:LYS:H	2:B:58:LYS:HD2	1.85	0.42
3:C:7:ALA:HB3	5:E:98:LEU:HB2	2.01	0.42
4:D:131:VAL:CG1	4:D:132:CYS:N	2.82	0.42
2:B:54:LEU:HD11	2:B:62:PHE:CD1	2.55	0.42
4:D:25:TYR:CE1	4:D:33:PHE:CE1	3.08	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:161:MET:HE1	5:E:195:ARG:HG2	2.02	0.42
4:D:159:LEU:O	4:D:167:LYS:HA	2.20	0.42
5:E:63:ASN:O	5:E:64:LEU:HD12	2.20	0.41
2:B:25:CYS:HB2	2:B:39:LEU:HD21	2.01	0.41
5:E:223:TRP:CB	5:E:229:LYS:HE2	2.50	0.41
1:A:258:THR:CG2	1:A:272:LEU:C	2.86	0.41
1:A:93:HIS:HD2	1:A:119:ASP:OD1	2.03	0.41
1:A:198:GLU:OE2	1:A:198:GLU:C	2.63	0.41
4:D:153:ILE:HD12	4:D:154:THR:N	2.35	0.41
1:A:14:ARG:NH2	1:A:21:ARG:HG2	2.35	0.41
5:E:34:TYR:CE1	5:E:49:TYR:CD1	3.09	0.41
4:D:117:ALA:CA	4:D:192:PRO:HB3	2.51	0.41
1:A:129:ASP:OD1	1:A:129:ASP:C	2.62	0.41
1:A:185:PRO:HD3	1:A:263:HIS:CD2	2.56	0.41
5:E:222:GLU:HA	5:E:222:GLU:OE1	2.21	0.41
4:D:195:THR:HG21	5:E:133:ALA:HB2	2.03	0.41
5:E:16:GLY:HA2	5:E:80:LYS:CG	2.46	0.41
5:E:37[A]:ARG:HD3	5:E:89:TYR:CZ	2.56	0.41
1:A:214:THR:HB	1:A:262:GLN:HB2	2.03	0.40
4:D:13:VAL:HA	4:D:14:PRO:HD3	1.96	0.40
4:D:131:VAL:HG23	5:E:128:PHE:CD2	2.56	0.40
1:A:248:VAL:C	1:A:250:PRO:CD	2.95	0.40
4:D:84:SER:O	4:D:85:ALA:HB2	2.21	0.40
4:D:190:ILE:CD1	4:D:191:ILE:HG12	2.51	0.40
5:E:184:ASN:C	5:E:184:ASN:ND2	2.78	0.40
4:D:182:CYS:SG	4:D:185:ALA:HB2	2.62	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	274/276 (99%)	239 (87%)	29 (11%)	6 (2%)	5	14
2	B	98/100 (98%)	91 (93%)	7 (7%)	0	100	100
3	C	8/10 (80%)	8 (100%)	0	0	100	100
4	D	195/197 (99%)	178 (91%)	10 (5%)	7 (4%)	2	6
5	E	246/245 (100%)	239 (97%)	7 (3%)	0	100	100
All	All	821/828 (99%)	755 (92%)	53 (6%)	13 (2%)	7	20

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	195	SER
4	D	163	SER
4	D	192	PRO
4	D	150	ASP
4	D	151	VAL
4	D	191	ILE
1	A	196	ASP
1	A	252	GLY
4	D	126	SER
1	A	194	VAL
1	A	198	GLU
4	D	149	SER
1	A	267	PRO

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	232/232 (100%)	221 (95%)	11 (5%)	23	51
2	B	95/95 (100%)	92 (97%)	3 (3%)	34	64
3	C	6/6 (100%)	5 (83%)	1 (17%)	2	6
4	D	175/175 (100%)	163 (93%)	12 (7%)	14	34
5	E	214/211 (101%)	206 (96%)	8 (4%)	30	59

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	722/719 (100%)	687 (95%)	35 (5%)	22	50

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	ARG
1	A	21	ARG
1	A	73	THR
1	A	129	ASP
1	A	154	GLU
1	A	181	ARG
1	A	198	GLU
1	A	215	LEU
1	A	222	GLU
1	A	253	GLN
1	A	266	LEU
2	B	0	MET
2	B	50	GLU
2	B	58	LYS
3	C	1	GLU
4	D	11	LEU
4	D	28	LEU
4	D	91	VAL
4	D	97	LEU
4	D	108	LYS
4	D	110	ASN
4	D	129	LYS
4	D	147	LYS
4	D	153	ILE
4	D	176	ASN
4	D	184	ASN
4	D	190	ILE
5	E	2	GLN
5	E	33	LEU
5	E	82	LEU
5	E	96	THR
5	E	98	LEU
5	E	107	GLU
5	E	132	GLU
5	E	184	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (35)

such sidechains are listed below:

Mol	Chain	Res	Type
1	A	32	GLN
1	A	86	ASN
1	A	93	HIS
1	A	145	HIS
1	A	174	ASN
1	A	180	GLN
1	A	188	HIS
1	A	191	HIS
1	A	192	HIS
1	A	218	GLN
1	A	226	GLN
1	A	242	GLN
1	A	255	GLN
1	A	263	HIS
2	B	8	GLN
2	B	21	ASN
2	B	83	ASN
2	B	89	GLN
4	D	38	GLN
4	D	110	ASN
4	D	112	GLN
4	D	140	GLN
4	D	145	GLN
4	D	176	ASN
4	D	184	ASN
4	D	187	ASN
5	E	2	GLN
5	E	38	GLN
5	E	45	GLN
5	E	62	GLN
5	E	70	GLN
5	E	119	ASN
5	E	180	GLN
5	E	184	ASN
5	E	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

6 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	$\# Z > 2$
7	GOL	E	301	-	5,5,5	0.58	0	5,5,5	0.09	0
6	SO4	D	202	-	4,4,4	0.30	0	6,6,6	0.09	0
6	SO4	A	301	-	4,4,4	0.28	0	6,6,6	0.10	0
6	SO4	A	302	-	4,4,4	0.32	0	6,6,6	0.16	0
6	SO4	E	302	-	4,4,4	0.26	0	6,6,6	0.08	0
7	GOL	D	201	-	5,5,5	0.41	0	5,5,5	0.26	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
7	GOL	D	201	-	-	3/4/4/4	-
7	GOL	E	301	-	-	0/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	D	201	GOL	C1-C2-C3-O3
7	D	201	GOL	O2-C2-C3-O3
7	D	201	GOL	O1-C1-C2-O2

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	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	276/276 (100%)	0.53	40 (14%) 6 5	22, 55, 199, 264	0
2	B	100/100 (100%)	0.26	1 (1%) 79 78	28, 67, 139, 158	0
3	C	10/10 (100%)	-0.27	0 100 100	26, 32, 36, 44	0
4	D	197/197 (100%)	0.29	18 (9%) 15 12	20, 44, 135, 210	0
5	E	245/245 (100%)	-0.17	5 (2%) 65 62	12, 41, 82, 113	3 (1%)
All	All	828/828 (100%)	0.22	64 (7%) 19 17	12, 48, 154, 264	3 (0%)

All (64) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	248	VAL	7.5
4	D	191	ILE	5.5
1	A	197	HIS	5.4
1	A	249	VAL	5.2
1	A	257	TYR	4.9
4	D	182	CYS	4.8
1	A	193	ALA	4.6
1	A	247	VAL	4.6
4	D	196	PHE	4.5
1	A	194	VAL	4.1
4	D	197	PHE	4.1
2	B	99	MET	4.0
1	A	228	THR	3.9
4	D	190	ILE	3.8
4	D	163	SER	3.6
4	D	195	THR	3.6
1	A	246	ALA	3.6
1	A	199	ALA	3.5
1	A	276	PRO	3.5
1	A	250	PRO	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	200	THR	3.2
1	A	274	TRP	3.2
4	D	147	LYS	3.2
4	D	149	SER	3.0
1	A	245	ALA	3.0
1	A	251	SER	3.0
5	E	167[A]	HIS	2.9
1	A	219	ARG	2.9
1	A	258	THR	2.9
5	E	154[A]	HIS	2.9
1	A	201	LEU	2.8
1	A	191	HIS	2.8
4	D	164	MET	2.8
1	A	268	LYS	2.7
1	A	225	THR	2.7
4	D	166	PHE	2.7
1	A	189	MET	2.6
4	D	183	ALA	2.6
1	A	227	ASP	2.6
1	A	255	GLN	2.6
5	E	226	ASP	2.6
1	A	273	ARG	2.6
5	E	0	MET	2.6
1	A	269	PRO	2.6
1	A	192	HIS	2.5
1	A	259	CYS	2.4
1	A	204	TRP	2.4
1	A	271	THR	2.4
1	A	190	THR	2.4
1	A	253	GLN	2.4
4	D	145	GLN	2.3
1	A	226	GLN	2.3
1	A	220	ASP	2.3
4	D	138	ASP	2.3
1	A	196	ASP	2.3
4	D	126	SER	2.2
1	A	216	THR	2.2
5	E	222	GLU	2.2
4	D	177	LYS	2.2
4	D	181	ALA	2.1
1	A	272	LEU	2.1
4	D	189	SER	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	195	SER	2.0
1	A	261	VAL	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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
6	SO4	A	301	5/5	0.45	0.13	129,129,130,131	0
6	SO4	D	202	5/5	0.58	0.17	123,124,124,125	0
6	SO4	E	302	5/5	0.75	0.14	157,158,158,158	0
6	SO4	A	302	5/5	0.80	0.18	101,102,106,106	0
7	GOL	E	301	6/6	0.80	0.14	87,89,90,90	0
7	GOL	D	201	6/6	0.91	0.12	63,70,72,72	0

6.5 Other polymers [i](#)

There are no such residues in this entry.