Full wwPDB X-ray Structure Validation Report (i)

Jun 15, 2024 – 02:52 PM EDT

PDB ID	:	2H3X
Title	:	Crystal Structure of an Electron Transfer Complex Between Aromatic Amine Dehydrogenase and Azurin from Alcaligenes Faecalis (Form 3)
Authors	:	Sukumar, N.; Chen, Z.; Leys, D.; Scrutton, N.S.; Ferrati, D.; Merli, A.; Rossi, G.L.; Bellamy, H.D.; Chistoserdov, A.; Davidson, V.L.; Mathews, F.S.
Deposited on	:	2006-05-23
Resolution	:	2.50 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 (i) 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 (i)) were used in the production of this report:

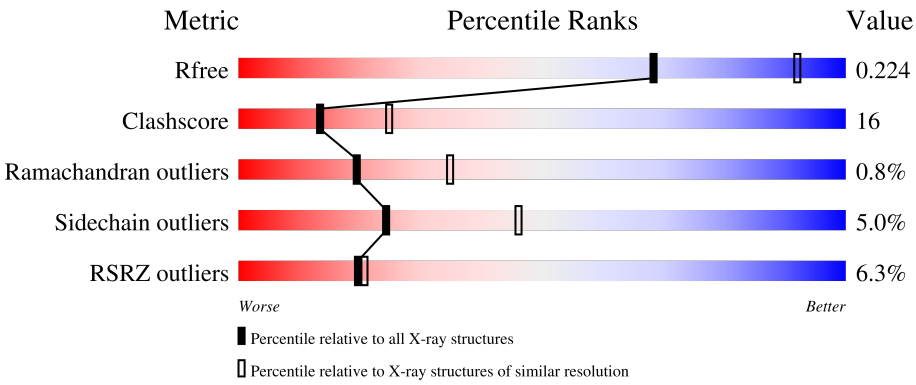
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

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 2.50 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}	130704	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

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 <=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	390	74% 16% • 8%
1	D	390	75% 16% • 8%
2	B	135	% 61% 25% 6% 7%
2	E	135	59% 27% 5% 8%
3	C	128	29% 56% 38% 6%

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Mol	Chain	Length	Quality of chain
3	F	128	30% 55% 38% 6%

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 10086 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 Aromatic Amine Dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	359	2805	1770	489	533	13	0	0	0
1	D	359	2805	1770	489	533	13	0	0	0

- Molecule 2 is a protein called Aromatic Amine Dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	125	956	582	168	191	15	0	0	0
2	E	124	951	579	167	190	15	0	0	0

- Molecule 3 is a protein called Azurin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	C	128	958	596	160	195	7	0	0	0
3	F	128	958	596	160	195	7	0	0	0

- Molecule 4 is COPPER (II) ION (three-letter code: CU) (formula: Cu).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Cu		
4	C	1	1	1	0	0
4	F	1	1	1	0	0

- Molecule 5 is water.

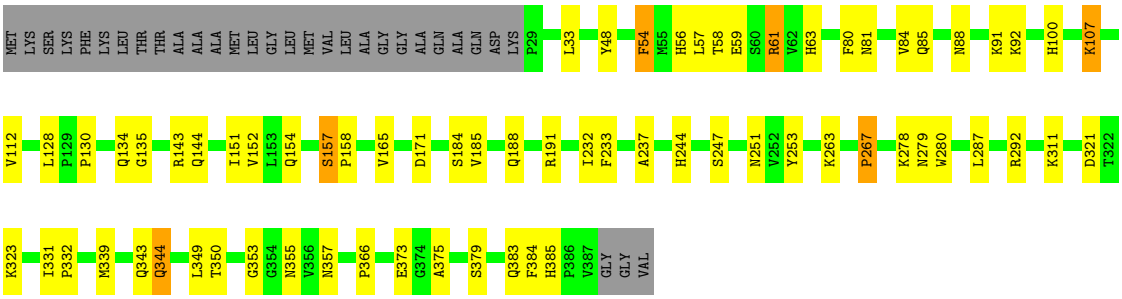
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
5	A	226	226	226	0	0
			Total	O		
5	B	76	76	76	0	0
			Total	O		
5	C	32	32	32	0	0
			Total	O		
5	D	228	228	228	0	0
			Total	O		
5	E	63	63	63	0	0
			Total	O		
5	F	26	26	26	0	0

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.

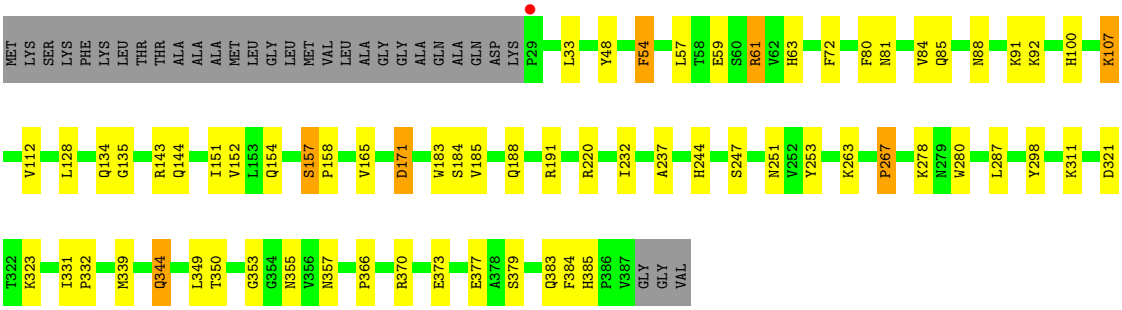
• Molecule 1: Aromatic Amine Dehydrogenase

Chain A:



• Molecule 1: Aromatic Amine Dehydrogenase

Chain D:

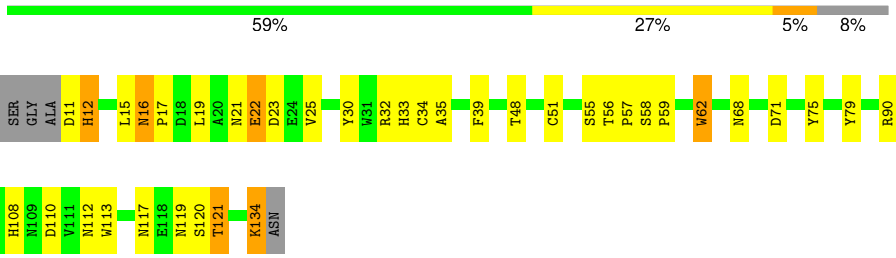


• Molecule 2: Aromatic Amine Dehydrogenase

Chain B:

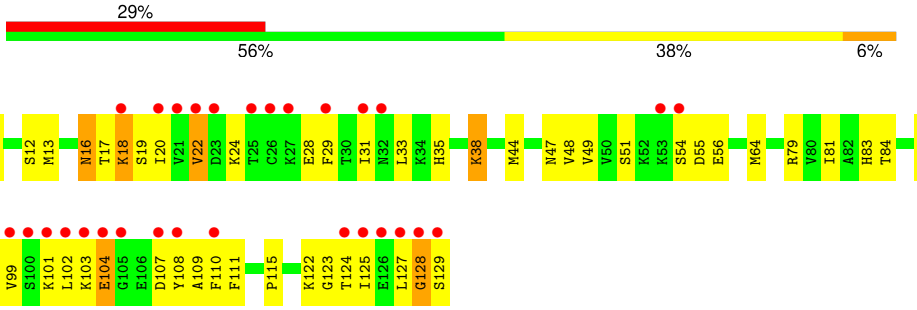
• Molecule 2: Aromatic Amine Dehydrogenase

Chain E:



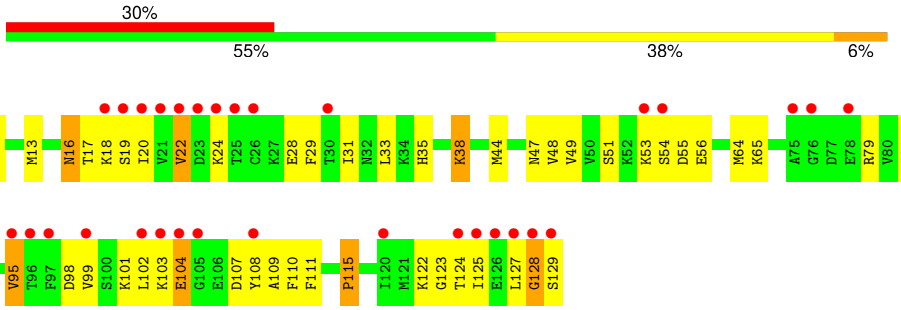
- Molecule 3: Azurin

Chain C:



- Molecule 3: Azurin

Chain F:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	64.95Å 94.76Å 211.48Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.00 – 2.50 47.79 – 2.50	Depositor EDS
% Data completeness (in resolution range)	71.5 (40.00-2.50) 71.9 (47.79-2.50)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.54 (at 2.51Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.183 , 0.230 0.175 , 0.224	Depositor DCC
R_{free} test set	3401 reflections (9.98%)	wwPDB-VP
Wilson B-factor (Å^2)	30.7	Xtriage
Anisotropy	0.320	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 57.5	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	10086	wwPDB-VP
Average B, all atoms (Å^2)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.86% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: CU, TRQ

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	0.32	0/2870	0.65	0/3890
1	D	0.33	0/2870	0.65	0/3890
2	B	0.35	0/964	0.67	0/1313
2	E	0.33	0/959	0.71	0/1306
3	C	0.31	0/973	0.62	0/1307
3	F	0.32	0/973	0.62	0/1307
All	All	0.33	0/9609	0.65	0/13013

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	2805	0	2749	57	0
1	D	2805	0	2749	64	0
2	B	956	0	842	42	0
2	E	951	0	837	50	0
3	C	958	0	932	57	0
3	F	958	0	932	57	0
4	C	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	F	1	0	0	0	0
5	A	226	0	0	2	0
5	B	76	0	0	0	0
5	C	32	0	0	2	0
5	D	228	0	0	2	0
5	E	63	0	0	0	0
5	F	26	0	0	1	0
All	All	10086	0	9041	297	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 16.

All (297) 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)
2:B:21:ASN:HD21	2:B:23:ASP:HB2	1.14	1.06
2:B:92:GLN:HE21	2:B:93:CYS:N	1.55	1.03
2:E:92:GLN:HE21	2:E:93:CYS:N	1.55	1.02
2:B:62:TRQ:HB2	2:B:113:TRP:NE1	1.74	1.00
2:E:62:TRQ:HB2	2:E:113:TRP:NE1	1.76	1.00
2:E:16:ASN:HD21	2:E:19:LEU:HG	1.29	0.94
2:B:62:TRQ:HB2	2:B:113:TRP:HE1	1.33	0.89
2:E:62:TRQ:HB2	2:E:113:TRP:HE1	1.35	0.89
2:E:21:ASN:HD21	2:E:23:ASP:HB2	1.37	0.87
1:A:112:VAL:HG23	1:A:128:LEU:HD11	1.60	0.83
2:E:92:GLN:NE2	2:E:93:CYS:N	2.27	0.82
1:D:112:VAL:HG23	1:D:128:LEU:HD11	1.61	0.82
2:B:92:GLN:NE2	2:B:93:CYS:N	2.27	0.82
2:B:119:ASN:OD1	2:B:121:THR:HG23	1.79	0.82
2:E:119:ASN:OD1	2:E:121:THR:HG23	1.79	0.81
2:B:21:ASN:ND2	2:B:23:ASP:HB2	1.96	0.81
1:D:54:PHE:CZ	2:E:92:GLN:HG3	2.16	0.80
2:E:16:ASN:ND2	2:E:19:LEU:HG	1.99	0.78
1:D:54:PHE:HZ	2:E:92:GLN:HG3	1.48	0.77
3:F:33:LEU:HG	3:F:87:ILE:HD11	1.67	0.76
2:B:92:GLN:HE21	2:B:92:GLN:C	1.89	0.76
2:E:92:GLN:HE21	2:E:92:GLN:C	1.90	0.75
2:B:11:ASP:O	2:B:12:HIS:HB2	1.88	0.73
3:C:33:LEU:HG	3:C:87:ILE:HD11	1.69	0.73
1:A:81:ASN:HD21	1:A:135:GLY:H	1.33	0.72
3:C:8:GLU:H	3:C:16:ASN:ND2	1.88	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:16:ASN:OD1	2:B:19:LEU:HD12	1.90	0.72
3:F:8:GLU:H	3:F:16:ASN:ND2	1.89	0.71
1:D:81:ASN:HD21	1:D:135:GLY:H	1.36	0.71
3:F:5:VAL:HG12	3:F:31:ILE:HG13	1.74	0.70
3:C:5:VAL:HG12	3:C:31:ILE:HG13	1.74	0.70
3:F:55:ASP:OD2	3:F:79:ARG:HD2	1.92	0.70
1:A:88:ASN:H	1:A:144:GLN:HE22	1.41	0.69
3:C:55:ASP:OD2	3:C:79:ARG:HD2	1.91	0.69
1:D:88:ASN:H	1:D:144:GLN:HE22	1.39	0.69
2:E:92:GLN:HE21	2:E:93:CYS:H	1.40	0.69
2:B:90:ARG:HG3	2:B:90:ARG:HH11	1.58	0.68
3:F:28:GLU:HA	3:F:98:ASP:HA	1.75	0.68
3:F:103:LYS:HB2	3:F:108:TYR:OH	1.93	0.68
2:E:21:ASN:ND2	2:E:23:ASP:HB2	2.09	0.68
2:E:90:ARG:HG3	2:E:90:ARG:HH11	1.57	0.68
3:C:28:GLU:HA	3:C:98:ASP:HA	1.74	0.68
3:C:103:LYS:HB2	3:C:108:TYR:OH	1.93	0.67
1:A:188:GLN:NE2	1:A:191:ARG:HH11	1.93	0.66
3:F:127:LEU:HD12	3:F:128:GLY:H	1.60	0.66
3:C:38:LYS:HG3	3:C:38:LYS:O	1.96	0.66
3:C:127:LEU:HD12	3:C:128:GLY:H	1.60	0.66
1:A:54:PHE:CZ	2:B:92:GLN:HG3	2.31	0.65
1:D:188:GLN:NE2	1:D:191:ARG:HH11	1.95	0.65
1:A:61:ARG:HH11	1:A:63:HIS:HE1	1.46	0.64
1:A:154:GLN:HE21	1:A:184:SER:H	1.45	0.64
1:A:54:PHE:HZ	2:B:92:GLN:HG3	1.62	0.64
1:D:61:ARG:HH11	1:D:63:HIS:HE1	1.46	0.63
1:D:143:ARG:HH22	1:D:383:GLN:NE2	1.96	0.63
3:F:38:LYS:O	3:F:38:LYS:HG3	1.96	0.63
2:B:92:GLN:HE21	2:B:93:CYS:H	1.42	0.63
2:E:39:PHE:CE2	2:E:92:GLN:HG2	2.34	0.63
1:D:311:LYS:HZ2	2:E:96:GLN:NE2	1.96	0.63
1:D:251:ASN:HB3	1:D:267:PRO:HB2	1.81	0.62
1:D:232:ILE:HA	1:D:247:SER:HA	1.82	0.62
2:B:59:PRO:O	3:C:12:SER:HA	2.00	0.62
1:D:154:GLN:HE21	1:D:184:SER:H	1.47	0.62
3:C:84:THR:HG22	3:C:95:VAL:HB	1.82	0.62
1:D:154:GLN:NE2	1:D:184:SER:H	1.98	0.62
1:A:143:ARG:HH22	1:A:383:GLN:NE2	1.98	0.61
2:B:39:PHE:CE2	2:B:92:GLN:HG2	2.35	0.61
3:F:84:THR:HG22	3:F:95:VAL:HB	1.82	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:154:GLN:NE2	1:A:184:SER:H	1.99	0.61
3:C:33:LEU:HD22	3:C:48:VAL:HG23	1.83	0.61
3:F:33:LEU:HD22	3:F:48:VAL:HG23	1.83	0.60
1:A:232:ILE:HA	1:A:247:SER:HA	1.82	0.60
1:A:134:GLN:HE22	2:B:120:SER:HB2	1.67	0.60
3:C:17:THR:HG22	3:C:19:SER:H	1.67	0.59
3:F:17:THR:HG22	3:F:19:SER:H	1.67	0.59
3:C:5:VAL:HG11	3:C:20:ILE:HD13	1.84	0.59
1:A:251:ASN:HB3	1:A:267:PRO:HB2	1.82	0.59
1:D:220:ARG:HH21	3:F:64:MET:HE1	1.67	0.59
2:E:92:GLN:HE21	2:E:92:GLN:CA	2.15	0.59
3:F:44:MET:O	3:F:44:MET:HG3	2.02	0.59
2:B:92:GLN:HE21	2:B:92:GLN:CA	2.15	0.58
3:F:53:LYS:HG3	5:F:343:HOH:O	2.03	0.58
2:B:30:TYR:HB3	2:B:33:HIS:CD2	2.38	0.58
3:F:51:SER:HA	3:F:81:ILE:HD13	1.85	0.58
3:F:5:VAL:HG11	3:F:20:ILE:HD13	1.85	0.58
3:C:44:MET:O	3:C:44:MET:HG3	2.04	0.57
3:F:102:LEU:HD22	3:F:108:TYR:CD2	2.39	0.57
1:D:311:LYS:NZ	2:E:96:GLN:NE2	2.52	0.57
2:E:30:TYR:HB3	2:E:33:HIS:CD2	2.40	0.57
3:C:102:LEU:HD22	3:C:108:TYR:CD2	2.39	0.57
2:E:11:ASP:O	2:E:12:HIS:HB2	2.04	0.57
1:D:88:ASN:H	1:D:144:GLN:NE2	2.03	0.56
3:F:104:GLU:OE1	3:F:104:GLU:N	2.38	0.56
3:C:104:GLU:OE1	3:C:104:GLU:N	2.38	0.56
1:A:84:VAL:HG22	1:A:85:GLN:N	2.21	0.56
3:C:51:SER:HA	3:C:81:ILE:HD13	1.88	0.56
3:F:110:PHE:CZ	3:F:123:GLY:HA3	2.41	0.55
2:E:58:SER:HA	2:E:117:ASN:HD21	1.72	0.55
1:A:237:ALA:HB3	1:A:244:HIS:HB2	1.88	0.55
2:E:62:TRQ:CB	2:E:113:TRP:NE1	2.63	0.55
2:B:108:HIS:CE1	2:B:110:ASP:HB2	2.42	0.55
2:E:108:HIS:CE1	2:E:110:ASP:HB2	2.41	0.55
1:D:48:TYR:OH	1:D:385:HIS:HD2	1.90	0.55
1:A:56:HIS:NE2	1:D:377:GLU:OE2	2.33	0.54
2:B:58:SER:HA	2:B:117:ASN:HD21	1.72	0.54
2:E:15:LEU:HD21	2:E:75:TYR:CZ	2.42	0.54
1:D:84:VAL:HG22	1:D:85:GLN:N	2.22	0.54
2:B:35:ALA:HB3	2:B:79:TYR:CE1	2.41	0.54
2:E:35:ALA:HB3	2:E:79:TYR:CE1	2.43	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:237:ALA:HB3	1:D:244:HIS:HB2	1.88	0.54
3:C:110:PHE:CZ	3:C:123:GLY:HA3	2.42	0.54
1:D:143:ARG:HH22	1:D:383:GLN:HE22	1.56	0.54
3:F:128:GLY:O	3:F:129:SER:HB2	2.07	0.54
1:D:134:GLN:HE22	2:E:120:SER:HB2	1.73	0.54
1:D:344:GLN:HB2	5:D:527:HOH:O	2.07	0.54
3:F:5:VAL:HG21	3:F:20:ILE:HG23	1.90	0.54
3:C:5:VAL:HG21	3:C:20:ILE:HG23	1.89	0.53
3:C:128:GLY:O	3:C:129:SER:HB2	2.07	0.53
3:C:22:VAL:HG12	3:C:29:PHE:CE1	2.43	0.53
3:C:127:LEU:HD12	3:C:128:GLY:N	2.23	0.53
2:E:106:PHE:CZ	3:F:13:MET:HE1	2.44	0.53
3:F:22:VAL:HG12	3:F:29:PHE:CE1	2.43	0.53
1:A:61:ARG:HH11	1:A:63:HIS:CE1	2.27	0.53
3:C:24:LYS:HG3	3:C:128:GLY:HA3	1.90	0.53
1:D:339:MET:HA	1:D:349:LEU:O	2.09	0.53
2:B:62:TRQ:CB	2:B:113:TRP:NE1	2.61	0.53
3:F:24:LYS:HG3	3:F:128:GLY:HA3	1.91	0.53
3:F:31:ILE:HG23	3:F:31:ILE:O	2.09	0.53
1:D:134:GLN:NE2	2:E:112:ASN:HD22	2.07	0.52
3:F:127:LEU:HD12	3:F:128:GLY:N	2.23	0.52
2:E:16:ASN:OD1	2:E:19:LEU:HD12	2.09	0.52
1:A:339:MET:HA	1:A:349:LEU:O	2.10	0.52
2:E:22:GLU:HA	2:E:25:VAL:HG22	1.90	0.52
2:E:90:ARG:HG3	2:E:90:ARG:NH1	2.23	0.52
2:E:92:GLN:NE2	2:E:93:CYS:H	2.03	0.52
2:B:90:ARG:HG3	2:B:90:ARG:NH1	2.24	0.52
3:C:31:ILE:HG23	3:C:31:ILE:O	2.09	0.52
2:B:106:PHE:CZ	3:C:13:MET:HE1	2.45	0.51
1:A:88:ASN:H	1:A:144:GLN:NE2	2.05	0.51
1:A:355:ASN:OD1	1:A:373:GLU:HA	2.10	0.51
1:A:33:LEU:CD2	2:E:57:PRO:HB3	2.41	0.51
1:A:81:ASN:HD21	1:A:135:GLY:N	2.05	0.51
3:C:5:VAL:CG1	3:C:31:ILE:HG13	2.40	0.51
3:C:8:GLU:H	3:C:16:ASN:HD21	1.56	0.51
1:A:278:LYS:HG2	5:A:545:HOH:O	2.10	0.51
1:A:48:TYR:OH	1:A:385:HIS:HD2	1.94	0.51
1:A:353:GLY:HA2	1:A:379:SER:O	2.11	0.51
3:C:24:LYS:HG2	3:C:127:LEU:HD11	1.93	0.51
1:A:152:VAL:HG12	1:A:185:VAL:HG21	1.93	0.51
1:D:355:ASN:OD1	1:D:373:GLU:HA	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:350:THR:OG1	1:A:357:ASN:HB2	2.11	0.50
3:F:24:LYS:HG2	3:F:127:LEU:HD11	1.94	0.50
1:A:331:ILE:HD13	1:A:366:PRO:HB2	1.94	0.50
1:D:350:THR:OG1	1:D:357:ASN:HB2	2.12	0.50
1:A:57:LEU:O	1:A:80:PHE:HA	2.12	0.50
3:C:7:ILE:HG23	3:C:33:LEU:HD13	1.92	0.49
3:F:5:VAL:CG1	3:F:31:ILE:HG13	2.39	0.49
1:D:321:ASP:OD1	1:D:323:LYS:HB2	2.13	0.49
3:F:49:VAL:HG22	3:F:83:HIS:HB2	1.94	0.49
1:A:143:ARG:HH22	1:A:383:GLN:HE22	1.60	0.49
1:D:61:ARG:HH11	1:D:63:HIS:CE1	2.27	0.49
1:D:91:LYS:O	1:D:92:LYS:HD3	2.12	0.49
1:D:220:ARG:HH21	3:F:64:MET:CE	2.25	0.49
1:D:157:SER:N	1:D:158:PRO:HA	2.27	0.49
1:A:157:SER:N	1:A:158:PRO:HA	2.27	0.49
2:B:51:CYS:HB3	2:B:55:SER:OG	2.13	0.49
2:E:51:CYS:HB3	2:E:55:SER:OG	2.13	0.49
3:C:16:ASN:N	3:C:16:ASN:HD22	2.11	0.48
3:C:99:VAL:C	3:C:101:LYS:H	2.16	0.48
3:F:109:ALA:HB1	3:F:123:GLY:O	2.12	0.48
1:A:134:GLN:NE2	2:B:112:ASN:HD22	2.11	0.48
2:B:57:PRO:HB3	1:D:33:LEU:CD2	2.44	0.48
1:D:152:VAL:HG12	1:D:185:VAL:HG21	1.94	0.48
3:F:7:ILE:HG23	3:F:33:LEU:HD13	1.94	0.48
3:C:49:VAL:HG22	3:C:83:HIS:HB2	1.96	0.48
3:C:109:ALA:HB1	3:C:123:GLY:O	2.14	0.48
1:D:331:ILE:HD13	1:D:366:PRO:HB2	1.96	0.48
1:A:33:LEU:HD23	2:E:57:PRO:HB3	1.95	0.48
1:A:233:PHE:HB3	5:A:451:HOH:O	2.14	0.48
3:F:104:GLU:CD	3:F:104:GLU:H	2.08	0.48
1:D:57:LEU:O	1:D:80:PHE:HA	2.14	0.47
3:F:16:ASN:N	3:F:16:ASN:HD22	2.12	0.47
3:C:7:ILE:O	3:C:33:LEU:HD12	2.13	0.47
1:A:253:TYR:CZ	1:A:267:PRO:HB3	2.50	0.47
1:D:353:GLY:HA2	1:D:379:SER:O	2.14	0.47
2:E:62:TRQ:HB2	2:E:113:TRP:CE2	2.49	0.47
1:A:349:LEU:HG	1:A:384:PHE:CE1	2.50	0.47
1:D:349:LEU:HG	1:D:384:PHE:CE1	2.50	0.47
3:F:13:MET:HE2	3:F:13:MET:HB2	1.75	0.47
3:F:122:LYS:C	3:F:122:LYS:HD3	2.36	0.47
3:C:122:LYS:HD3	3:C:122:LYS:C	2.35	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:7:ILE:O	3:F:33:LEU:HD12	2.14	0.47
1:A:321:ASP:OD1	1:A:323:LYS:HB2	2.15	0.46
2:E:134:LYS:HE3	2:E:134:LYS:HB3	1.74	0.46
3:C:104:GLU:CD	3:C:104:GLU:H	2.08	0.46
1:D:54:PHE:CE1	2:E:92:GLN:CG	2.98	0.46
1:A:100:HIS:HA	1:A:107:LYS:O	2.15	0.46
2:B:86:THR:HG21	1:D:59:GLU:HG3	1.97	0.46
1:D:81:ASN:HD21	1:D:135:GLY:N	2.08	0.46
1:A:91:LYS:O	1:A:92:LYS:HD3	2.15	0.46
2:B:32:ARG:HD2	2:B:48:THR:HG21	1.96	0.46
3:C:47:ASN:C	3:C:47:ASN:OD1	2.54	0.46
1:D:220:ARG:NH2	3:F:64:MET:CE	2.78	0.46
2:E:32:ARG:HD2	2:E:48:THR:HG21	1.98	0.46
3:F:99:VAL:C	3:F:101:LYS:H	2.18	0.46
2:B:16:ASN:OD1	2:B:19:LEU:HB2	2.16	0.46
3:F:47:ASN:OD1	3:F:47:ASN:C	2.54	0.46
3:C:13:MET:HB2	3:C:13:MET:HE2	1.64	0.46
1:D:253:TYR:CZ	1:D:267:PRO:HB3	2.51	0.46
3:C:18:LYS:HE2	5:C:145:HOH:O	2.15	0.45
1:A:349:LEU:HG	1:A:384:PHE:CZ	2.51	0.45
1:A:344:GLN:HE21	1:A:344:GLN:HB3	1.58	0.45
3:C:5:VAL:HG21	3:C:20:ILE:CG2	2.47	0.45
3:C:64:MET:HB2	5:C:143:HOH:O	2.15	0.45
1:D:349:LEU:HG	1:D:384:PHE:CZ	2.51	0.45
2:E:17:PRO:HD3	2:E:32:ARG:NH2	2.31	0.45
3:C:102:LEU:HD22	3:C:108:TYR:CG	2.52	0.45
1:A:59:GLU:O	1:A:59:GLU:HG2	2.17	0.45
3:F:35:HIS:ND1	3:F:89:GLY:HA2	2.32	0.45
3:F:8:GLU:H	3:F:16:ASN:HD21	1.59	0.45
3:C:87:ILE:HB	3:C:91:GLU:HB2	1.99	0.45
1:D:100:HIS:HA	1:D:107:LYS:O	2.16	0.45
1:A:157:SER:N	1:A:158:PRO:CA	2.80	0.44
3:F:81:ILE:HD12	3:F:81:ILE:N	2.32	0.44
3:F:102:LEU:HD22	3:F:108:TYR:CG	2.52	0.44
3:F:125:ILE:HG23	3:F:125:ILE:O	2.17	0.44
2:B:60:ILE:HG22	3:C:44:MET:HE3	1.98	0.44
2:B:62:TRQ:HB2	2:B:113:TRP:CE2	2.48	0.44
3:C:81:ILE:HD12	3:C:81:ILE:N	2.32	0.44
3:F:22:VAL:HG22	3:F:127:LEU:HA	1.99	0.44
3:F:87:ILE:HB	3:F:91:GLU:HB2	2.00	0.44
2:B:92:GLN:NE2	2:B:93:CYS:H	2.05	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:108:HIS:NE2	2:E:110:ASP:HB2	2.33	0.44
1:D:183:TRP:CD1	2:E:102:GLY:HA3	2.53	0.44
1:A:151:ILE:HB	1:A:165:VAL:HB	2.00	0.44
2:E:59:PRO:HD2	2:E:117:ASN:HD21	1.83	0.44
3:C:56:GLU:HG3	3:C:111:PHE:CE1	2.52	0.44
1:D:157:SER:N	1:D:158:PRO:CA	2.81	0.44
2:B:108:HIS:NE2	2:B:110:ASP:HB2	2.33	0.43
3:F:5:VAL:HG21	3:F:20:ILE:CG2	2.48	0.43
2:B:16:ASN:HD21	2:B:19:LEU:HG	1.82	0.43
2:B:59:PRO:HD2	2:B:117:ASN:HD21	1.83	0.43
3:C:35:HIS:ND1	3:C:89:GLY:HA2	2.32	0.43
1:A:63:HIS:HD2	1:A:375:ALA:O	2.02	0.43
3:F:56:GLU:HG3	3:F:111:PHE:CE1	2.52	0.43
1:A:84:VAL:CG2	1:A:85:GLN:N	2.82	0.43
1:D:61:ARG:NH1	1:D:63:HIS:HE1	2.15	0.43
1:D:151:ILE:HB	1:D:165:VAL:HB	2.01	0.43
1:D:370:ARG:HD3	5:D:392:HOH:O	2.18	0.43
3:C:22:VAL:HG22	3:C:127:LEU:HA	2.00	0.43
3:C:122:LYS:HD3	3:C:123:GLY:N	2.34	0.43
3:C:125:ILE:O	3:C:125:ILE:HG23	2.17	0.43
2:E:11:ASP:O	2:E:12:HIS:CB	2.66	0.43
3:F:22:VAL:O	3:F:22:VAL:CG2	2.67	0.43
1:D:154:GLN:HE21	1:D:184:SER:N	2.16	0.42
3:C:19:SER:HA	3:C:124:THR:O	2.18	0.42
2:B:44:CYS:HA	1:D:72:PHE:CD2	2.54	0.42
1:A:61:ARG:HD2	1:A:61:ARG:N	2.35	0.42
1:A:188:GLN:HE22	1:A:191:ARG:HH11	1.66	0.42
3:C:22:VAL:O	3:C:22:VAL:CG2	2.67	0.42
1:D:54:PHE:CE1	2:E:92:GLN:HG3	2.53	0.42
3:C:109:ALA:HB1	3:C:122:LYS:HE2	2.02	0.42
1:D:59:GLU:O	1:D:59:GLU:HG2	2.18	0.42
1:D:61:ARG:N	1:D:61:ARG:HD2	2.34	0.42
3:F:19:SER:HA	3:F:124:THR:O	2.20	0.42
3:F:35:HIS:HB3	3:F:88:GLY:O	2.19	0.42
1:A:278:LYS:HB2	1:A:280:TRP:CD1	2.54	0.42
2:E:16:ASN:HA	2:E:17:PRO:HD3	1.82	0.41
2:E:56:THR:HA	2:E:57:PRO:HD3	1.93	0.41
3:C:35:HIS:HB3	3:C:88:GLY:O	2.19	0.41
3:F:109:ALA:HA	3:F:124:THR:HA	2.02	0.41
1:A:61:ARG:NH1	1:A:63:HIS:HE1	2.15	0.41
3:C:102:LEU:HD23	3:C:102:LEU:HA	1.92	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:109:ALA:HB1	3:F:122:LYS:HE2	2.01	0.41
1:A:311:LYS:NZ	2:B:96:GLN:NE2	2.69	0.41
1:A:292:ARG:H	1:A:343:GLN:HE22	1.69	0.41
1:D:54:PHE:CZ	2:E:92:GLN:CG	2.97	0.41
1:D:278:LYS:HB2	1:D:280:TRP:CD1	2.55	0.41
2:E:68:ASN:HB3	2:E:71:ASP:OD1	2.20	0.41
3:F:122:LYS:HD3	3:F:123:GLY:N	2.35	0.41
2:B:16:ASN:HA	2:B:17:PRO:HD3	1.75	0.41
1:A:278:LYS:O	1:A:279:ASN:HB2	2.21	0.41
2:B:99:GLU:OE2	2:B:108:HIS:HD2	2.04	0.41
3:C:99:VAL:C	3:C:101:LYS:N	2.74	0.41
2:E:16:ASN:OD1	2:E:19:LEU:CD1	2.69	0.41
1:A:188:GLN:NE2	1:A:191:ARG:HD2	2.36	0.41
1:D:220:ARG:CD	3:F:65:LYS:HD3	2.51	0.41
1:D:331:ILE:HB	1:D:332:PRO:CD	2.51	0.41
1:A:188:GLN:NE2	1:A:191:ARG:NH1	2.65	0.40
3:C:98:ASP:OD1	3:C:98:ASP:N	2.48	0.40
1:D:165:VAL:HA	1:D:171:ASP:O	2.21	0.40
1:D:298:TYR:CD1	1:D:298:TYR:N	2.89	0.40
1:A:58:THR:OG1	1:D:59:GLU:OE1	2.29	0.40
1:D:84:VAL:CG2	1:D:85:GLN:N	2.84	0.40
1:A:331:ILE:HB	1:A:332:PRO:CD	2.52	0.40
2:B:68:ASN:HB3	2:B:71:ASP:OD1	2.21	0.40
1:D:188:GLN:NE2	1:D:191:ARG:HD2	2.37	0.40
3:F:99:VAL:C	3:F:101:LYS:N	2.75	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	357/390 (92%)	331 (93%)	24 (7%)	2 (1%)	25 43

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	357/390 (92%)	334 (94%)	22 (6%)	1 (0%)	41	61
2	B	122/135 (90%)	112 (92%)	8 (7%)	2 (2%)	9	17
2	E	121/135 (90%)	113 (93%)	7 (6%)	1 (1%)	19	35
3	C	126/128 (98%)	112 (89%)	12 (10%)	2 (2%)	9	17
3	F	126/128 (98%)	112 (89%)	12 (10%)	2 (2%)	9	17
All	All	1209/1306 (93%)	1114 (92%)	85 (7%)	10 (1%)	19	35

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	11	ASP
2	B	12	HIS
2	E	12	HIS
3	C	128	GLY
3	F	128	GLY
3	C	115	PRO
3	F	115	PRO
1	D	157	SER
1	A	157	SER
1	A	130	PRO

5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	305/325 (94%)	297 (97%)	8 (3%)	46	72
1	D	305/325 (94%)	297 (97%)	8 (3%)	46	72
2	B	108/112 (96%)	100 (93%)	8 (7%)	13	27
2	E	108/112 (96%)	101 (94%)	7 (6%)	17	33
3	C	106/106 (100%)	96 (91%)	10 (9%)	8	17
3	F	106/106 (100%)	95 (90%)	11 (10%)	7	13
All	All	1038/1086 (96%)	986 (95%)	52 (5%)	24	46

All (52) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	54	PHE
1	A	61	ARG
1	A	107	LYS
1	A	171	ASP
1	A	263	LYS
1	A	267	PRO
1	A	287	LEU
1	A	344	GLN
2	B	16	ASN
2	B	21	ASN
2	B	23	ASP
2	B	34	CYS
2	B	92	GLN
2	B	98	ARG
2	B	121	THR
2	B	134	LYS
3	C	4	ASP
3	C	5	VAL
3	C	16	ASN
3	C	18	LYS
3	C	22	VAL
3	C	38	LYS
3	C	54	SER
3	C	95	VAL
3	C	104	GLU
3	C	107	ASP
1	D	54	PHE
1	D	61	ARG
1	D	107	LYS
1	D	171	ASP
1	D	263	LYS
1	D	267	PRO
1	D	287	LEU
1	D	344	GLN
2	E	16	ASN
2	E	22	GLU
2	E	34	CYS
2	E	92	GLN
2	E	98	ARG
2	E	121	THR
2	E	134	LYS
3	F	4	ASP

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Mol	Chain	Res	Type
3	F	5	VAL
3	F	16	ASN
3	F	18	LYS
3	F	22	VAL
3	F	38	LYS
3	F	54	SER
3	F	95	VAL
3	F	104	GLU
3	F	107	ASP
3	F	115	PRO

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (43) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	43	GLN
1	A	63	HIS
1	A	81	ASN
1	A	85	GLN
1	A	100	HIS
1	A	134	GLN
1	A	144	GLN
1	A	154	GLN
1	A	188	GLN
1	A	223	GLN
1	A	343	GLN
1	A	344	GLN
1	A	381	GLN
1	A	383	GLN
1	A	385	HIS
2	B	21	ASN
2	B	92	GLN
2	B	96	GLN
2	B	108	HIS
2	B	117	ASN
3	C	16	ASN
1	D	43	GLN
1	D	63	HIS
1	D	81	ASN
1	D	85	GLN
1	D	100	HIS
1	D	134	GLN
1	D	144	GLN

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Mol	Chain	Res	Type
1	D	154	GLN
1	D	188	GLN
1	D	223	GLN
1	D	343	GLN
1	D	344	GLN
1	D	381	GLN
1	D	383	GLN
1	D	385	HIS
2	E	21	ASN
2	E	92	GLN
2	E	94	ASN
2	E	96	GLN
2	E	108	HIS
2	E	117	ASN
3	F	16	ASN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains (i)

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

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
2	TRQ	E	62	2	13,17,18	4.54	3 (23%)	13,24,26	2.88	6 (46%)
2	TRQ	B	62	2	13,17,18	4.48	3 (23%)	13,24,26	3.40	6 (46%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	TRQ	E	62	2	-	0/4/19/21	0/2/2/2
2	TRQ	B	62	2	-	0/4/19/21	0/2/2/2

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	62	TRQ	CH2-CZ2	-13.70	1.37	1.53
2	B	62	TRQ	CH2-CZ2	-13.41	1.37	1.53
2	E	62	TRQ	CE2-CZ2	-8.03	1.40	1.50
2	B	62	TRQ	CE2-CZ2	-7.93	1.40	1.50
2	B	62	TRQ	CZ3-CE3	3.33	1.40	1.34
2	E	62	TRQ	CZ3-CE3	2.95	1.39	1.34

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	62	TRQ	O7-CZ2-CH2	6.09	125.76	119.01
2	B	62	TRQ	CZ2-CE2-NE1	5.61	128.90	119.94
2	E	62	TRQ	CZ2-CE2-NE1	5.60	128.89	119.94
2	B	62	TRQ	O6-CH2-CZ2	5.37	122.43	118.45
2	B	62	TRQ	O7-CZ2-CE2	-4.71	116.38	121.76
2	E	62	TRQ	O7-CZ2-CH2	4.68	124.20	119.01
2	E	62	TRQ	O6-CH2-CZ2	3.98	121.40	118.45
2	E	62	TRQ	O7-CZ2-CE2	-3.29	118.00	121.76
2	B	62	TRQ	O6-CH2-CZ3	-3.00	116.18	121.58
2	E	62	TRQ	CD1-CG-CD2	2.97	106.49	104.79
2	B	62	TRQ	CD1-CG-CD2	2.91	106.46	104.79
2	E	62	TRQ	O6-CH2-CZ3	-2.44	117.20	121.58

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	62	TRQ	4	0
2	B	62	TRQ	4	0

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

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 (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 (i)

6.1 Protein, DNA and RNA chains (i)

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	359/390 (92%)	-0.35	0 100 100	14, 26, 44, 69	0
1	D	359/390 (92%)	-0.43	1 (0%) 94 94	13, 26, 42, 67	0
2	B	124/135 (91%)	-0.31	1 (0%) 86 87	12, 25, 54, 76	0
2	E	123/135 (91%)	-0.29	0 100 100	16, 26, 62, 76	0
3	C	128/128 (100%)	1.45	37 (28%) 0 0	31, 56, 93, 100	0
3	F	128/128 (100%)	1.52	38 (29%) 0 0	38, 61, 93, 100	0
All	All	1221/1306 (93%)	0.02	77 (6%) 20 21	12, 28, 76, 100	0

All (77) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	C	127	LEU	7.3
3	F	129	SER	6.9
3	C	107	ASP	6.6
3	F	128	GLY	6.1
3	C	21	VAL	6.0
3	C	128	GLY	5.9
3	F	22	VAL	5.5
3	C	20	ILE	5.5
3	C	5	VAL	5.4
3	F	103	LYS	5.1
3	C	105	GLY	5.1
3	C	125	ILE	4.9
3	C	22	VAL	4.8
3	C	129	SER	4.7
3	F	105	GLY	4.7
3	F	95	VAL	4.5
3	C	2	ALA	4.4
3	F	127	LEU	4.4
3	C	31	ILE	4.3

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Mol	Chain	Res	Type	RSRZ
3	F	21	VAL	4.3
3	C	98	ASP	4.2
3	F	97	PHE	4.1
3	C	104	GLU	4.1
3	F	23	ASP	4.1
3	C	29	PHE	3.9
3	F	18	LYS	3.9
3	C	108	TYR	3.9
3	C	54	SER	3.8
3	C	102	LEU	3.7
3	F	20	ILE	3.7
2	B	134	LYS	3.7
3	F	75	ALA	3.6
3	C	3	CYS	3.5
3	C	97	PHE	3.5
3	C	126	GLU	3.4
3	C	99	VAL	3.3
3	F	99	VAL	3.3
3	F	92	THR	3.3
3	F	7	ILE	3.3
3	F	102	LEU	3.2
3	C	110	PHE	3.2
3	C	27	LYS	3.1
3	F	96	THR	3.1
3	F	19	SER	3.1
3	C	103	LYS	3.1
3	F	124	THR	3.1
3	F	104	GLU	3.0
3	C	23	ASP	3.0
3	C	4	ASP	3.0
3	F	54	SER	3.0
3	C	26	CYS	2.9
3	F	76	GLY	2.9
3	C	18	LYS	2.9
3	F	5	VAL	2.8
3	C	25	THR	2.8
3	C	124	THR	2.7
3	F	2	ALA	2.7
1	D	29	PRO	2.7
3	C	101	LYS	2.7
3	F	53	LYS	2.6
3	C	92	THR	2.6

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Mol	Chain	Res	Type	RSRZ
3	F	26	CYS	2.6
3	C	53	LYS	2.6
3	F	126	GLU	2.6
3	F	24	LYS	2.6
3	F	108	TYR	2.5
3	F	90	GLY	2.5
3	F	25	THR	2.5
3	F	30	THR	2.4
3	C	100	SER	2.3
3	F	120	ILE	2.3
3	F	78	GLU	2.3
3	C	32	ASN	2.2
3	F	3	CYS	2.2
3	F	87	ILE	2.2
3	F	125	ILE	2.2
3	C	7	ILE	2.1

6.2 Non-standard residues in protein, DNA, RNA chains (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(Å^2)	Q<0.9
2	TRQ	B	62	16/17	0.96	0.13	17,22,28,29	0
2	TRQ	E	62	16/17	0.98	0.12	15,23,26,28	0

6.3 Carbohydrates (i)

There are no monosaccharides 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(Å^2)	Q<0.9
4	CU	C	130	1/1	0.99	0.06	49,49,49,49	0
4	CU	F	130	1/1	0.99	0.05	50,50,50,50	0

6.5 Other polymers (i)

There are no such residues in this entry.