Full wwPDB X-ray Structure Validation Report (i)

Jun 16, 2024 – 10:11 AM EDT

PDB ID	:	2FPI
Title	:	Crystal structure of pig GTP-specific succinyl-CoA synthetase from polyethylene glycol
Authors	:	Fraser, M.E.
Deposited on	:	2006-01-16
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 (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:

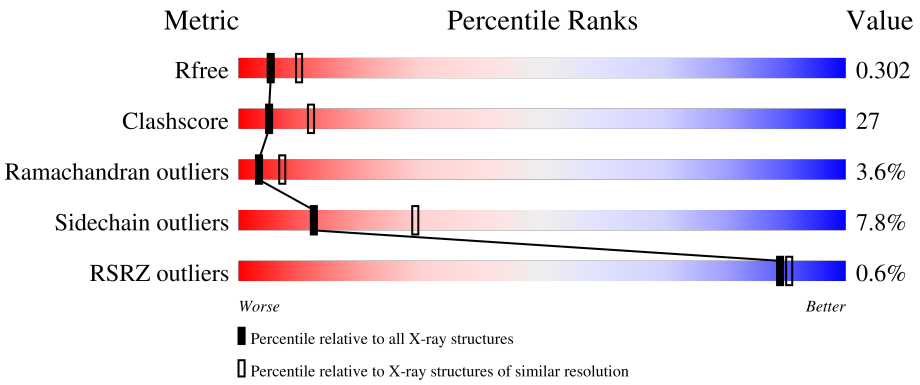
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, 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 2.70 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}	130704	2808 (2.70-2.70)
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)
RSRZ outliers	127900	2737 (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 <=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	305	64% 31% 6%
2	B	395	% 46% 46% 7% •

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 5299 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 Succinyl-CoA ligase [GDP-forming] alpha-chain, mitochondrial.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	P	S			
1	A	305	2255	1418	400	425	1	11	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	2	SER	-	CLONING ARTIFACT	UNP O19069-2
A	259	NEP	HIS	MODIFIED RESIDUE	UNP O19069-2

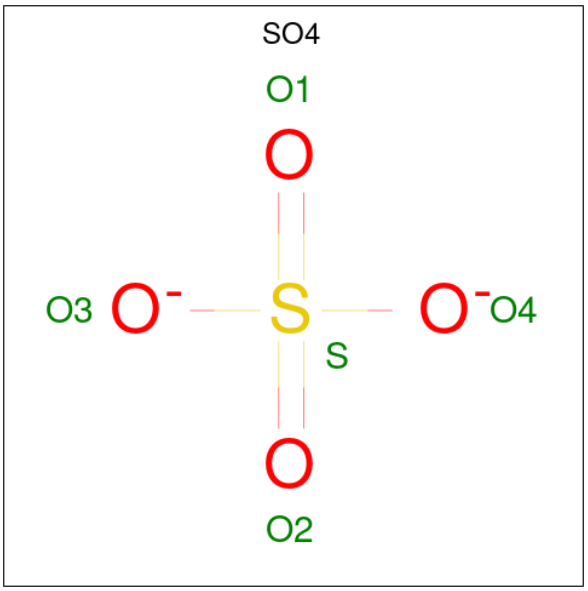
- Molecule 2 is a protein called Succinyl-CoA ligase [GDP-forming] beta-chain, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	393	2970	1876	502	578	14	0	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	1	MET	-	INITIATING METHIONINE	UNP P53590

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



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

- Molecule 4 is water.

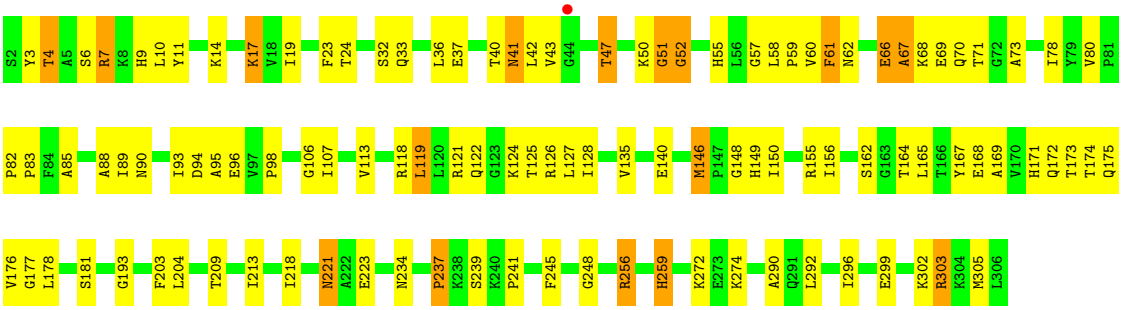
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	39	Total O 39 39	0	0
4	B	25	Total O 25 25	0	0

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.

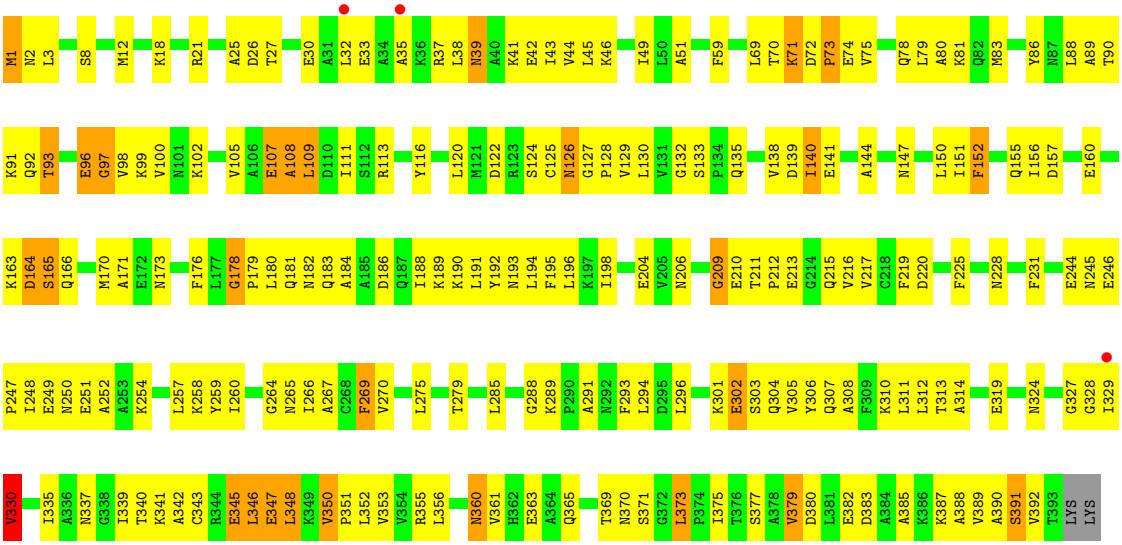
- Molecule 1: Succinyl-CoA ligase [GDP-forming] alpha-chain, mitochondrial

Chain A:



- Molecule 2: Succinyl-CoA ligase [GDP-forming] beta-chain, mitochondrial

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, α, β, γ	135.79Å 135.79Å 77.15Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	100.00 – 2.70 34.95 – 2.68	Depositor EDS
% Data completeness (in resolution range)	99.8 (100.00-2.70) 99.2 (34.95-2.68)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	5.88 (at 2.68Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.215 , 0.310 0.208 , 0.302	Depositor DCC
R_{free} test set	1143 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å^2)	50.0	Xtriage
Anisotropy	0.021	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.30 , 47.1	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.034 for h,-h-k,-l	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	5299	wwPDB-VP
Average B, all atoms (Å^2)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.83% 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: SO4, NEP

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	0.77	0/2281	0.91	1/3081 (0.0%)
2	B	0.73	0/3007	0.90	1/4059 (0.0%)
All	All	0.75	0/5288	0.91	2/7140 (0.0%)

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
2	B	0	1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	7	ARG	NE-CZ-NH1	-5.88	117.36	120.30
2	B	209	GLY	N-CA-C	5.08	125.80	113.10

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	86	TYR	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	2255	0	2295	95	0
2	B	2970	0	3023	199	0
3	A	5	0	0	0	0
3	B	5	0	0	0	0
4	A	39	0	0	0	0
4	B	25	0	0	2	0
All	All	5299	0	5318	287	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 27.

All (287) 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:70:THR:HG21	2:B:75:VAL:HB	1.47	0.97
1:A:89:ILE:HG22	1:A:93:ILE:HD11	1.53	0.89
2:B:113:ARG:HG3	2:B:113:ARG:HH11	1.37	0.88
2:B:180:LEU:HD13	2:B:210:GLU:OE1	1.72	0.88
1:A:90:ASN:HA	1:A:93:ILE:HD12	1.55	0.87
2:B:360:ASN:HD21	2:B:363:GLU:HG2	1.39	0.87
2:B:329:ILE:O	2:B:330:VAL:HB	1.75	0.87
2:B:352:LEU:HD12	2:B:353:VAL:N	1.91	0.86
2:B:360:ASN:ND2	2:B:363:GLU:HG2	1.91	0.86
2:B:330:VAL:HG11	2:B:335:ILE:HD11	1.58	0.83
2:B:335:ILE:O	2:B:339:ILE:HG13	1.79	0.82
2:B:352:LEU:HD12	2:B:353:VAL:H	1.45	0.81
1:A:256:ARG:HH12	2:B:258:LYS:NZ	1.81	0.79
2:B:392:VAL:HG23	2:B:392:VAL:O	1.83	0.79
2:B:329:ILE:O	2:B:329:ILE:HG13	1.82	0.77
2:B:310:LYS:O	2:B:313:THR:HB	1.83	0.77
2:B:35:ALA:HB1	2:B:43:ILE:HD13	1.66	0.76
2:B:212:PRO:HD2	2:B:213:GLU:OE2	1.85	0.76
2:B:122:ASP:HB3	2:B:125:CYS:SG	2.26	0.75
2:B:269:PHE:CE1	2:B:296:LEU:HD11	2.21	0.75
2:B:348:LEU:HD21	2:B:373:LEU:CD2	2.16	0.75
1:A:23:PHE:CE2	1:A:55:HIS:HB3	2.24	0.73
2:B:21:ARG:HD3	2:B:107:GLU:OE2	1.89	0.73
2:B:270:VAL:HG22	2:B:324:ASN:HB3	1.69	0.72
2:B:269:PHE:HE1	2:B:296:LEU:HD11	1.52	0.72

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:122:ASP:OD2	2:B:125:CYS:SG	2.46	0.72
2:B:135:GLN:O	2:B:135:GLN:HG2	1.88	0.72
1:A:171:HIS:O	1:A:175:GLN:HG2	1.91	0.71
2:B:1:MET:HE1	2:B:3:LEU:HG	1.73	0.71
2:B:179:PRO:HG2	2:B:180:LEU:H	1.56	0.70
2:B:38:LEU:O	2:B:39:ASN:HB3	1.92	0.70
2:B:120:LEU:HD23	2:B:120:LEU:N	2.07	0.69
2:B:348:LEU:HD21	2:B:373:LEU:HD23	1.75	0.69
1:A:90:ASN:HA	1:A:93:ILE:CD1	2.23	0.68
1:A:6:SER:OG	1:A:126:ARG:HB2	1.94	0.68
2:B:144:ALA:HA	2:B:151:ILE:CD1	2.24	0.67
1:A:241:PRO:HB3	1:A:303:ARG:CZ	2.25	0.67
2:B:111:ILE:HD12	2:B:111:ILE:H	1.59	0.67
2:B:1:MET:HE3	2:B:2:ASN:C	2.15	0.66
2:B:350:VAL:HG12	2:B:351:PRO:HD2	1.78	0.66
1:A:10:LEU:HD23	1:A:128:ILE:HD11	1.78	0.66
1:A:221:ASN:H	1:A:221:ASN:HD22	1.44	0.66
1:A:61:PHE:CD1	1:A:61:PHE:N	2.66	0.64
2:B:88:LEU:HD12	2:B:89:ALA:N	2.12	0.64
2:B:176:PHE:O	2:B:181:GLN:HG3	1.98	0.64
2:B:251:GLU:OE2	2:B:254:LYS:HD2	1.98	0.63
2:B:330:VAL:CG1	2:B:335:ILE:HD11	2.27	0.63
2:B:342:ALA:O	2:B:346:LEU:HB2	1.99	0.63
2:B:41:LYS:O	2:B:71:LYS:HB3	1.98	0.63
1:A:82:PRO:HB2	1:A:83:PRO:HD3	1.80	0.63
1:A:17:LYS:HA	1:A:41:ASN:HB3	1.81	0.62
1:A:221:ASN:H	1:A:221:ASN:ND2	1.97	0.61
2:B:111:ILE:HD12	2:B:111:ILE:N	2.15	0.61
2:B:138:VAL:HG12	2:B:139:ASP:N	2.15	0.61
2:B:267:ALA:HA	2:B:291:ALA:HB3	1.83	0.61
1:A:168:GLU:OE2	2:B:355:ARG:NH1	2.34	0.60
2:B:59:PHE:CZ	2:B:83:MET:HG2	2.36	0.60
1:A:94:ASP:OD1	1:A:118:ARG:NH2	2.20	0.60
1:A:127:LEU:HD23	1:A:193:GLY:HA3	1.82	0.60
1:A:90:ASN:CA	1:A:93:ILE:HD12	2.29	0.60
1:A:256:ARG:HH12	2:B:258:LYS:HZ3	1.48	0.60
2:B:90:THR:C	2:B:92:GLN:H	2.04	0.60
2:B:38:LEU:O	2:B:39:ASN:CB	2.50	0.59
1:A:122:GLN:NE2	1:A:125:THR:HB	2.17	0.59
2:B:107:GLU:O	2:B:107:GLU:HG3	2.02	0.59
2:B:144:ALA:HA	2:B:151:ILE:HD11	1.85	0.59

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:269:PHE:HB2	2:B:294:LEU:HD23	1.85	0.59
2:B:189:LYS:O	2:B:193:ASN:ND2	2.36	0.58
2:B:113:ARG:HH11	2:B:113:ARG:CG	2.12	0.58
2:B:387:LYS:O	2:B:388:ALA:C	2.42	0.58
1:A:17:LYS:HE2	1:A:41:ASN:OD1	2.03	0.58
2:B:257:LEU:HD12	2:B:311:LEU:HD11	1.87	0.56
2:B:369:THR:HG22	2:B:370:ASN:ND2	2.20	0.56
1:A:245:PHE:HB2	1:A:292:LEU:HD21	1.86	0.56
2:B:140:ILE:HG23	2:B:141:GLU:N	2.20	0.56
2:B:129:VAL:HG22	2:B:155:GLN:HE21	1.69	0.56
1:A:36:LEU:HD11	1:A:58:LEU:HD11	1.88	0.56
2:B:178:GLY:O	2:B:181:GLN:HB2	2.06	0.56
1:A:156:ILE:HD12	1:A:296:ILE:HD13	1.87	0.55
2:B:303:SER:O	2:B:306:TYR:HB3	2.06	0.55
2:B:245:ASN:HB2	2:B:250:ASN:HD21	1.72	0.55
1:A:61:PHE:CG	1:A:67:ALA:HB2	2.42	0.55
2:B:144:ALA:HA	2:B:151:ILE:HD12	1.89	0.55
2:B:120:LEU:HD23	2:B:120:LEU:H	1.70	0.54
2:B:319:GLU:O	2:B:350:VAL:HG12	2.07	0.54
1:A:173:THR:O	1:A:176:VAL:HG22	2.08	0.54
2:B:90:THR:OG1	2:B:93:THR:HG23	2.07	0.54
2:B:288:GLY:O	2:B:289:LYS:HD3	2.07	0.54
1:A:19:ILE:HG13	1:A:19:ILE:O	2.06	0.54
1:A:36:LEU:HD21	1:A:42:LEU:HB3	1.90	0.54
2:B:164:ASP:HB2	4:B:403:HOH:O	2.07	0.54
2:B:186:ASP:O	2:B:190:LYS:HG3	2.08	0.54
2:B:198:ILE:HD12	2:B:225:PHE:CE1	2.43	0.53
2:B:113:ARG:HG3	2:B:113:ARG:NH1	2.15	0.53
2:B:192:TYR:OH	2:B:196:LEU:HD11	2.09	0.53
1:A:14:LYS:H	1:A:140:GLU:HB3	1.74	0.53
1:A:155:ARG:NH2	1:A:305:MET:O	2.29	0.53
2:B:144:ALA:CA	2:B:151:ILE:HD11	2.38	0.53
1:A:119:LEU:HD21	1:A:127:LEU:HB3	1.89	0.53
2:B:365:GLN:HG2	2:B:377:SER:OG	2.08	0.53
2:B:163:LYS:N	2:B:166:GLN:OE1	2.32	0.53
2:B:90:THR:C	2:B:92:GLN:N	2.61	0.53
2:B:122:ASP:CB	2:B:125:CYS:SG	2.97	0.53
2:B:194:LEU:O	2:B:198:ILE:HG12	2.09	0.53
1:A:36:LEU:HD11	1:A:58:LEU:CD1	2.39	0.52
2:B:135:GLN:HE21	2:B:150:LEU:HD13	1.75	0.52
2:B:107:GLU:C	2:B:108:ALA:O	2.47	0.52

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:174:THR:O	1:A:177:GLY:N	2.36	0.52
2:B:32:LEU:HD13	2:B:73:PRO:O	2.09	0.52
2:B:157:ASP:HB3	2:B:160:GLU:HB3	1.92	0.52
2:B:188:ILE:HA	2:B:191:LEU:HD12	1.92	0.52
2:B:217:VAL:HB	2:B:219:PHE:CZ	2.46	0.51
2:B:109:LEU:HD22	2:B:211:THR:HG22	1.93	0.51
1:A:290:ALA:HB3	2:B:382:GLU:OE1	2.10	0.51
2:B:182:ASN:HD22	2:B:182:ASN:H	1.57	0.51
2:B:246:GLU:HG2	2:B:247:PRO:HD2	1.93	0.51
1:A:57:GLY:O	1:A:58:LEU:HG	2.10	0.51
2:B:98:VAL:CG1	2:B:244:GLU:HG2	2.41	0.51
2:B:343:CYS:HB3	2:B:373:LEU:HD22	1.92	0.51
1:A:113:VAL:HG11	2:B:127:GLY:HA2	1.93	0.51
1:A:24:THR:OG1	1:A:47:THR:HG21	2.11	0.51
2:B:379:VAL:HG12	2:B:383:ASP:HB2	1.93	0.51
2:B:379:VAL:HG12	2:B:383:ASP:CB	2.41	0.50
1:A:223:GLU:HB2	1:A:274:LYS:HE3	1.93	0.50
1:A:122:GLN:HE22	1:A:125:THR:CA	2.24	0.50
1:A:14:LYS:N	1:A:140:GLU:HB3	2.27	0.50
1:A:43:VAL:O	1:A:58:LEU:HD22	2.11	0.50
1:A:94:ASP:C	1:A:96:GLU:H	2.14	0.50
2:B:308:ALA:O	2:B:312:LEU:HG	2.12	0.50
2:B:113:ARG:CG	2:B:113:ARG:NH1	2.74	0.50
2:B:340:THR:HG22	2:B:373:LEU:HD13	1.94	0.50
1:A:4:THR:O	1:A:7:ARG:HG3	2.12	0.49
1:A:162:SER:HB2	1:A:259:NEP:O1P	2.11	0.49
1:A:299:GLU:OE2	1:A:302:LYS:NZ	2.36	0.49
2:B:116:TYR:CD2	2:B:140:ILE:HD13	2.47	0.49
2:B:270:VAL:HA	2:B:324:ASN:O	2.12	0.49
1:A:164:THR:HA	1:A:167:TYR:CD2	2.47	0.49
2:B:388:ALA:O	2:B:391:SER:HB3	2.12	0.49
2:B:138:VAL:CG1	2:B:139:ASP:N	2.75	0.49
2:B:147:ASN:HB3	2:B:150:LEU:CD1	2.43	0.49
2:B:120:LEU:N	2:B:120:LEU:CD2	2.76	0.49
2:B:125:CYS:O	2:B:126:ASN:O	2.31	0.49
2:B:152:PHE:CD1	2:B:152:PHE:N	2.80	0.49
1:A:36:LEU:CD2	1:A:42:LEU:HB3	2.42	0.48
2:B:170:MET:O	2:B:171:ALA:C	2.50	0.48
2:B:348:LEU:O	2:B:348:LEU:HG	2.13	0.48
2:B:192:TYR:O	2:B:195:PHE:HB3	2.13	0.48
1:A:66:GLU:O	1:A:68:LYS:N	2.47	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:58:LEU:HD23	1:A:59:PRO:HD2	1.95	0.48
2:B:192:TYR:CZ	2:B:196:LEU:HD11	2.48	0.48
1:A:43:VAL:HG21	1:A:71:THR:CB	2.44	0.48
1:A:89:ILE:O	1:A:93:ILE:HD12	2.13	0.48
1:A:221:ASN:ND2	1:A:221:ASN:N	2.57	0.48
2:B:304:GLN:O	2:B:305:VAL:C	2.50	0.48
2:B:327:GLY:O	2:B:329:ILE:N	2.46	0.48
2:B:245:ASN:HB2	2:B:250:ASN:ND2	2.28	0.48
2:B:275:LEU:HD11	2:B:324:ASN:HD21	1.79	0.48
1:A:122:GLN:HE22	1:A:125:THR:HB	1.77	0.48
1:A:234:ASN:O	1:A:239:SER:HB2	2.14	0.48
2:B:329:ILE:O	2:B:329:ILE:CG1	2.55	0.48
1:A:106:GLY:O	2:B:228:ASN:HB3	2.15	0.47
1:A:156:ILE:HD12	1:A:296:ILE:CD1	2.43	0.47
2:B:135:GLN:NE2	2:B:150:LEU:CD1	2.77	0.47
2:B:360:ASN:HD21	2:B:363:GLU:CG	2.19	0.47
2:B:18:LYS:HE2	2:B:183:GLN:OE1	2.14	0.47
2:B:163:LYS:HG2	2:B:165:SER:OG	2.14	0.47
2:B:206:ASN:O	2:B:220:ASP:HB2	2.15	0.47
1:A:60:VAL:C	1:A:61:PHE:CD1	2.88	0.47
1:A:203:PHE:HB3	1:A:209:THR:HG21	1.96	0.47
1:A:173:THR:O	1:A:176:VAL:CG2	2.63	0.47
2:B:371:SER:HB2	2:B:373:LEU:HD12	1.95	0.47
2:B:42:GLU:HB3	2:B:108:ALA:HB3	1.96	0.47
2:B:109:LEU:HD22	2:B:211:THR:CG2	2.45	0.46
2:B:231:PHE:CD1	2:B:231:PHE:C	2.88	0.46
2:B:140:ILE:HG23	2:B:141:GLU:HG3	1.97	0.46
1:A:121:ARG:HG3	1:A:121:ARG:HH11	1.79	0.46
2:B:90:THR:O	2:B:92:GLN:N	2.49	0.46
2:B:163:LYS:HE3	2:B:165:SER:OG	2.16	0.46
2:B:270:VAL:HG21	2:B:279:THR:OG1	2.15	0.46
2:B:265:ASN:OD1	2:B:266:ILE:HG13	2.15	0.46
1:A:50:LYS:O	1:A:52:GLY:N	2.48	0.46
2:B:210:GLU:HA	2:B:215:GLN:O	2.15	0.45
2:B:361:VAL:HG12	2:B:365:GLN:OE1	2.16	0.45
2:B:307:GLN:OE1	2:B:310:LYS:HD3	2.17	0.45
2:B:330:VAL:CG1	2:B:335:ILE:CD1	2.93	0.45
2:B:152:PHE:CE2	2:B:173:ASN:ND2	2.84	0.45
2:B:51:ALA:CB	2:B:93:THR:HG22	2.46	0.45
2:B:217:VAL:HB	2:B:219:PHE:CE1	2.51	0.45
2:B:125:CYS:SG	2:B:155:GLN:NE2	2.90	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:156:ILE:HG21	1:A:213:ILE:HG13	1.98	0.45
2:B:385:ALA:O	2:B:389:VAL:HG23	2.17	0.45
2:B:49:ILE:O	2:B:49:ILE:HG13	2.16	0.45
1:A:32:SER:HB3	1:A:42:LEU:HD21	1.99	0.45
2:B:27:THR:HG23	2:B:30:GLU:OE2	2.17	0.45
2:B:248:ILE:HG21	2:B:259:TYR:CZ	2.52	0.45
2:B:259:TYR:O	2:B:260:ILE:HG13	2.17	0.45
1:A:127:LEU:HD23	1:A:193:GLY:CA	2.47	0.44
1:A:82:PRO:CB	1:A:83:PRO:HD3	2.47	0.44
2:B:128:PRO:C	2:B:129:VAL:HG23	2.37	0.44
1:A:51:GLY:HA3	1:A:62:ASN:CG	2.38	0.44
2:B:380:ASP:OD1	2:B:380:ASP:C	2.55	0.44
1:A:89:ILE:C	1:A:93:ILE:HD12	2.38	0.44
2:B:80:ALA:O	2:B:81:LYS:C	2.54	0.44
2:B:107:GLU:O	2:B:108:ALA:O	2.36	0.44
2:B:192:TYR:CE1	2:B:196:LEU:HD11	2.53	0.44
2:B:156:ILE:HG21	2:B:192:TYR:CE2	2.52	0.44
2:B:182:ASN:H	2:B:182:ASN:ND2	2.15	0.44
2:B:116:TYR:OH	2:B:204:GLU:OE2	2.17	0.44
2:B:147:ASN:HB3	2:B:150:LEU:HG	1.99	0.44
1:A:89:ILE:HG22	1:A:93:ILE:CD1	2.36	0.44
2:B:198:ILE:HD12	2:B:225:PHE:HE1	1.83	0.44
2:B:304:GLN:C	2:B:306:TYR:N	2.68	0.44
1:A:78:ILE:HG22	1:A:80:VAL:H	1.83	0.43
1:A:85:ALA:HB1	1:A:107:ILE:HD11	2.00	0.43
2:B:264:GLY:HA3	2:B:289:LYS:O	2.19	0.43
2:B:35:ALA:HB1	2:B:43:ILE:CD1	2.42	0.43
2:B:380:ASP:OD1	2:B:382:GLU:N	2.52	0.43
1:A:113:VAL:CG1	2:B:127:GLY:HA2	2.48	0.43
1:A:176:VAL:HG23	1:A:178:LEU:HG	2.01	0.43
2:B:1:MET:CE	2:B:3:LEU:HG	2.47	0.43
2:B:98:VAL:HG11	2:B:244:GLU:HG2	2.00	0.43
2:B:345:GLU:O	2:B:345:GLU:HG2	2.19	0.43
1:A:14:LYS:HG2	1:A:140:GLU:OE1	2.19	0.42
2:B:33:GLU:HB3	2:B:37:ARG:HH12	1.84	0.42
2:B:392:VAL:O	2:B:392:VAL:CG2	2.54	0.42
1:A:162:SER:OG	1:A:165:LEU:HB3	2.19	0.42
2:B:8:SER:O	2:B:12:MET:HG3	2.20	0.42
2:B:128:PRO:HB2	2:B:156:ILE:HB	2.01	0.42
2:B:370:ASN:ND2	2:B:370:ASN:N	2.67	0.42
1:A:85:ALA:CB	1:A:107:ILE:HD11	2.49	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:204:LEU:HD23	1:A:204:LEU:HA	1.78	0.42
2:B:72:ASP:O	2:B:74:GLU:N	2.52	0.42
2:B:360:ASN:C	2:B:360:ASN:HD22	2.21	0.42
1:A:165:LEU:O	1:A:168:GLU:HB2	2.18	0.42
2:B:1:MET:C	2:B:1:MET:HE2	2.40	0.42
2:B:45:LEU:HD21	2:B:83:MET:CE	2.50	0.42
2:B:337:ASN:HD22	2:B:337:ASN:HA	1.69	0.42
1:A:10:LEU:HD23	1:A:128:ILE:CD1	2.48	0.42
1:A:169:ALA:O	1:A:173:THR:HG23	2.19	0.42
1:A:256:ARG:HH11	1:A:256:ARG:HB3	1.85	0.42
1:A:148:GLY:O	1:A:150:ILE:N	2.53	0.42
1:A:168:GLU:O	1:A:172:GLN:HG3	2.19	0.42
2:B:135:GLN:NE2	2:B:150:LEU:HD13	2.33	0.42
2:B:387:LYS:O	2:B:390:ALA:N	2.52	0.42
1:A:89:ILE:CG2	1:A:93:ILE:HD11	2.37	0.42
2:B:44:VAL:HG12	2:B:46:LYS:HG2	2.01	0.42
1:A:218:ILE:HG23	1:A:248:GLY:HA3	2.02	0.41
2:B:44:VAL:HG22	2:B:69:LEU:CD2	2.50	0.41
2:B:93:THR:OG1	2:B:97:GLY:HA2	2.19	0.41
2:B:139:ASP:O	2:B:141:GLU:N	2.53	0.41
2:B:176:PHE:CD2	2:B:184:ALA:CB	3.03	0.41
2:B:179:PRO:CG	2:B:180:LEU:H	2.29	0.41
1:A:88:ALA:O	1:A:89:ILE:C	2.55	0.41
2:B:192:TYR:CE1	2:B:196:LEU:CD1	3.03	0.41
2:B:191:LEU:O	2:B:194:LEU:HB3	2.20	0.41
2:B:45:LEU:HD13	2:B:105:VAL:HG22	2.02	0.41
2:B:75:VAL:HG12	2:B:79:LEU:HD12	2.03	0.41
2:B:163:LYS:HG2	2:B:165:SER:H	1.84	0.41
2:B:257:LEU:HD22	2:B:296:LEU:HD23	2.02	0.41
2:B:341:LYS:O	2:B:345:GLU:HB3	2.21	0.41
1:A:221:ASN:HD22	1:A:221:ASN:N	2.08	0.41
2:B:43:ILE:O	2:B:69:LEU:HA	2.20	0.41
2:B:96:GLU:H	2:B:96:GLU:CD	2.24	0.41
2:B:307:GLN:O	2:B:310:LYS:HB3	2.21	0.41
2:B:25:ALA:O	2:B:102:LYS:HA	2.21	0.41
1:A:9:HIS:ND1	1:A:98:PRO:HB3	2.35	0.41
2:B:301:LYS:O	2:B:302:GLU:C	2.59	0.41
2:B:343:CYS:HB3	2:B:373:LEU:CD2	2.51	0.41
2:B:373:LEU:HB3	2:B:375:ILE:HG13	2.03	0.41
1:A:122:GLN:HE22	1:A:125:THR:C	2.24	0.41
1:A:135:VAL:O	1:A:146:MET:HE1	2.21	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:186:ASP:HB3	4:B:409:HOH:O	2.21	0.41
2:B:209:GLY:O	2:B:216:VAL:HA	2.21	0.41
2:B:356:LEU:HD23	2:B:356:LEU:HA	1.96	0.41
1:A:125:THR:HG22	1:A:126:ARG:N	2.36	0.40
2:B:133:SER:HB2	2:B:150:LEU:O	2.21	0.40
2:B:138:VAL:CG1	2:B:139:ASP:H	2.35	0.40
2:B:360:ASN:CG	2:B:363:GLU:HG2	2.40	0.40
1:A:23:PHE:CE1	1:A:55:HIS:HD2	2.39	0.40
1:A:122:GLN:CD	1:A:125:THR:HB	2.41	0.40
1:A:167:TYR:N	1:A:167:TYR:CD1	2.83	0.40
2:B:99:LYS:HG2	2:B:100:VAL:N	2.36	0.40
2:B:116:TYR:O	2:B:132:GLY:HA2	2.22	0.40
1:A:3:TYR:CE1	1:A:119:LEU:HG	2.57	0.40
1:A:61:PHE:N	1:A:61:PHE:HD1	2.17	0.40
2:B:44:VAL:HG23	2:B:108:ALA:HA	2.03	0.40
2:B:252:ALA:O	2:B:257:LEU:HB2	2.21	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	302/305 (99%)	266 (88%)	25 (8%)	11 (4%)	3 7
2	B	391/395 (99%)	334 (85%)	43 (11%)	14 (4%)	3 7
All	All	693/700 (99%)	600 (87%)	68 (10%)	25 (4%)	3 7

All (25) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	51	GLY
1	A	67	ALA

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	73	ALA
2	B	140	ILE
2	B	314	ALA
2	B	330	VAL
1	A	66	GLU
2	B	39	ASN
2	B	97	GLY
2	B	126	ASN
2	B	178	GLY
2	B	328	GLY
2	B	391	SER
1	A	52	GLY
2	B	108	ALA
1	A	69	GLU
1	A	119	LEU
1	A	149	HIS
1	A	237	PRO
2	B	91	LYS
1	A	95	ALA
2	B	302	GLU
2	B	347	GLU
2	B	345	GLU
1	A	47	THR

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	237/237 (100%)	220 (93%)	17 (7%)	14	34
2	B	317/319 (99%)	291 (92%)	26 (8%)	11	26
All	All	554/556 (100%)	511 (92%)	43 (8%)	12	29

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	THR
1	A	11	TYR
1	A	17	LYS
1	A	33	GLN
1	A	37	GLU
1	A	40	THR
1	A	41	ASN
1	A	61	PHE
1	A	70	GLN
1	A	124	LYS
1	A	146	MET
1	A	181	SER
1	A	221	ASN
1	A	237	PRO
1	A	256	ARG
1	A	272	LYS
1	A	303	ARG
2	B	1	MET
2	B	26	ASP
2	B	71	LYS
2	B	73	PRO
2	B	78	GLN
2	B	93	THR
2	B	96	GLU
2	B	107	GLU
2	B	109	LEU
2	B	124	SER
2	B	130	LEU
2	B	152	PHE
2	B	164	ASP
2	B	165	SER
2	B	249	GLU
2	B	269	PHE
2	B	285	LEU
2	B	293	PHE
2	B	330	VAL
2	B	346	LEU
2	B	347	GLU
2	B	348	LEU
2	B	350	VAL
2	B	360	ASN
2	B	373	LEU
2	B	379	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	33	GLN
1	A	55	HIS
1	A	221	ASN
1	A	225	ASN
1	A	233	HIS
2	B	20	GLN
2	B	135	GLN
2	B	155	GLN
2	B	168	GLN
2	B	173	ASN
2	B	182	ASN
2	B	250	ASN
2	B	324	ASN
2	B	337	ASN
2	B	370	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)

1 non-standard protein/DNA/RNA residue 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	$\#\|Z\| > 2$
1	NEP	A	259	1	11,14,15	2.18	4 (36%)	4,20,22	2.98	2 (50%)

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
1	NEP	A	259	1	-	2/5/12/14	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	259	NEP	P-O3P	4.78	1.51	1.47
1	A	259	NEP	P-O1P	-3.10	1.48	1.54
1	A	259	NEP	P-O2P	-2.70	1.49	1.54
1	A	259	NEP	P-NE2	2.43	1.89	1.76

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	259	NEP	O1P-P-O3P	-5.20	102.84	113.76
1	A	259	NEP	CB-CG-CD2	2.39	132.38	127.89

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	259	NEP	CA-CB-CG-ND1
1	A	259	NEP	CA-CB-CG-CD2

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	259	NEP	1	0

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

2 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond

length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
3	SO4	A	307	-	4,4,4	0.45	0	6,6,6	0.45	0
3	SO4	B	396	-	4,4,4	0.47	0	6,6,6	0.22	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 (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	304/305 (99%)	-0.41	1 (0%)	94	95	16, 35, 63, 75	0
2	B	393/395 (99%)	-0.25	3 (0%)	86	87	18, 50, 72, 83	0
All	All	697/700 (99%)	-0.32	4 (0%)	89	91	16, 44, 70, 83	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	329	ILE	2.8
2	B	35	ALA	2.4
1	A	44	GLY	2.2
2	B	32	LEU	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
1	NEP	A	259	14/15	0.98	0.17	16,20,21,24	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
3	SO4	B	396	5/5	0.96	0.12	71,71,72,72	0
3	SO4	A	307	5/5	0.98	0.15	50,50,51,52	0

6.5 Other polymers (i)

There are no such residues in this entry.