



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 09:22 PM UTC

PDB ID : 1C7J / pdb_00001c7j
Title : PNB ESTERASE 56C8
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Deposited on : 2000-02-21
Resolution : 1.60 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

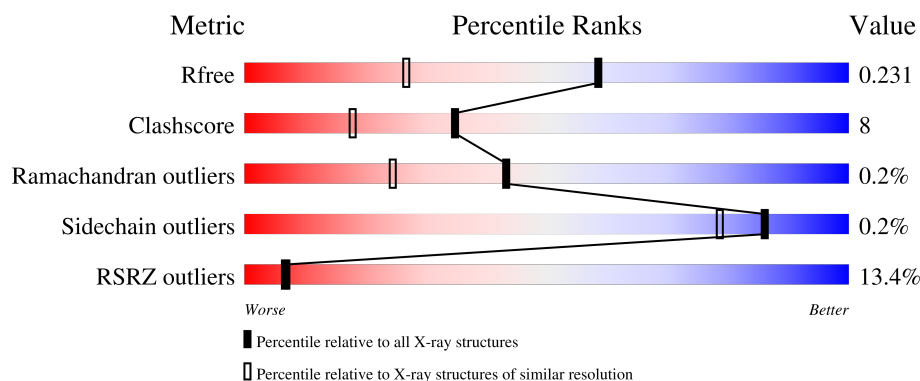
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.60 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	4673 (1.60-1.60)
Clashscore	190562	4931 (1.60-1.60)
Ramachandran outliers	187476	4831 (1.60-1.60)
Sidechain outliers	187428	4830 (1.60-1.60)
RSRZ outliers	180081	4672 (1.60-1.60)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	489	13% 83% 16%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 4055 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.

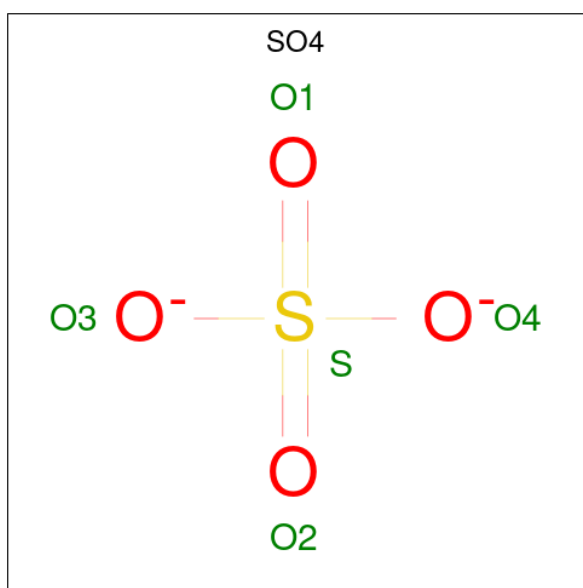
- Molecule 1 is a protein called PROTEIN (PARA-NITROBENZYL ESTERASE).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	485	3779	2420	621	727	11	0	0	0

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	60	VAL	ILE	engineered mutation	UNP P37967
A	144	MET	LEU	engineered mutation	UNP P37967
A	317	SER	PRO	engineered mutation	UNP P37967
A	322	ARG	HIS	engineered mutation	UNP P37967
A	334	SER	LEU	engineered mutation	UNP P37967
A	358	VAL	MET	engineered mutation	UNP P37967
A	370	PHE	TYR	engineered mutation	UNP P37967

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

- Molecule 3 is POTASSIUM ION (CCD ID: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	K	0	0
			1	1		

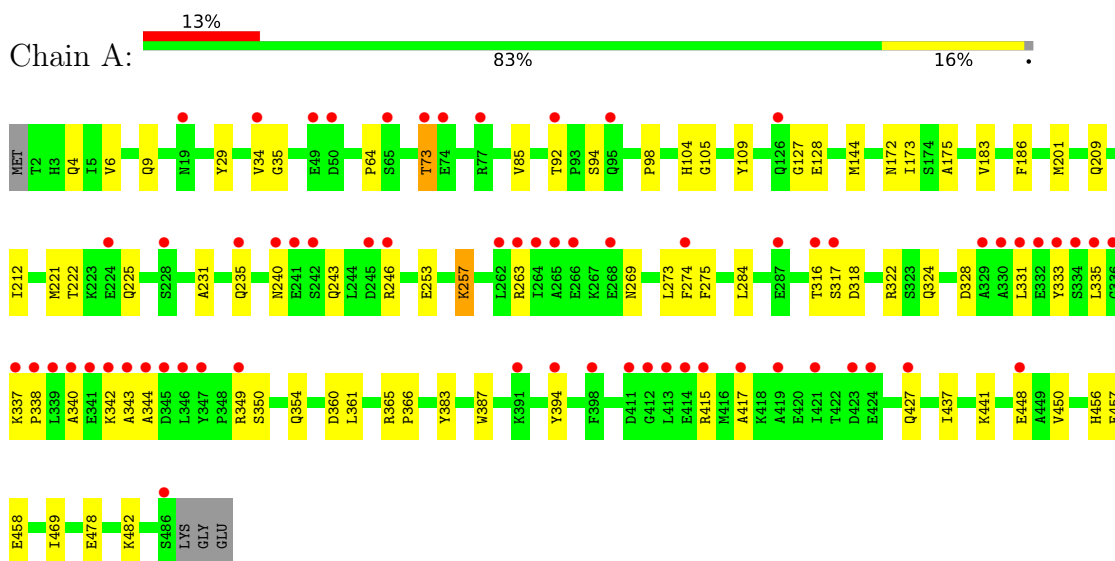
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	270	Total	O	0	0
			270	270		

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: PROTEIN (PARA-NITROBENZYL ESTERASE)



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	67.51Å 81.83Å 98.84Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	25.00 – 1.60 25.00 – 1.60	Depositor EDS
% Data completeness (in resolution range)	89.0 (25.00-1.60) 91.5 (25.00-1.60)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.04 (at 1.60Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.214 , 0.228 0.218 , 0.231	Depositor DCC
R_{free} test set	6762 reflections (10.13%)	wwPDB-VP
Wilson B-factor (Å ²)	15.6	Xtriage
Anisotropy	0.526	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 34.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4055	wwPDB-VP
Average B, all atoms (Å ²)	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.59% 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: SO4, K

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	0.34	0/3882	0.87	12/5291 (0.2%)

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	4	GLN	N-CA-C	7.87	122.36	111.74
1	A	73	THR	N-CA-C	-7.02	103.70	112.90
1	A	201	MET	N-CA-C	6.58	118.18	109.83
1	A	284	LEU	CA-C-N	5.81	125.94	119.32
1	A	284	LEU	C-N-CA	5.81	125.94	119.32
1	A	387	TRP	N-CA-C	5.38	117.66	110.35
1	A	221	MET	N-CA-C	-5.33	101.88	110.14
1	A	457	GLU	N-CA-C	5.33	119.40	113.01
1	A	360	ASP	N-CA-C	5.29	116.85	111.14
1	A	173	ILE	N-CA-C	5.22	117.57	111.05
1	A	64	PRO	N-CA-C	-5.17	103.66	111.41
1	A	144	MET	N-CA-C	-5.07	99.03	107.99

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	3779	0	3667	56	0
2	A	5	0	0	0	0
3	A	1	0	0	0	0
4	A	270	0	0	3	0
All	All	4055	0	3667	56	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 8.

All (56) 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:324:GLN:H	1:A:349:ARG:HE	1.08	0.96
1:A:92:THR:HG22	1:A:94:SER:H	1.31	0.93
1:A:342:LYS:HD3	1:A:343:ALA:H	1.38	0.88
1:A:342:LYS:HD3	1:A:343:ALA:N	1.88	0.88
1:A:324:GLN:N	1:A:349:ARG:HE	1.71	0.87
1:A:337:LYS:HB2	1:A:338:PRO:HD3	1.60	0.81
1:A:324:GLN:H	1:A:349:ARG:NE	1.81	0.78
1:A:175:ALA:HB1	4:A:547:HOH:O	1.83	0.77
1:A:337:LYS:HB2	1:A:338:PRO:CD	2.16	0.75
1:A:92:THR:HG21	1:A:128:GLU:OE1	1.87	0.75
1:A:9:GLN:HE21	1:A:172:ASN:HD21	1.36	0.74
1:A:183:VAL:H	1:A:209:GLN:HE21	1.41	0.67
1:A:73:THR:HG22	1:A:415:ARG:NH2	2.11	0.66
1:A:183:VAL:H	1:A:209:GLN:NE2	1.93	0.66
1:A:73:THR:HG22	1:A:415:ARG:HH22	1.60	0.65
1:A:322:ARG:H	1:A:354:GLN:HE22	1.43	0.64
1:A:394:TYR:CD2	1:A:417:ALA:HA	2.32	0.63
1:A:269:ASN:HD21	1:A:273:LEU:H	1.46	0.63
1:A:263:ARG:HH21	1:A:275:PHE:HD2	1.46	0.61
1:A:342:LYS:O	1:A:343:ALA:HB3	2.04	0.58
1:A:104:HIS:HD2	1:A:105:GLY:O	1.87	0.57
1:A:450:VAL:HG22	1:A:469:ILE:HG21	1.86	0.56
1:A:257:LYS:HE3	1:A:257:LYS:HA	1.88	0.56
1:A:34:VAL:HG22	1:A:35:GLY:N	2.22	0.54
1:A:340:ALA:O	1:A:344:ALA:HB2	2.08	0.54
1:A:104:HIS:HE1	4:A:521:HOH:O	1.90	0.53
1:A:456:HIS:HD2	1:A:458:GLU:H	1.57	0.52
1:A:127:GLY:CA	1:A:437:ILE:HD13	2.41	0.51
1:A:263:ARG:HG2	1:A:274:PHE:CE2	2.46	0.50
1:A:365:ARG:HB3	1:A:366:PRO:HD3	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:231:ALA:O	1:A:235:GLN:HG3	2.10	0.50
1:A:349:ARG:O	1:A:350:SER:HB3	2.10	0.50
1:A:127:GLY:HA3	1:A:437:ILE:HD13	1.94	0.49
1:A:243:GLN:HB3	1:A:246:ARG:HD3	1.94	0.49
1:A:331:LEU:HD23	1:A:361:LEU:HD11	1.95	0.49
1:A:333:TYR:CD1	1:A:333:TYR:C	2.91	0.48
1:A:6:VAL:HG13	4:A:547:HOH:O	2.13	0.48
1:A:478:GLU:O	1:A:482:LYS:HG3	2.13	0.48
1:A:316:THR:HG22	1:A:317:SER:N	2.31	0.46
1:A:98:PRO:HG2	1:A:441:LYS:HG2	1.97	0.46
1:A:448:GLU:HA	1:A:448:GLU:OE1	2.16	0.45
1:A:324:GLN:HG3	1:A:328:ASP:OD2	2.18	0.44
1:A:240:ASN:N	1:A:240:ASN:HD22	2.15	0.44
1:A:29:TYR:CD1	1:A:85:VAL:HG13	2.53	0.44
1:A:263:ARG:HG2	1:A:274:PHE:CD2	2.53	0.44
1:A:253:GLU:H	1:A:253:GLU:CD	2.26	0.43
1:A:222:THR:OG1	1:A:225:GLN:HG3	2.19	0.43
1:A:9:GLN:NE2	1:A:172:ASN:HD21	2.09	0.43
1:A:316:THR:HG22	1:A:318:ASP:H	1.84	0.42
1:A:383:TYR:CD1	1:A:383:TYR:C	2.98	0.42
1:A:324:GLN:N	1:A:349:ARG:NE	2.50	0.41
1:A:427:GLN:NE2	1:A:427:GLN:HA	2.35	0.41
1:A:109:TYR:CG	1:A:263:ARG:NH2	2.88	0.41
1:A:342:LYS:O	1:A:343:ALA:CB	2.69	0.41
1:A:186:PHE:CB	1:A:212:ILE:HB	2.50	0.40
1:A:186:PHE:HB2	1:A:212:ILE:HB	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	483/489 (99%)	459 (95%)	23 (5%)	1 (0%)	43 24

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	335	LEU

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	400/403 (99%)	399 (100%)	1 (0%)	86 78

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	257	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	4	GLN
1	A	9	GLN
1	A	19	ASN
1	A	104	HIS
1	A	182	ASN
1	A	209	GLN
1	A	240	ASN
1	A	269	ASN
1	A	276	GLN
1	A	354	GLN
1	A	427	GLN
1	A	456	HIS
1	A	481	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	SO4	A	490	-	4,4,4	0.24	0	6,6,6	0.13	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 ⓘ

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	485/489 (99%)	0.63	65 (13%) 7 7	7, 16, 35, 46	0

All (65) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	335	LEU	7.2
1	A	339	LEU	7.0
1	A	336	GLY	6.7
1	A	343	ALA	5.9
1	A	240	ASN	5.9
1	A	333	TYR	5.3
1	A	338	PRO	5.2
1	A	263	ARG	5.0
1	A	414	GLU	4.6
1	A	415	ARG	4.5
1	A	394	TYR	4.3
1	A	337	LYS	4.3
1	A	331	LEU	4.2
1	A	349	ARG	4.1
1	A	266	GLU	4.0
1	A	268	GLU	3.9
1	A	340	ALA	3.9
1	A	486	SER	3.8
1	A	342	LYS	3.7
1	A	264	ILE	3.7
1	A	412	GLY	3.6
1	A	126	GLN	3.5
1	A	344	ALA	3.5
1	A	413	LEU	3.4
1	A	341	GLU	3.4
1	A	345	ASP	3.3
1	A	235	GLN	3.2

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Mol	Chain	Res	Type	RSRZ
1	A	411	ASP	3.2
1	A	50	ASP	3.1
1	A	346	LEU	3.1
1	A	317	SER	3.1
1	A	329	ALA	3.0
1	A	316	THR	2.9
1	A	332	GLU	2.9
1	A	262	LEU	2.9
1	A	65	SER	2.9
1	A	241	GLU	2.8
1	A	287	GLU	2.8
1	A	421	ILE	2.7
1	A	242	SER	2.7
1	A	398	PHE	2.7
1	A	92	THR	2.7
1	A	347	TYR	2.7
1	A	246	ARG	2.6
1	A	424	GLU	2.6
1	A	448	GLU	2.6
1	A	334	SER	2.6
1	A	419	ALA	2.6
1	A	74	GLU	2.5
1	A	49	GLU	2.5
1	A	34	VAL	2.4
1	A	19	ASN	2.3
1	A	245	ASP	2.3
1	A	265	ALA	2.3
1	A	427	GLN	2.3
1	A	228	SER	2.3
1	A	391	LYS	2.3
1	A	95	GLN	2.2
1	A	77	ARG	2.2
1	A	417	ALA	2.2
1	A	330	ALA	2.2
1	A	423	ASP	2.2
1	A	274	PHE	2.1
1	A	224	GLU	2.1
1	A	73	THR	2.1

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

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

6.3 Carbohydrates [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
2	SO4	A	490	5/5	0.92	0.11	31,31,32,32	0
3	K	A	500	1/1	0.98	0.09	7,7,7,7	0

6.5 Other polymers [i](#)

There are no such residues in this entry.