



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 05:34 PM UTC

PDB ID : 6YYN / pdb_00006yyn
Title : Structure of Cathepsin S in complex with Compound 14
Authors : Wagener, M.; Schade, M.; Merla, B.; Hars, U.; Kueckelhaus, S.Q.
Deposited on : 2020-05-05
Resolution : 2.22 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
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
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

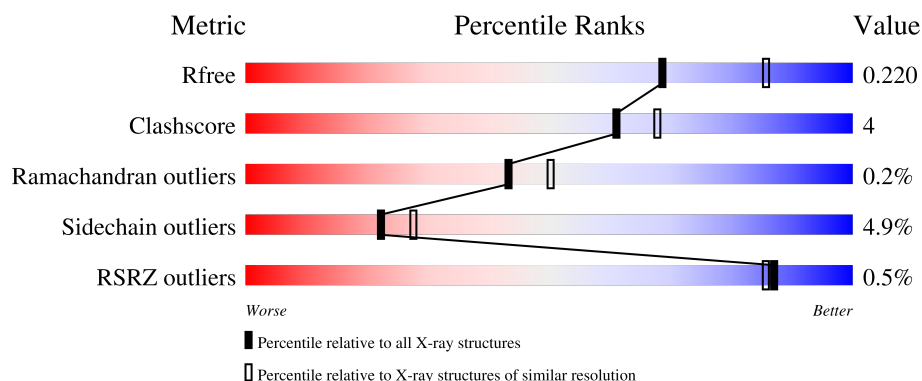
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.22 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	7682 (2.24-2.20)
Clashscore	190562	8402 (2.24-2.20)
Ramachandran outliers	187476	8303 (2.24-2.20)
Sidechain outliers	187428	8304 (2.24-2.20)
RSRZ outliers	180081	7683 (2.24-2.20)

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	AAA	225	 87% 10% .
1	BBB	225	 83% 14% ..

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 4029 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.

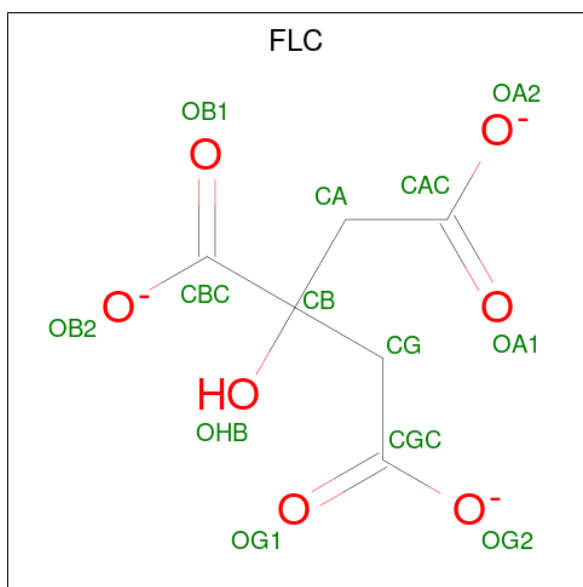
- Molecule 1 is a protein called Cathepsin S.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	AAA	220	Total	C	N	O	S	0	0	0
			1713	1082	293	326	12			
1	BBB	222	Total	C	N	O	S	0	0	0
			1734	1094	300	328	12			

There are 12 discrepancies between the modelled and reference sequences:

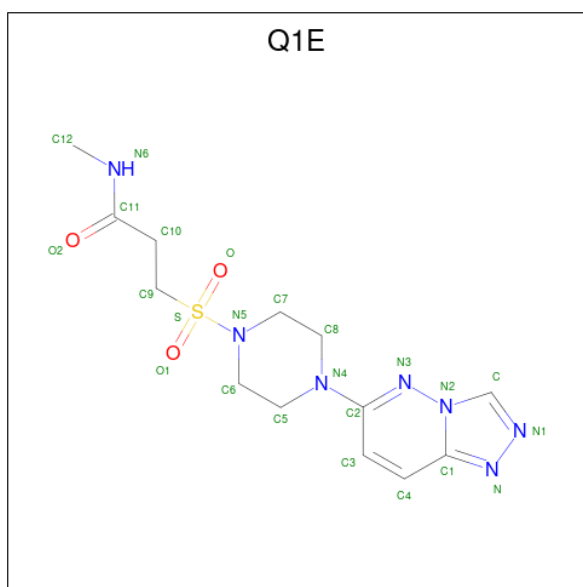
Chain	Residue	Modelled	Actual	Comment	Reference
AAA	218	HIS	-	expression tag	UNP P25774
AAA	219	HIS	-	expression tag	UNP P25774
AAA	220	HIS	-	expression tag	UNP P25774
AAA	221	HIS	-	expression tag	UNP P25774
AAA	222	HIS	-	expression tag	UNP P25774
AAA	223	HIS	-	expression tag	UNP P25774
BBB	218	HIS	-	expression tag	UNP P25774
BBB	219	HIS	-	expression tag	UNP P25774
BBB	220	HIS	-	expression tag	UNP P25774
BBB	221	HIS	-	expression tag	UNP P25774
BBB	222	HIS	-	expression tag	UNP P25774
BBB	223	HIS	-	expression tag	UNP P25774

- Molecule 2 is CITRATE ANION (CCD ID: FLC) (formula: C₆H₅O₇).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	AAA	1	Total	C	O	0	0
			13	6	7		
2	AAA	1	Total	C	O	0	0
			13	6	7		

- Molecule 3 is {N}-methyl-3-[4-([1,2,4]triazolo[4,3-b]pyridazin-6-yl)piperazin-1-yl]sulfonylpropanamide (CCD ID: Q1E) (formula: C₁₃H₁₉N₇O₃S) (labeled as "Ligand of Interest" by depositor).



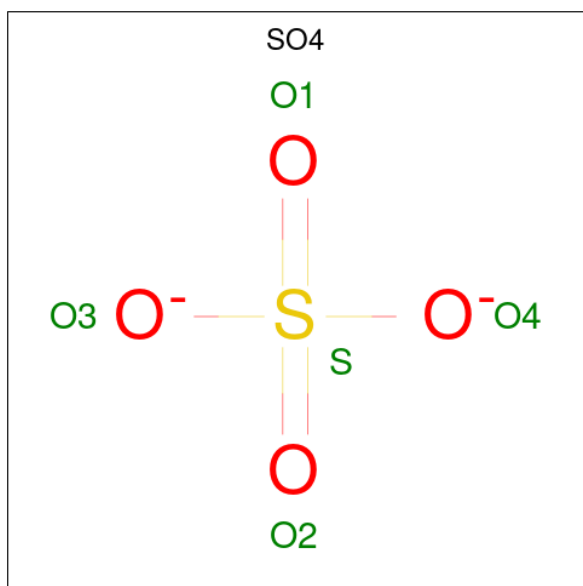
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	AAA	1	Total	C	N	O	S	0	0
			24	13	7	3	1		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	BBB	1	Total	C	N	O	S	0	0
			24	13	7	3	1		

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



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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	AAA	248	Total	O	0	0
			248	248		
5	BBB	230	Total	O	0	0
			230	230		

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.

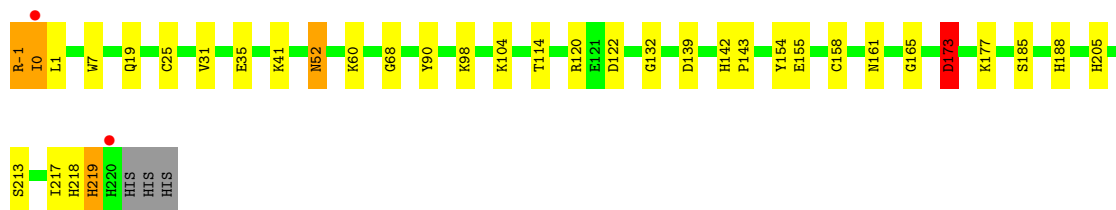
- Molecule 1: Cathepsin S

Chain AAA:  87% 10% .



- Molecule 1: Cathepsin S

Chain BBB:  % 83% 14% ..



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 2 2	Depositor
Cell constants a, b, c, α , β , γ	85.67Å 85.67Å 150.92Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.51 – 2.22 19.51 – 2.22	Depositor EDS
% Data completeness (in resolution range)	99.0 (19.51-2.22) 99.0 (19.51-2.22)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.11 (at 2.21Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.158 , 0.217 0.166 , 0.220	Depositor DCC
R_{free} test set	1436 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	31.0	Xtriage
Anisotropy	0.258	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 52.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4029	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.25% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, FLC, Q1E

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	AAA	1.17	2/1757 (0.1%)	1.34	4/2376 (0.2%)
1	BBB	1.17	2/1779 (0.1%)	1.34	5/2405 (0.2%)
All	All	1.17	4/3536 (0.1%)	1.34	9/4781 (0.2%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	BBB	-1	ARG	N-CA	6.30	1.58	1.46
1	AAA	117	PRO	C-O	-5.68	1.17	1.23
1	BBB	218	HIS	CE1-NE2	5.10	1.37	1.32
1	AAA	35	GLU	C-O	5.06	1.30	1.24

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AAA	141	ARG	CB-CA-C	8.64	121.85	111.22
1	BBB	68	GLY	CA-C-O	-6.91	117.73	122.22
1	BBB	218	HIS	CA-C-N	6.63	129.45	120.44
1	BBB	218	HIS	C-N-CA	6.63	129.45	120.44
1	AAA	145	PHE	CA-C-N	5.78	128.02	120.28
1	AAA	145	PHE	C-N-CA	5.78	128.02	120.28
1	BBB	98	LYS	CB-CA-C	5.42	118.91	109.65
1	AAA	141	ARG	CB-CG-CD	5.37	123.64	111.30
1	BBB	173	ASP	CB-CA-C	5.29	118.80	109.80

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	AAA	1713	0	1632	11	0
1	BBB	1734	0	1652	18	0
2	AAA	26	0	10	0	0
3	AAA	24	0	0	0	0
3	BBB	24	0	0	0	0
4	BBB	30	0	0	2	0
5	AAA	248	0	0	2	0
5	BBB	230	0	0	2	0
All	All	4029	0	3294	29	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 4.

All (29) 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:BBB:139:ASP:OD2	1:BBB:142:HIS:HE1	1.74	0.71
1:AAA:155:GLU:H	1:AAA:205:HIS:CE1	2.20	0.59
1:AAA:218:HIS:O	1:AAA:219:HIS:HB2	2.02	0.58
1:AAA:31:VAL:O	1:AAA:35:GLU:HG3	2.04	0.58
1:AAA:116:LEU:HD22	1:AAA:212:PRO:HB2	1.84	0.58
1:BBB:155:GLU:H	1:BBB:205:HIS:CE1	2.23	0.57
1:BBB:188:HIS:HD2	5:BBB:1249:HOH:O	1.88	0.56
1:BBB:31:VAL:O	1:BBB:35:GLU:HG3	2.05	0.56
1:AAA:141:ARG:HG2	1:AAA:163:ASN:ND2	2.22	0.55
1:AAA:141:ARG:HG2	1:AAA:163:ASN:CG	2.32	0.54
1:BBB:188:HIS:N	4:BBB:1003:SO4:O2	2.40	0.53
1:AAA:154:TYR:HA	1:AAA:205:HIS:CE1	2.46	0.49
1:AAA:188:HIS:O	1:AAA:193:GLU:HA	2.12	0.49
1:BBB:154:TYR:HA	1:BBB:205:HIS:CE1	2.48	0.49
1:AAA:7:TRP:CZ2	1:AAA:132:GLY:HA2	2.50	0.47
1:BBB:139:ASP:OD2	1:BBB:142:HIS:CE1	2.62	0.46
1:BBB:7:TRP:CE2	1:BBB:132:GLY:HA2	2.52	0.45
1:BBB:25:CYS:HB2	1:BBB:165:GLY:O	2.18	0.43
1:AAA:187:GLY:HA2	5:AAA:1257:HOH:O	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BBB:114:THR:O	1:BBB:213:SER:HA	2.20	0.42
1:BBB:52:ASN:HD22	1:BBB:52:ASN:C	2.28	0.42
1:BBB:173:ASP:HA	1:BBB:177:LYS:O	2.20	0.42
1:BBB:219:HIS:CD2	1:BBB:219:HIS:N	2.87	0.42
1:BBB:120:ARG:HB3	4:BBB:1004:SO4:O3	2.20	0.41
1:AAA:25:CYS:SG	5:AAA:1243:HOH:O	2.61	0.41
1:BBB:0:ILE:O	1:BBB:0:ILE:HG13	2.21	0.41
1:BBB:122:ASP:HB2	5:BBB:1203:HOH:O	2.21	0.40
1:BBB:19:GLN:NE2	1:BBB:185:SER:OG	2.55	0.40
1:BBB:41:LYS:HE3	1:BBB:217:ILE:HG22	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	AAA	218/225 (97%)	211 (97%)	7 (3%)	0	100	100
1	BBB	220/225 (98%)	213 (97%)	6 (3%)	1 (0%)	24	26
All	All	438/450 (97%)	424 (97%)	13 (3%)	1 (0%)	43	50

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	BBB	0	ILE

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	AAA	181/186 (97%)	174 (96%)	7 (4%)	28	37
1	BBB	183/186 (98%)	172 (94%)	11 (6%)	17	20
All	All	364/372 (98%)	346 (95%)	18 (5%)	22	27

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	AAA	0	ILE
1	AAA	41	LYS
1	AAA	52	ASN
1	AAA	90	TYR
1	AAA	141	ARG
1	AAA	158	CYS
1	AAA	177	LYS
1	BBB	-1	ARG
1	BBB	1	LEU
1	BBB	52	ASN
1	BBB	60	LYS
1	BBB	90	TYR
1	BBB	104	LYS
1	BBB	143	PRO
1	BBB	158	CYS
1	BBB	161	ASN
1	BBB	173	ASP
1	BBB	219	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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](#)

10 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	$\# Z > 2$
4	SO4	BBB	1006	-	4,4,4	0.46	0	6,6,6	0.39	0
4	SO4	BBB	1005	-	4,4,4	0.32	0	6,6,6	0.26	0
4	SO4	BBB	1001	-	4,4,4	0.31	0	6,6,6	0.18	0
2	FLC	AAA	1001	-	12,12,12	1.50	1 (8%)	17,17,17	1.12	1 (5%)
4	SO4	BBB	1002	-	4,4,4	0.27	0	6,6,6	0.08	0
4	SO4	BBB	1003	-	4,4,4	0.42	0	6,6,6	0.13	0
4	SO4	BBB	1004	-	4,4,4	0.31	0	6,6,6	0.14	0
3	Q1E	AAA	1003	-	25,26,26	2.16	8 (32%)	31,37,37	2.23	12 (38%)
2	FLC	AAA	1002	-	12,12,12	1.68	1 (8%)	17,17,17	2.16	7 (41%)
3	Q1E	BBB	1007	-	25,26,26	2.26	8 (32%)	31,37,37	2.54	12 (38%)

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	FLC	AAA	1001	-	-	6/16/16/16	-
3	Q1E	BBB	1007	-	-	1/18/28/28	0/3/3/3
3	Q1E	AAA	1003	-	-	0/18/28/28	0/3/3/3
2	FLC	AAA	1002	-	-	4/16/16/16	-

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	BBB	1007	Q1E	N1-N	-6.66	1.25	1.39
3	AAA	1003	Q1E	N2-N3	-5.77	1.27	1.38
3	AAA	1003	Q1E	N1-N	-5.36	1.28	1.39
3	BBB	1007	Q1E	N2-N3	-4.72	1.29	1.38
2	AAA	1002	FLC	CB-CBC	4.19	1.57	1.53
3	AAA	1003	Q1E	C9-S	4.13	1.82	1.78
2	AAA	1001	FLC	CB-CBC	3.65	1.57	1.53
3	BBB	1007	Q1E	C4-C1	-3.48	1.35	1.41
3	BBB	1007	Q1E	O1-S	3.15	1.46	1.43
3	AAA	1003	Q1E	C6-N5	-2.85	1.45	1.47
3	BBB	1007	Q1E	C7-N5	-2.78	1.45	1.47
3	BBB	1007	Q1E	C1-N2	-2.56	1.33	1.38
3	BBB	1007	Q1E	O2-C11	-2.39	1.18	1.23
3	BBB	1007	Q1E	C-N1	2.39	1.36	1.31
3	AAA	1003	Q1E	C3-C4	2.29	1.40	1.35
3	AAA	1003	Q1E	O1-S	2.13	1.45	1.43
3	AAA	1003	Q1E	C3-C2	-2.08	1.37	1.41
3	AAA	1003	Q1E	C1-N2	-2.00	1.34	1.38

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	BBB	1007	Q1E	C7-N5-C6	-7.14	103.91	112.12
3	BBB	1007	Q1E	O1-S-C9	6.11	118.00	107.97
3	AAA	1003	Q1E	O-S-C9	-5.12	99.54	107.97
3	BBB	1007	Q1E	O-S-C9	-4.95	99.83	107.97
3	AAA	1003	Q1E	O1-S-C9	4.59	115.50	107.97
2	AAA	1002	FLC	OB2-CBC-CB	4.57	121.92	113.14
2	AAA	1002	FLC	OB1-CBC-CB	-4.34	113.69	122.09
3	AAA	1003	Q1E	O1-S-O	-4.19	114.77	118.99
3	BBB	1007	Q1E	O1-S-N5	-3.87	103.51	107.28
3	BBB	1007	Q1E	N2-C-N1	-3.78	105.23	109.83
3	AAA	1003	Q1E	C8-N4-C5	3.59	120.00	112.68
2	AAA	1002	FLC	CB-CG-CGC	3.45	123.34	113.92
3	AAA	1003	Q1E	C7-N5-C6	-3.35	108.26	112.12
3	AAA	1003	Q1E	C6-N5-S	-3.04	112.75	118.52
3	BBB	1007	Q1E	C1-N-N1	-2.90	104.35	107.15
3	AAA	1003	Q1E	C9-S-N5	2.77	113.26	105.82
3	AAA	1003	Q1E	C10-C11-N6	-2.72	112.58	116.39
3	AAA	1003	Q1E	O1-S-N5	-2.69	104.66	107.28
2	AAA	1002	FLC	CG-CB-CBC	2.60	115.79	110.03
3	BBB	1007	Q1E	C-N1-N	2.58	112.79	107.66
3	AAA	1003	Q1E	C7-N5-S	-2.58	113.62	118.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	BBB	1007	Q1E	C4-C1-N	-2.51	126.65	132.71
3	BBB	1007	Q1E	C9-S-N5	2.44	112.36	105.82
2	AAA	1002	FLC	OHB-CB-CA	-2.43	103.84	109.38
3	BBB	1007	Q1E	C10-C11-N6	2.41	119.78	116.39
2	AAA	1002	FLC	CB-CA-CAC	2.35	120.33	113.92
3	BBB	1007	Q1E	O1-S-O	-2.29	116.68	118.99
3	AAA	1003	Q1E	N2-C-N1	-2.21	107.14	109.83
2	AAA	1001	FLC	OB2-CBC-CB	2.15	117.25	113.14
3	BBB	1007	Q1E	N2-C1-N	2.08	111.78	108.67
2	AAA	1002	FLC	OG1-CGC-CG	-2.06	117.12	122.95
3	AAA	1003	Q1E	O-S-N5	2.06	109.28	107.28

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	AAA	1001	FLC	CAC-CA-CB-CBC
2	AAA	1001	FLC	CAC-CA-CB-CG
3	BBB	1007	Q1E	C7-N5-S-C9
2	AAA	1002	FLC	CA-CB-CG-CGC
2	AAA	1002	FLC	CBC-CB-CG-CGC
2	AAA	1002	FLC	OHB-CB-CG-CGC
2	AAA	1001	FLC	CAC-CA-CB-OHB
2	AAA	1002	FLC	CAC-CA-CB-OHB
2	AAA	1001	FLC	CB-CG-CGC-OG1
2	AAA	1001	FLC	CB-CA-CAC-OA1
2	AAA	1001	FLC	CB-CA-CAC-OA2

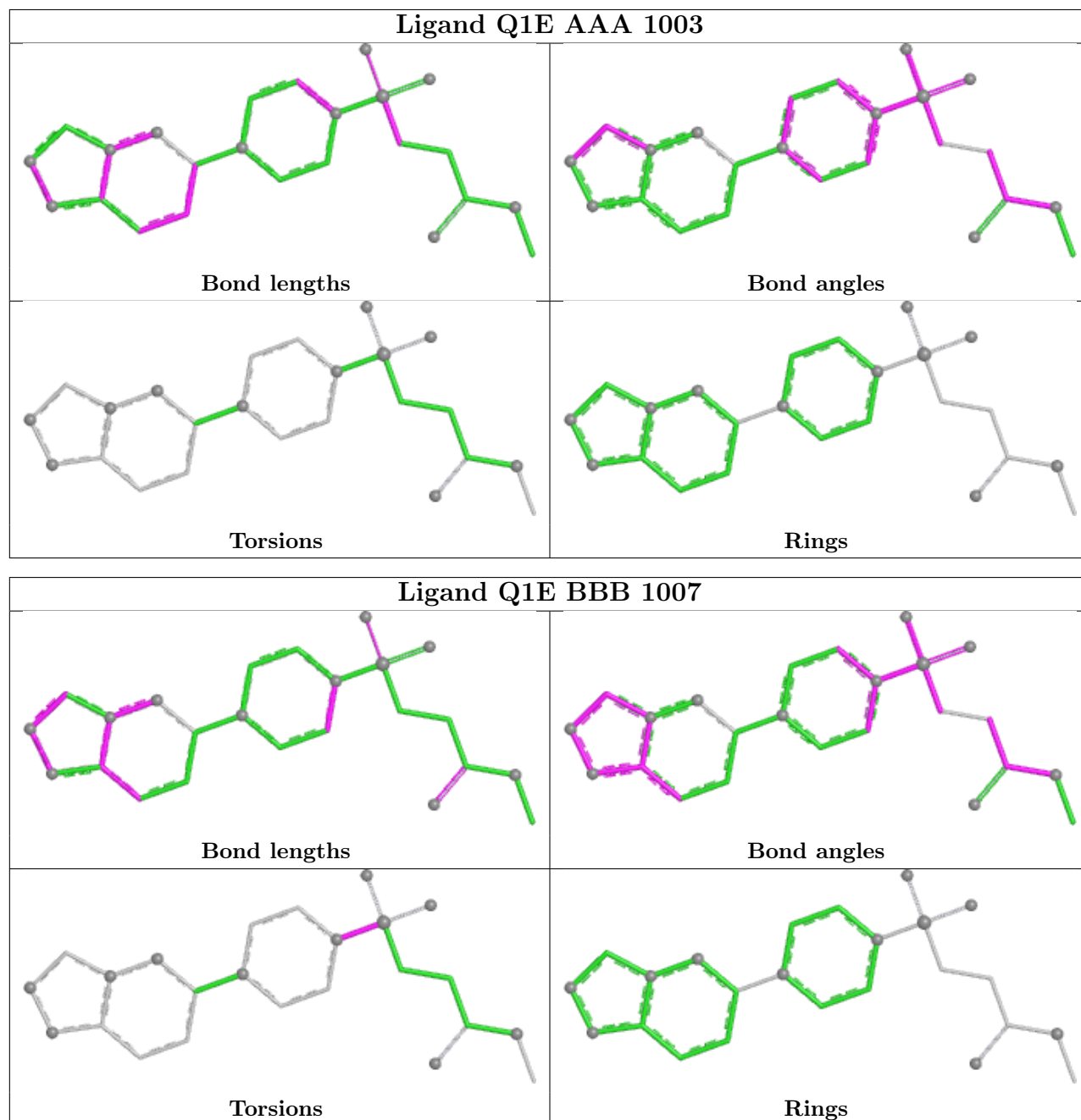
There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	BBB	1003	SO4	1	0
4	BBB	1004	SO4	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring

in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers ⓘ

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 [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	AAA	220/225 (97%)	-0.66	0	100 100	23, 32, 46, 93	0
1	BBB	222/225 (98%)	-0.60	2 (0%)	81 79	23, 32, 47, 117	0
All	All	442/450 (98%)	-0.63	2 (0%)	87 86	23, 32, 47, 117	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	BBB	220	HIS	3.5
1	BBB	0	ILE	3.2

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(Å ²)	Q<0.9
4	SO4	BBB	1001	5/5	0.62	0.13	147,150,156,160	0
4	SO4	BBB	1006	5/5	0.74	0.20	64,86,104,113	0
2	FLC	AAA	1002	13/13	0.78	0.16	58,71,81,97	0

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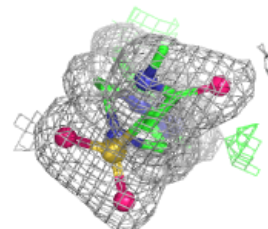
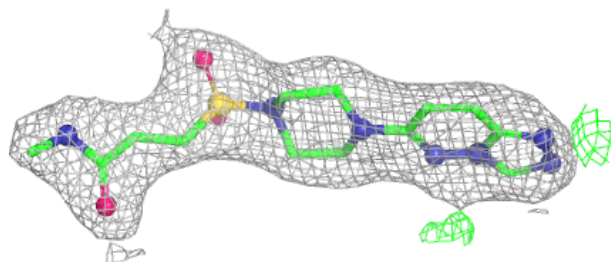
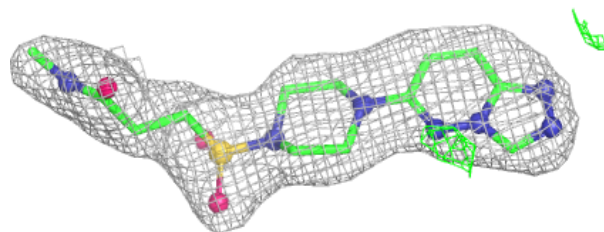
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	SO4	BBB	1002	5/5	0.80	0.12	87,92,103,105	0
4	SO4	BBB	1004	5/5	0.82	0.10	94,103,103,108	0
4	SO4	BBB	1005	5/5	0.90	0.12	71,71,73,76	5
2	FLC	AAA	1001	13/13	0.91	0.10	45,48,49,51	13
4	SO4	BBB	1003	5/5	0.94	0.09	47,48,49,50	5
3	Q1E	AAA	1003	24/24	0.97	0.06	32,36,47,50	0
3	Q1E	BBB	1007	24/24	0.98	0.05	29,34,45,48	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

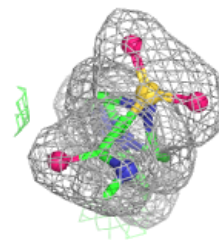
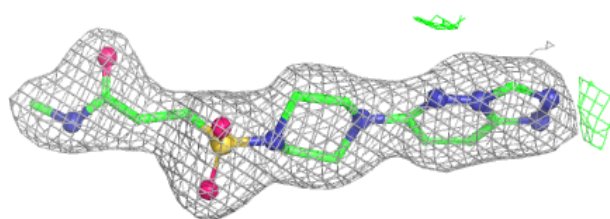
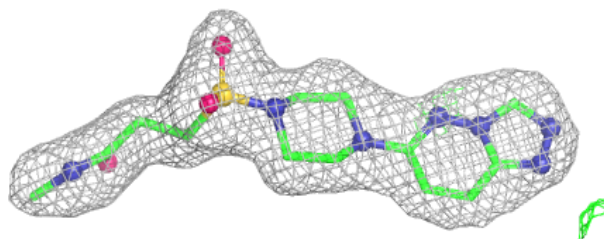
Electron density around Q1E AAA 1003:

2mF_o-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Q1E BBB 1007:

$2mF_o - DF_c$ (at 0.7 rmsd) in gray
 $mF_o - DF_c$ (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.